Full wwPDB X-ray Structure Validation Report (i)

Jun 11, 2024 – 09:58 PM EDT

PDB ID	:	1ADE
Title	:	STRUCTURE OF ADENYLOSUCCINATE SYNTHETASE PH 7 AT 25 DEGREES CELSIUS
Authors	:	Silva, M.M.; Poland, B.W.; Hoffman, C.M.; Fromm, H.J.; Honzatko, R.B.
Deposited on	:	1995-09-14
Resolution	:	2.00 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

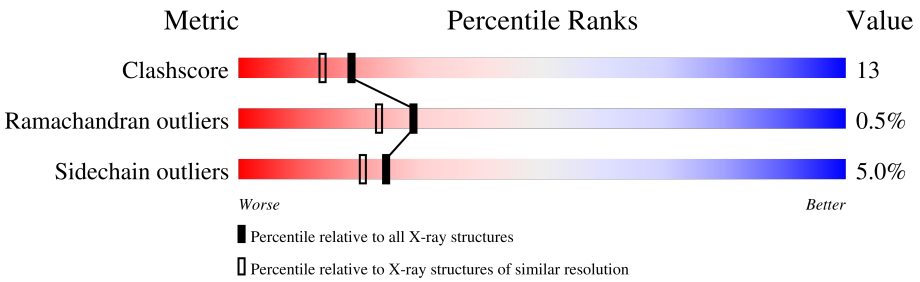
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.00 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	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)

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	A	431	72% 25% •
1	B	431	71% 28% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 9988 atoms, of which 2677 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 ADENYLOSUCCINATE SYNTHETASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	431	4150	2128	771	592	646	13	0	6	0
1	B	431	4116	2114	758	585	646	13	0	4	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	H	O		
2	A	314	942	628	314	0	0
2	B	260	780	520	260	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.

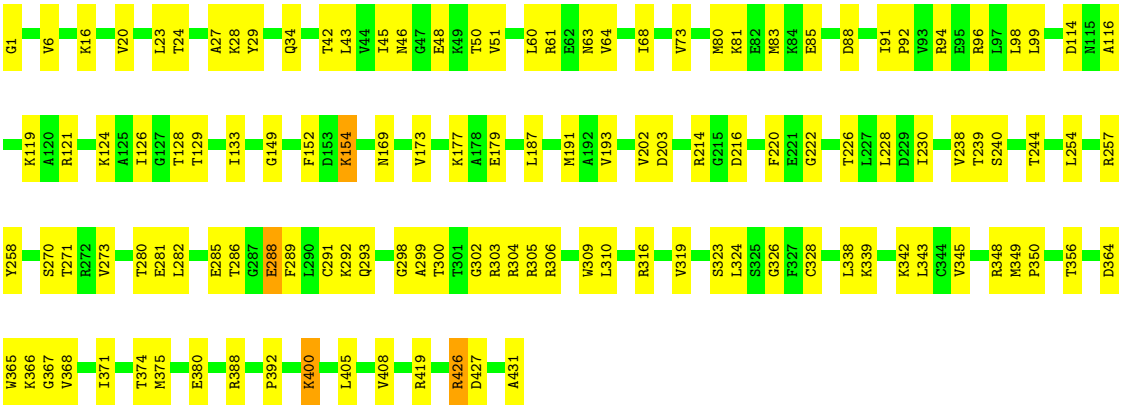
Note EDS was not executed.

• Molecule 1: ADENYLOSUCCINATE SYNTHETASE

Chain A:

• Molecule 1: ADENYLOSUCCINATE SYNTHETASE

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	73.26Å 72.18Å 82.90Å 90.00° 108.17° 90.00°	Depositor
Resolution (Å)	5.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (5.00-2.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.199 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9988	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

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	0.54	0/3437	0.79	2/4649 (0.0%)
1	B	0.54	0/3416	0.79	1/4624 (0.0%)
All	All	0.54	0/6853	0.79	3/9273 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	302	GLY	N-CA-C	7.87	132.78	113.10
1	A	343	LEU	CA-CB-CG	5.38	127.69	115.30
1	B	328	CYS	N-CA-C	-5.02	97.45	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	A	3379	771	3398	91	4
1	B	3358	758	3365	90	2
2	A	314	628	0	7	0
2	B	260	520	0	15	0
All	All	7311	2677	6763	178	4

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 13.

All (178) 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:296:GLU:O	1:A:305:ARG:HG2	1.77	0.84
1:B:273:VAL:HG11	1:B:300:THR:OG1	1.79	0.82
1:A:272:ARG:HH11	1:A:272:ARG:HG2	1.46	0.81
1:A:303:ARG:HD3	1:A:303:ARG:H	1.50	0.77
1:A:28:LYS:NZ	1:A:28:LYS:HB2	2.02	0.75
1:B:350:PRO:HG3	1:B:367:GLY:HA3	1.69	0.74
1:B:375:MET:SD	2:B:625:HOH:O	2.45	0.74
1:A:288:GLU:O	1:A:292[A]:LYS:HG2	1.88	0.74
1:B:46:ASN:HD21	1:B:91:ILE:HD11	1.51	0.73
1:A:290:LEU:HB2	1:A:307:THR:HG21	1.70	0.73
1:B:42:THR:HG22	1:B:51:VAL:HG22	1.70	0.72
1:B:285:GLU:O	1:B:288:GLU:HG3	1.89	0.72
1:A:228:LEU:O	1:A:239:THR:HG22	1.91	0.69
1:A:25:GLU:HG2	1:A:61[B]:ARG:NH1	2.09	0.68
1:A:274:GLY:H	1:A:304:ARG:HB3	1.60	0.66
1:B:73:VAL:HG21	1:B:133:ILE:HG23	1.76	0.66
1:A:83:MET:O	1:A:87:GLU:HG2	1.96	0.66
1:B:126:ILE:O	1:B:128:THR:HG23	1.95	0.65
1:B:280:THR:HG21	1:B:309:TRP:O	1.96	0.65
1:B:28:LYS:HD2	1:B:29:TYR:CE2	2.31	0.65
1:A:280:THR:HG21	1:A:309:TRP:O	1.96	0.65
1:B:298:GLY:HA2	1:B:303:ARG:O	1.98	0.64
1:A:98:LEU:HD13	1:A:202:VAL:HG11	1.81	0.63
1:B:230:ILE:HD12	1:B:240:SER:HA	1.82	0.61
1:A:273:VAL:HG11	1:A:302:GLY:HA3	1.81	0.61
1:A:289:PHE:HB2	2:A:693:HOH:O	2.00	0.60
1:B:400:LYS:O	1:B:400:LYS:HE3	2.02	0.59
1:B:342:LYS:HG2	1:B:374:THR:HG22	1.85	0.59
1:B:273:VAL:CG2	1:B:300:THR:HG21	2.32	0.58
1:A:16[B]:LYS:HE2	1:A:222:GLY:O	2.04	0.58
1:A:419[A]:ARG:HD2	1:A:419[A]:ARG:H	1.68	0.58
1:B:426:ARG:HH11	1:B:426:ARG:HB2	1.69	0.57
1:A:191:MET:HA	1:A:191:MET:CE	2.35	0.57
1:B:349:MET:CE	1:B:368:VAL:HG12	2.35	0.57
1:A:173:VAL:O	1:A:177[A]:LYS:HA	2.05	0.57
1:A:206:ASP:O	1:A:210:GLN:HG2	2.05	0.57
1:A:293:GLN:HB2	1:A:334:VAL:HG13	1.87	0.56
1:B:116:ALA:HA	1:B:119:LYS:HD3	1.88	0.56
1:B:345:VAL:HG12	1:B:405:LEU:HD13	1.87	0.56
1:B:365:TRP:O	1:B:368:VAL:HG22	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:ARG:HB3	2:B:565:HOH:O	2.05	0.56
1:B:426:ARG:O	1:B:426:ARG:HG3	2.05	0.55
1:A:227:LEU:HD11	2:A:565:HOH:O	2.06	0.55
1:B:173:VAL:O	1:B:177[A]:LYS:HA	2.06	0.55
1:A:43:LEU:HD13	1:A:45:ILE:HG13	1.89	0.55
1:A:128:THR:HG23	1:A:130:GLY:H	1.72	0.54
1:B:83:MET:SD	2:B:549:HOH:O	2.59	0.54
1:B:46:ASN:ND2	1:B:91:ILE:HD11	2.19	0.54
1:B:426:ARG:HB2	1:B:426:ARG:NH1	2.22	0.54
1:B:316:ARG:NH1	1:B:356:THR:O	2.39	0.54
1:A:323:SER:O	1:B:257:ARG:NH1	2.41	0.53
1:A:128:THR:HG23	1:A:130:GLY:N	2.23	0.53
1:A:161:LEU:O	1:A:165:MET:HG3	2.08	0.53
1:A:248:VAL:HB	2:A:663:HOH:O	2.07	0.52
1:A:282:LEU:HD22	1:A:308:GLY:HA2	1.91	0.52
1:A:61[A]:ARG:NE	1:A:64:VAL:HG23	2.25	0.52
1:A:281:GLU:OE2	1:A:306:ARG:HD2	2.10	0.51
1:B:80:MET:HG3	1:B:193:VAL:HG11	1.91	0.51
1:A:16[B]:LYS:NZ	1:A:221:GLU:OE2	2.43	0.51
1:A:28:LYS:HB2	1:A:28:LYS:HZ2	1.73	0.51
1:A:61[A]:ARG:HE	1:A:64:VAL:HG23	1.76	0.51
1:B:173:VAL:O	1:B:177[B]:LYS:HA	2.09	0.51
1:B:119:LYS:NZ	2:B:641:HOH:O	2.42	0.51
1:A:92:PRO:HB2	1:A:95:GLU:HG2	1.91	0.51
1:B:350:PRO:HG2	1:B:364:ASP:OD2	2.10	0.51
1:A:272:ARG:HH11	1:A:272:ARG:CG	2.16	0.50
1:B:302:GLY:O	1:B:303:ARG:HG3	2.11	0.50
1:A:303:ARG:O	1:A:304:ARG:HB2	2.11	0.50
1:B:364:ASP:OD1	1:B:366:LYS:HG3	2.11	0.50
1:A:37:HIS:O	1:A:131:ARG:NH1	2.44	0.50
1:B:220:PHE:CZ	1:B:254:LEU:HD22	2.48	0.49
1:A:280:THR:HG23	1:A:280:THR:O	2.12	0.49
1:B:20:VAL:O	1:B:24:THR:HG23	2.11	0.49
1:B:350:PRO:HG3	1:B:367:GLY:CA	2.42	0.49
1:A:85:GLU:O	1:A:89:ARG:HD3	2.13	0.49
1:B:63:ASN:HA	2:B:511:HOH:O	2.12	0.49
1:B:88:ASP:HA	2:B:562:HOH:O	2.13	0.49
1:A:289:PHE:CE1	1:A:293:GLN:NE2	2.81	0.49
1:A:255:GLY:HA3	1:B:323:SER:HB3	1.95	0.49
1:B:282:LEU:HD13	1:B:286:THR:HB	1.94	0.49
1:A:61[A]:ARG:HH22	1:A:63:ASN:HB2	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:380:GLU:HG3	1:B:392:PRO:HB3	1.95	0.48
1:B:187:LEU:O	1:B:191:MET:HG2	2.13	0.48
1:A:274:GLY:N	1:A:304:ARG:HB3	2.26	0.48
1:B:419[A]:ARG:NH1	2:B:663:HOH:O	2.46	0.48
1:B:214:ARG:HD2	1:B:216:ASP:OD2	2.13	0.48
1:B:45:ILE:N	1:B:48:GLU:O	2.46	0.47
1:B:68:ILE:HB	1:B:99:LEU:HD23	1.97	0.47
1:B:326:GLY:HA2	1:B:408:VAL:CG1	2.44	0.47
1:B:273:VAL:HG22	1:B:305:ARG:HG3	1.97	0.47
1:A:177[A]:LYS:HG2	2:A:487:HOH:O	2.15	0.47
1:B:203:ASP:HB3	2:B:559:HOH:O	2.14	0.47
1:B:273:VAL:HG22	1:B:300:THR:HG21	1.97	0.47
1:A:73:VAL:HG13	1:A:137:TYR:CE1	2.50	0.47
1:A:250:THR:HA	2:B:486:HOH:O	2.14	0.47
1:B:281:GLU:OE2	1:B:306:ARG:HD2	2.14	0.47
1:B:273:VAL:HG21	1:B:300:THR:HG21	1.97	0.46
1:A:296:GLU:C	1:A:298:GLY:H	2.18	0.46
1:A:365:TRP:HA	1:A:368:VAL:HG13	1.97	0.46
1:B:98:LEU:HD13	1:B:202:VAL:HG11	1.97	0.46
1:A:88:ASP:HB2	1:A:89:ARG:HH11	1.79	0.46
1:A:272:ARG:HG2	1:A:272:ARG:NH1	2.23	0.46
1:A:359:PRO:HG2	1:A:365:TRP:CE2	2.50	0.46
1:B:292:LYS:HG3	1:B:293:GLN:N	2.30	0.46
1:B:319:VAL:HG22	1:B:324:LEU:HD12	1.98	0.46
1:B:126:ILE:O	1:B:126:ILE:HG23	2.16	0.46
1:B:348:ARG:NH1	1:B:371:ILE:HD11	2.30	0.46
1:A:191:MET:HA	1:A:191:MET:HE2	1.98	0.46
1:A:261:TYR:CE1	1:A:428:PRO:HB3	2.50	0.46
1:B:16:LYS:HE2	1:B:222:GLY:O	2.16	0.45
1:B:61:ARG:HD3	2:B:572:HOH:O	2.15	0.45
1:B:28:LYS:O	1:B:28:LYS:HD3	2.16	0.45
1:B:149:GLY:O	1:B:152:PHE:HD2	1.99	0.45
1:A:410:ILE:O	1:A:425:LEU:HD12	2.17	0.44
1:A:22:LEU:CD1	1:A:419[B]:ARG:HG2	2.48	0.44
1:A:73:VAL:HG23	1:A:106:ILE:HD11	2.00	0.44
1:A:290:LEU:CB	1:A:307:THR:HG21	2.42	0.44
1:B:228:LEU:O	1:B:238:VAL:HB	2.17	0.44
1:B:281:GLU:OE2	1:B:306:ARG:HB3	2.18	0.44
1:A:267:LYS:NZ	1:A:296:GLU:OE2	2.48	0.44
1:A:255:GLY:HA2	1:A:256:PRO:HD3	1.87	0.43
1:B:1:GLY:N	1:B:431:ALA:OXT	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:LEU:N	1:A:343:LEU:HD23	2.34	0.43
1:B:50:THR:HG22	1:B:51:VAL:H	1.82	0.43
1:A:271:THR:HA	1:A:306:ARG:O	2.18	0.43
1:B:169:ASN:O	1:B:173:VAL:HG22	2.19	0.43
1:A:29:TYR:CD1	1:A:65:THR:HB	2.54	0.43
1:A:144:ARG:HD3	2:A:649:HOH:O	2.17	0.43
1:A:22:LEU:HD12	1:A:419[B]:ARG:HG2	2.01	0.43
1:A:286:THR:HG21	1:A:309:TRP:CH2	2.53	0.43
1:B:81[A]:LYS:NZ	2:B:531:HOH:O	2.51	0.43
1:B:129:THR:HA	2:B:480:HOH:O	2.18	0.43
1:A:92:PRO:HB2	1:A:95:GLU:CG	2.48	0.42
1:B:1:GLY:H1	1:B:431:ALA:C	2.19	0.42
1:B:280:THR:HG22	2:B:619:HOH:O	2.17	0.42
1:B:6:VAL:HG21	1:B:23:LEU:HD12	2.02	0.42
1:A:11:TRP:CZ2	1:A:227:LEU:HD13	2.54	0.42
1:A:202:VAL:HG22	1:A:203:ASP:N	2.35	0.42
1:A:393:GLN:HB2	2:A:712:HOH:O	2.19	0.42
1:A:43:LEU:HD12	1:A:50:THR:O	2.18	0.42
1:A:349:MET:CE	1:A:368:VAL:HG12	2.50	0.42
1:B:92:PRO:O	1:B:96:ARG:HD3	2.19	0.42
1:A:16[A]:LYS:NZ	1:A:223:ALA:O	2.52	0.42
1:B:27:ALA:O	1:B:64:VAL:HG22	2.19	0.42
1:A:412:ILE:HG22	1:A:413:ILE:N	2.35	0.42
1:B:226:THR:HG23	1:B:244:THR:HG22	2.01	0.42
1:B:345:VAL:HG12	1:B:405:LEU:CD1	2.49	0.42
1:A:160:LYS:O	1:A:164:VAL:HG23	2.20	0.42
1:A:323:SER:HB3	1:B:258:TYR:CE2	2.55	0.42
1:B:380:GLU:CG	1:B:392:PRO:HB3	2.49	0.42
1:A:202:VAL:HG23	2:A:608:HOH:O	2.19	0.42
1:B:45:ILE:HG23	1:B:60:LEU:HD21	2.01	0.41
1:A:32:ARG:HG2	1:A:56:PRO:HG3	2.02	0.41
1:B:121:ARG:HB3	1:B:121:ARG:NH1	2.36	0.41
1:A:26:ARG:HH12	1:A:430:ASP:CG	2.24	0.41
1:A:286:THR:HG21	1:A:309:TRP:CZ2	2.56	0.41
1:A:289:PHE:O	1:A:293:GLN:HG2	2.19	0.41
1:A:401:ARG:HA	1:A:401:ARG:HD2	1.60	0.41
1:A:61[A]:ARG:HH11	1:A:61[A]:ARG:HG2	1.86	0.41
1:A:303:ARG:H	1:A:303:ARG:CD	2.26	0.41
1:A:343:LEU:HD21	1:A:375:MET:CE	2.49	0.41
1:B:292:LYS:HD3	2:B:529:HOH:O	2.21	0.41
1:B:310:LEU:HB3	1:B:343:LEU:HD23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ASP:O	1:A:238:VAL:CG2	2.69	0.41
1:B:85:GLU:O	1:B:88:ASP:HB2	2.21	0.41
1:B:291:CYS:HB2	1:B:304:ARG:HH21	1.86	0.41
1:B:349:MET:HE2	1:B:368:VAL:HG12	2.03	0.41
1:A:88:ASP:CB	1:A:89:ARG:HH11	2.34	0.40
1:A:382:THR:HG22	1:A:392:PRO:HD2	2.04	0.40
1:A:28:LYS:HG3	1:A:29:TYR:CD2	2.56	0.40
1:B:289:PHE:CD2	1:B:338:LEU:HD21	2.55	0.40
1:B:380:GLU:CD	1:B:380:GLU:H	2.24	0.40
1:B:154:LYS:NZ	2:B:655:HOH:O	2.55	0.40
1:B:285:GLU:H	1:B:285:GLU:CD	2.25	0.40
1:A:351:ASP:OD1	1:A:353:ARG:HG2	2.22	0.40
1:B:271:THR:HA	1:B:306:ARG:O	2.22	0.40

All (4) 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:119:LYS:HZ3	1:A:354:GLU:OE2[2_556]	1.38	0.22
1:A:188:ASP:OD2	1:B:292:LYS:HZ3[2_455]	1.39	0.21
1:A:62:GLU:OE1	1:B:388:ARG:HH22[2_555]	1.51	0.09
1:A:182:ASP:OD2	1:A:339:LYS:HZ2[2_556]	1.55	0.05

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	435/431 (101%)	414 (95%)	20 (5%)	1 (0%)	47	44
1	B	433/431 (100%)	411 (95%)	19 (4%)	3 (1%)	22	16
All	All	868/862 (101%)	825 (95%)	39 (4%)	4 (0%)	29	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	124	LYS
1	A	427	ASP
1	B	299	ALA
1	B	427	ASP

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	359/353 (102%)	335 (93%)	24 (7%)	16	11
1	B	357/353 (101%)	346 (97%)	11 (3%)	40	40
All	All	716/706 (101%)	681 (95%)	35 (5%)	24	21

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	28	LYS
1	A	34	GLN
1	A	43	LEU
1	A	46	ASN
1	A	79	LEU
1	A	113	LEU
1	A	114	ASP
1	A	128	THR
1	A	147	ARG
1	A	179	GLU
1	A	191	MET
1	A	207	LEU
1	A	238	VAL
1	A	278	PHE
1	A	280	THR
1	A	282	LEU
1	A	285	GLU
1	A	303	ARG

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Mol	Chain	Res	Type
1	A	307	THR
1	A	340	GLU
1	A	343	LEU
1	A	389	SER
1	A	401	ARG
1	B	34	GLN
1	B	43	LEU
1	B	114	ASP
1	B	154	LYS
1	B	179	GLU
1	B	239	THR
1	B	270	SER
1	B	288	GLU
1	B	339	LYS
1	B	400	LYS
1	B	426	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	224	GLN
1	B	2	ASN
1	B	46	ASN
1	B	171	GLN
1	B	184	GLN
1	B	293	GLN
1	B	295	ASN

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)

There are no ligands in this entry.

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.