

Full wwPDB X-ray Structure Validation Report (i)

Aug 22, 2023 – 08:45 PM EDT

PDB ID	:	3CSU
Title	:	CATALYTIC TRIMER OF ESCHERICHIA COLI ASPARTATE TRAN-
		SCARBAMOYLASE
Authors	:	Beernink, P.T.; Endrizzi, J.A.; Alber, T.; Schachman, H.K.
Deposited on	:	1999-04-22
Resolution	:	1.88 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 (1)) 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)	:	2.35

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 1.88 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	10282 (1.90-1.86)		
Ramachandran outliers	138981	10152 (1.90-1.86)		
Sidechain outliers	138945	10152 (1.90-1.86)		

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	310	57%	30%	5%• 7%				
1	В	310	61%	26%	5% • 7%				
1	С	310	63%	29%	6% ••				



3CSU

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7601 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 PROTEIN (ASPARTATE CARBAMOYLTRANSFERASE).
 Mol Chain Residues Atoms ZeroOcc AltConf Trace
 A 288 Total C N O S

1	1 A	288	Total	С	N	0	S	0	0	0
			2225	1413	385	418	9			
1	1 B	266	Total	С	Ν	0	S	0	1	0
1		200	2219	1409	384	417	9			
1 C	304	Total	С	Ν	0	S	0	0	0	
		2366	1494	415	448	9		0	0	

• Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	1	Total Ca 1 1	0	0
2	С	1	Total Ca 1 1	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	234	Total O 234 234	0	0
3	В	233	Total O 233 233	0	0
3	С	322	Total O 322 322	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: PROTEIN (ASPARTATE CARBAMOYLTRANSFERASE)







4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	P 21 21 21	Depositor	
Cell constants	56.12Å 82.54 Å 210.52 Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	20.00 - 1.88	Depositor	
% Data completeness	94.6 (20.00-1.88)	Depositor	
(in resolution range)	51.0 (20.00 1.00)	Depositor	
R_{merge}	0.09	Depositor	
R_{sym}	(Not available)	Depositor	
Refinement program	TNT 5E	Depositor	
R, R_{free}	0.190 , 0.291	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	7601	wwPDB-VP	
Average B, all atoms $(Å^2)$	29.0	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: CA

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	Bo	nd lengths	Bond angles		
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.77	10/2267~(0.4%)	1.22	33/3078~(1.1%)	
1	В	0.76	9/2266~(0.4%)	1.22	34/3081~(1.1%)	
1	С	0.79	13/2410~(0.5%)	1.25	39/3272~(1.2%)	
All	All	0.77	32/6943~(0.5%)	1.23	$106/9431 \ (1.1\%)$	

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	60	GLU	CD-OE2	8.20	1.34	1.25
1	В	50	GLU	CD-OE2	7.18	1.33	1.25
1	В	117	GLU	CD-OE2	6.36	1.32	1.25
1	С	147	GLU	CD-OE2	6.25	1.32	1.25
1	С	216	GLU	CD-OE2	6.18	1.32	1.25
1	С	60	GLU	CD-OE2	6.18	1.32	1.25
1	С	204	GLU	CD-OE2	6.06	1.32	1.25
1	А	109	GLU	CD-OE2	5.85	1.32	1.25
1	С	109	GLU	CD-OE2	5.85	1.32	1.25
1	С	272	GLU	CD-OE2	5.80	1.32	1.25
1	В	217	GLU	CD-OE2	5.79	1.32	1.25
1	С	117	GLU	CD-OE2	5.75	1.31	1.25
1	С	86	GLU	CD-OE2	5.66	1.31	1.25
1	С	221	GLU	CD-OE2	5.65	1.31	1.25
1	А	117	GLU	CD-OE2	5.63	1.31	1.25
1	С	37	GLU	CD-OE2	5.57	1.31	1.25
1	С	239	GLU	CD-OE2	5.55	1.31	1.25
1	В	60	GLU	CD-OE2	5.51	1.31	1.25
1	А	221	GLU	CD-OE2	5.43	1.31	1.25
1	А	204	GLU	CD-OE2	5.42	1.31	1.25
1	В	204	GLU	CD-OE2	5.42	1.31	1.25
1	С	50	GLU	CD-OE1	5.40	1.31	1.25
1	В	216	GLU	CD-OE2	5.32	1.31	1.25



Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	В	272	GLU	CD-OE2	5.28	1.31	1.25
1	А	86	GLU	CD-OE2	5.25	1.31	1.25
1	В	37	GLU	CD-OE2	5.24	1.31	1.25
1	А	217	GLU	CD-OE2	5.23	1.31	1.25
1	А	216	GLU	CD-OE2	5.20	1.31	1.25
1	А	37	GLU	CD-OE2	5.14	1.31	1.25
1	В	221	GLU	CD-OE2	5.08	1.31	1.25
1	С	217	GLU	CD-OE2	5.06	1.31	1.25
1	А	50	GLU	CD-OE2	5.01	1.31	1.25

All (106) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	С	162	ASP	CB-CG-OD2	-9.57	109.69	118.30
1	В	90	ASP	CB-CG-OD2	-8.32	110.81	118.30
1	А	19	ASP	CB-CG-OD2	-7.90	111.19	118.30
1	В	162	ASP	CB-CG-OD1	7.82	125.34	118.30
1	А	19	ASP	CB-CG-OD1	7.69	125.22	118.30
1	А	278	ASP	CB-CG-OD2	-7.61	111.45	118.30
1	С	278	ASP	CB-CG-OD1	7.51	125.06	118.30
1	А	129	ASP	CB-CG-OD2	-7.39	111.65	118.30
1	В	203	ASP	CB-CG-OD1	7.37	124.93	118.30
1	С	14	ASP	CB-CG-OD2	-7.22	111.80	118.30
1	С	162	ASP	CB-CG-OD1	7.14	124.72	118.30
1	С	276	ASP	CB-CG-OD2	-7.13	111.88	118.30
1	В	14	ASP	CB-CG-OD2	-7.12	111.89	118.30
1	В	203	ASP	CB-CG-OD2	-7.11	111.90	118.30
1	А	90	ASP	CB-CG-OD2	-7.08	111.93	118.30
1	В	162	ASP	CB-CG-OD2	-7.05	111.96	118.30
1	В	200	ASP	CB-CG-OD2	-7.05	111.96	118.30
1	В	129	ASP	CB-CG-OD2	-7.01	111.99	118.30
1	С	129	ASP	CB-CG-OD2	-6.93	112.06	118.30
1	С	183	ARG	NE-CZ-NH2	-6.87	116.87	120.30
1	С	306	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	С	153	ASP	CB-CG-OD2	-6.82	112.16	118.30
1	А	18	ASP	CB-CG-OD1	6.79	124.41	118.30
1	В	153	ASP	CB-CG-OD1	6.78	124.40	118.30
1	В	276	ASP	CB-CG-OD2	-6.74	112.23	118.30
1	С	278	ASP	CB-CG-OD2	-6.68	112.29	118.30
1	С	271	ASP	CB-CG-OD1	6.62	124.26	118.30
1	А	18	ASP	CB-CG-OD2	-6.62	112.34	118.30
1	В	153	ASP	CB-CG-OD2	-6.60	112.36	118.30



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	$Ideal(^{o})$
1	С	223	ASP	CB-CG-OD2	-6.53	112.42	118.30
1	С	276	ASP	CB-CG-OD1	6.53	124.17	118.30
1	А	162	ASP	CB-CG-OD2	-6.50	112.45	118.30
1	А	75	ASP	CB-CG-OD2	-6.47	112.48	118.30
1	С	180	ASP	CB-CG-OD2	-6.46	112.49	118.30
1	А	14	ASP	CB-CG-OD1	6.41	124.07	118.30
1	А	307	ASP	CB-CG-OD2	-6.38	112.55	118.30
1	А	14	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	В	19	ASP	CB-CG-OD2	-6.37	112.57	118.30
1	В	307	ASP	CB-CG-OD2	-6.36	112.58	118.30
1	А	271	ASP	CB-CG-OD2	-6.35	112.58	118.30
1	А	162	ASP	CB-CG-OD1	6.35	124.02	118.30
1	В	14	ASP	CB-CG-OD1	6.33	124.00	118.30
1	С	296	ARG	NE-CZ-NH2	-6.32	117.14	120.30
1	В	141	ASP	CB-CG-OD2	-6.32	112.61	118.30
1	В	90	ASP	CB-CG-OD1	6.30	123.97	118.30
1	В	190	ASP	CB-CG-OD2	-6.29	112.64	118.30
1	С	75	ASP	CB-CG-OD2	-6.28	112.64	118.30
1	А	296	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	В	253	ASP	CB-CG-OD2	-6.24	112.69	118.30
1	В	75	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	А	278	ASP	CB-CG-OD1	6.20	123.88	118.30
1	С	14	ASP	CB-CG-OD1	6.18	123.86	118.30
1	А	190	ASP	CB-CG-OD2	-6.17	112.75	118.30
1	А	200	ASP	CB-CG-OD2	-6.14	112.77	118.30
1	С	271	ASP	CB-CG-OD2	-6.13	112.79	118.30
1	С	98	TYR	CB-CA-C	-6.12	98.16	110.40
1	С	307	ASP	CB-CG-OD2	-6.06	112.85	118.30
1	А	75	ASP	CB-CG-OD1	6.04	123.74	118.30
1	А	276	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	А	271	ASP	CB-CG-OD1	5.99	123.69	118.30
1	С	236	ASP	CB-CG-OD2	-5.95	112.94	118.30
1	А	183	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	А	223	ASP	CB-CG-OD2	-5.93	112.96	118.30
1	В	190	ASP	CB-CG-OD1	5.92	123.63	118.30
1	В	200	ASP	CB-CG-OD1	5.91	123.62	118.30
1	В	307	ASP	CB-CG-OD1	5.90	123.61	118.30
1	С	269	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	С	190	ASP	CB-CG-OD2	-5.89	113.00	118.30
1	А	276	ASP	CB-CG-OD1	5.87	123.59	118.30
1	С	18	ASP	CB-CG-OD2	-5.86	113.02	118.30
1	A	153	ASP	CB-CG-OD2	-5.85	113.04	118.30

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	75	ASP	CB-CG-OD1	5.83	123.55	118.30
1	С	190	ASP	CB-CG-OD1	5.79	123.52	118.30
1	С	203	ASP	CB-CG-OD2	-5.74	113.14	118.30
1	С	90	ASP	CB-CG-OD2	-5.74	113.14	118.30
1	С	141	ASP	CB-CG-OD2	-5.69	113.18	118.30
1	В	18	ASP	CB-CG-OD2	-5.60	113.26	118.30
1	С	223	ASP	CB-CG-OD1	5.58	123.32	118.30
1	С	17	ARG	NE-CZ-NH1	5.57	123.09	120.30
1	В	18	ASP	CB-CG-OD1	5.54	123.29	118.30
1	В	19	ASP	CB-CG-OD1	5.52	123.27	118.30
1	В	278	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	С	141	ASP	CB-CG-OD1	5.49	123.24	118.30
1	С	236	ASP	CB-CG-OD1	5.41	123.17	118.30
1	С	153	ASP	CB-CG-OD1	5.34	123.11	118.30
1	А	253	ASP	CB-CG-OD1	5.33	123.10	118.30
1	С	100	ASP	CB-CG-OD2	-5.33	113.50	118.30
1	В	100	ASP	CB-CG-OD1	5.32	123.09	118.30
1	С	307	ASP	CB-CG-OD1	5.31	123.08	118.30
1	А	129	ASP	CB-CG-OD1	5.27	123.04	118.30
1	С	75	ASP	CB-CG-OD1	5.26	123.03	118.30
1	С	100	ASP	CB-CG-OD1	5.25	123.03	118.30
1	А	200	ASP	CB-CG-OD1	5.24	123.01	118.30
1	В	100	ASP	CB-CG-OD2	-5.23	113.59	118.30
1	В	121	ASN	N-CA-CB	5.23	120.01	110.60
1	А	180	ASP	CB-CG-OD1	5.22	123.00	118.30
1	С	180	ASP	CB-CG-OD1	5.18	122.96	118.30
1	С	200	ASP	CB-CG-OD2	-5.15	113.67	118.30
1	В	180	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	А	203	ASP	CB-CG-OD1	5.12	122.90	118.30
1	В	278	ASP	CB-CG-OD1	5.10	122.89	118.30
1	А	190	ASP	CB-CG-OD1	5.09	122.88	118.30
1	А	253	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	В	105	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	А	65	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	В	253	ASP	CB-CG-OD1	5.05	122.85	118.30

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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	2225	0	2213	89	0
1	В	2219	0	2207	64	0
1	С	2366	0	2362	70	0
2	В	1	0	0	0	0
2	С	1	0	0	0	0
3	А	234	0	0	16	0
3	В	233	0	0	7	0
3	С	322	0	0	11	0
All	All	7601	0	6782	222	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 16.

All (222) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:C:8:HIS:HD2	1:C:124:VAL:H	1.11	0.95
1:C:47:CYS:HB3	1:C:73:PHE:CZ	2.05	0.92
1:B:114:LEU:HD11	1:B:118:PHE:HE2	1.32	0.91
1:C:114:LEU:HD11	1:C:118:PHE:HE2	1.42	0.84
1:A:8:HIS:HD2	1:A:124:VAL:H	1.22	0.83
1:B:279:LYS:H	1:B:279:LYS:HZ2	1.28	0.82
1:A:164:LYS:HD3	1:A:165:TYR:CE1	2.15	0.82
1:A:1:ALA:HB2	1:A:306:ARG:NH1	1.95	0.81
1:C:256:ASN:HD22	1:C:256:ASN:H	1.28	0.81
1:B:12:ILE:HD12	1:B:15:LEU:HD12	1.62	0.80
1:C:8:HIS:CD2	1:C:124:VAL:H	1.99	0.78
1:A:50:GLU:HB3	3:A:448:HOH:O	1.84	0.77
1:A:154:ASN:OD1	1:A:181:GLY:HA3	1.84	0.77
1:A:12:ILE:HG13	1:A:15:LEU:HD12	1.67	0.77
1:B:307:ASP:HB2	3:B:444:HOH:O	1.85	0.76
1:A:8:HIS:CD2	1:A:124:VAL:H	2.05	0.75
1:C:164:LYS:HE2	1:C:165:TYR:CZ	2.22	0.74
1:C:279:LYS:N	1:C:279:LYS:HE3	2.03	0.73
1:A:214:SER:OG	1:A:217:GLU:HG2	1.89	0.72



	lo de pagem	Interatomic	Clash	
Atom-1	Atom-2	distance (Å)	overlap (Å)	
1:A:267:LEU:HB3	1:A:268:PRO:HA	1.71	0.71	
1:C:121:ASN:HD22	1:C:121:ASN:H	1.39	0.71	
1:B:279:LYS:H	1:B:279:LYS:NZ	1.86	0.71	
1:C:25:ALA:O	1:C:29:LYS:HG2	1.91	0.71	
1:A:174:GLN:HG2	1:A:201:MET:HE3	1.74	0.70	
1:A:174:GLN:HG2	1:A:201:MET:CE	2.21	0.69	
1:B:158:ALA:HB2	1:B:222:VAL:HG11	1.73	0.69	
1:B:31:LYS:HE3	1:B:294:PHE:CE2	2.27	0.69	
1:B:279:LYS:HZ2	1:B:279:LYS:N	1.89	0.69	
1:C:13:ASN:HA	3:C:448:HOH:O	1.92	0.68	
1:C:155:LEU:H	1:C:182:ASN:HD22	1.41	0.68	
1:B:188:ALA:HB1	1:B:189:PRO:HD2	1.75	0.68	
1:A:31:LYS:HG3	1:A:294:PHE:CE2	2.28	0.68	
1:A:133:GLN:HG2	3:A:370:HOH:O	1.93	0.68	
1:A:305:ASN:ND2	3:A:339:HOH:O	2.28	0.66	
1:B:270:VAL:HG12	1:B:271:ASP:H	1.59	0.66	
1:A:248:VAL:HG22	1:A:271:ASP:O	1.97	0.65	
1:A:8:HIS:HD2	1:A:124:VAL:N	1.93	0.65	
1:B:265:HIS:NE2	1:B:272:GLU:OE1	2.29	0.64	
1:A:105:ARG:NH2	3:A:448:HOH:O	2.29	0.64	
1:B:144:THR:O	1:B:148:THR:HG23	1.97	0.64	
1:C:114:LEU:CD1	1:C:118:PHE:HE2	2.10	0.64	
1:A:12:ILE:O	1:A:15:LEU:HB2	1.97	0.64	
1:C:114:LEU:HD12	1:C:114:LEU:O	1.99	0.63	
1:C:275:THR:O	1:C:279:LYS:NZ	2.31	0.63	
1:A:41:HIS:ND1	3:A:447:HOH:O	2.30	0.63	
1:C:131:SER:HA	3:C:630:HOH:O	1.98	0.62	
1:B:24:LEU:HD22	1:B:143:PHE:HA	1.81	0.62	
1:B:270:VAL:HG12	1:B:271:ASP:N	2.14	0.62	
1:B:114:LEU:CD1	1:B:118:PHE:HE2	2.09	0.61	
1:B:247:PHE:N	3:B:492:HOH:O	2.33	0.61	
1:A:14:ASP:HB3	3:A:313:HOH:O	2.01	0.60	
1:C:8:HIS:CD2	1:C:123:PRO:HA	2.37	0.60	
1:B:265:HIS:ND1	1:B:266:PRO:HD2	2.17	0.60	
1:C:8:HIS:HD2	1:C:124:VAL:N	1.92	0.60	
1:A:14:ASP:OD1	1:A:15:LEU:HG	2.02	0.60	
1:A:106:HIS:ND1	1:A:107:PRO:HD2	2.17	0.60	
1:C:121:ASN:HD22	1:C:121:ASN:N	1.98	0.59	
1:B:212:HIS:HB3	1:B:217:GLU:OE1	2.02	0.59	
1:C:125:LEU:HD23	1:C:299:LEU:HD23	1.83	0.59	
1:C:132:ASN:OD1	1:C:133:GLN:HG2	2.02	0.59	



		Interatomic	Clash	
Atom-1	Atom-2	distance (Å)	overlap (Å)	
1:C:114:LEU:HD11	1:C:118:PHE:CE2	2.31	0.59	
1:C:206:GLY:HA2	3:C:441:HOH:O	2.02	0.59	
1:B:114:LEU:HD11	1:B:118:PHE:CE2	2.25	0.58	
1:A:219:MET:HE3	1:A:222:VAL:HG22	1.84	0.58	
1:C:227:MET:HE3	3:C:447:HOH:O	2.02	0.58	
1:B:3:PRO:HG2	1:B:22:LEU:HD22	1.85	0.58	
1:C:155:LEU:H	1:C:182:ASN:ND2	2.02	0.57	
1:A:197:TYR:HE2	3:A:346:HOH:O	1.87	0.57	
1:C:276:ASP:O	1:C:279:LYS:HD2	2.05	0.57	
1:A:261:MET:O	1:A:282:HIS:HD2	1.88	0.56	
1:A:1:ALA:HB2	1:A:306:ARG:CZ	2.36	0.56	
1:A:219:MET:HE3	1:A:219:MET:O	2.06	0.56	
1:A:29:LYS:HB2	3:A:530:HOH:O	2.05	0.56	
1:A:164:LYS:NZ	1:A:165:TYR:OH	2.29	0.56	
1:A:33:ASN:HB3	3:A:361:HOH:O	2.06	0.55	
1:B:164:LYS:HE2	1:B:165:TYR:CZ	2.42	0.55	
1:B:66:LEU:HG	1:B:297:GLN:HG2	1.87	0.55	
1:C:64:HIS:HD2	3:C:343:HOH:O	1.90	0.55	
1:C:265:HIS:ND1	1:C:266:PRO:HD2	2.21	0.55	
1:B:205:LYS:HB2	1:B:207:ILE:HD12	1.89	0.55	
1:C:129:ASP:O	1:C:132:ASN:HB3	2.06	0.55	
1:C:174:GLN:HA	1:C:201:MET:HE1	1.87	0.55	
1:B:301:ALA:O	1:B:305:ASN:HB2	2.07	0.55	
1:B:88:LEU:HD13	3:B:410:HOH:O	2.06	0.55	
1:C:270:VAL:HG12	1:C:271:ASP:N	2.22	0.54	
1:B:265:HIS:HE2	1:B:272:GLU:CD	2.09	0.54	
1:C:37:GLU:HG2	3:C:502:HOH:O	2.07	0.54	
1:A:265:HIS:NE2	1:A:272:GLU:OE2	2.37	0.54	
1:B:78:ASN:N	1:B:78:ASN:OD1	2.33	0.53	
1:A:247:PHE:HA	3:A:469:HOH:O	2.08	0.53	
1:C:106:HIS:CG	1:C:107:PRO:HD2	2.43	0.53	
1:C:271:ASP:N	1:C:271:ASP:OD1	2.41	0.53	
1:B:306:ARG:HB2	3:B:496:HOH:O	2.08	0.53	
1:B:38:LEU:HB3	1:B:66:LEU:HD22	1.91	0.53	
1:C:195:PRO:HA	3:C:478:HOH:O	2.09	0.52	
1:B:129:ASP:HB3	1:B:132:ASN:OD1	2.10	0.52	
1:A:258:LYS:NZ	3:A:471:HOH:O	2.28	0.52	
1:A:14:ASP:HB3	3:A:476:HOH:O	2.10	0.52	
1:A:105:ARG:HD2	1:A:128:GLY:O	2.10	0.52	
1:B:174:GLN:NE2	3:B:423:HOH:O	2.43	0.52	
1:A:87:THR:HB	3:C:606:HOH:O	2.09	0.52	



Interatomic Clash						
Atom-1	Atom-2	distance (Å)	overlap (Å)			
1·A·254·LEU·HD22	1·A·261·MET·CE	2.39	0.52			
1:B:50:GLU:HA	3:B:544:HOH:O	2.09	0.52			
1:C:50:GLU:OE1	1·C·105·ABG·NE	2.35	0.51			
1:A:69:SEB:HB3	1:C:64:HIS:CD2	2.44	0.51			
1:A:174:GLN:HG2	1:A:201:MET:HE1	1.92	0.51			
1:B:275:THR:O	1:B:279:LYS:NZ	2.36	0.51			
1:C:235:LEU:HD12	1:C:243:VAL:HG22	1.91	0.51			
1:A:247:PHE:N	1:A:247:PHE:CD1	2.79	0.51			
1:C:233:GLU:H	1:C:233:GLU:CD	2.13	0.51			
1:A:247:PHE:N	1:A:247:PHE:HD1	2.09	0.51			
1:B:50:GLU:O	1:B:50:GLU:HG3	2.09	0.51			
1:B:114:LEU:HD12	1:B:114:LEU:O	2.11	0.51			
1:C:164:LYS:HE2	1:C:165:TYR:OH	2.11	0.50			
1:C:256:ASN:H	1:C:256:ASN:ND2	2.05	0.50			
1:A:187:ILE:HG22	1:A:247:PHE:CE2	2.47	0.50			
1:B:75:ASP:OD1	1:B:76:SER:N	2.45	0.50			
1:A:31:LYS:HG3	1:A:294:PHE:CD2	2.47	0.50			
1:C:131:SER:O	1:C:131:SER:OG	2.29	0.50			
1:C:231:GLN:NE2	1:C:231:GLN:HA	2.27	0.49			
1:A:45:ALA:HB2	1:A:99:VAL:HG11	1.94	0.49			
1:B:39:LEU:HD23	1:B:305:ASN:HD21	1.77	0.49			
1:A:156:HIS:ND1	1:A:185:TYR:OH	2.32	0.49			
1:C:16:SER:O	1:C:20:LEU:HG	2.12	0.49			
1:C:261:MET:O	1:C:282:HIS:HD2	1.96	0.49			
1:B:12:ILE:CD1	1:B:15:LEU:HD12	2.38	0.48			
1:C:231:GLN:HB3	1:C:234:ARG:HG3	1.95	0.48			
1:A:196:GLN:HE21	1:A:200:ASP:CG	2.17	0.48			
1:B:59:PHE:O	1:B:63:MET:HG3	2.13	0.48			
1:B:162:ASP:OD2	1:B:165:TYR:HB2	2.14	0.48			
1:B:161:GLY:HA3	1:B:228:THR:OG1	2.14	0.48			
1:B:188:ALA:HB1	1:B:189:PRO:CD	2.44	0.48			
1:C:110:GLY:HA2	3:C:332:HOH:O	2.14	0.48			
1:A:163:LEU:O	1:A:195:PRO:HD3	2.14	0.47			
1:C:59:PHE:O	1:C:63:MET:HG3	2.14	0.47			
1:C:75:ASP:O	1:C:79:THR:HG23	2.14	0.47			
1:B:96:SER:OG	1:B:119:SER:HA	2.14	0.47			
1:C:164:LYS:HB2	1:C:192:LEU:HA	1.97	0.47			
1:A:153:ASP:HB2	1:A:180:ASP:O	2.14	0.47			
1:C:146:GLN:HB2	1:C:152:LEU:HG	1.97	0.47			
1:B:270:VAL:CG1	1:B:271:ASP:N	2.77	0.46			
1:A:186:PHE:CE1	1:A:194:MET:CE	2.99	0.46			



		Interatomic	Clash	
Atom-1	Atom-2	distance $(Å)$	overlap (Å)	
1:C:269:ARG:O	1:C:269:ARG:HG3	2.15	0.46	
1:A:36:PRO:HD2	3:A:386:HOH:O	2.15	0.46	
1:C:61:THR:O	1:C:65:ARG:HG2	2.15	0.46	
1:C:53:THR:O	1:C:57:LEU:HG	2.15	0.46	
1:B:254:LEU:HD13	1:B:282:HIS:CD2	2.51	0.46	
1:B:309:VAL:O	1:B:310:LEU:C	2.54	0.46	
1:A:54:ARG:HB2	1:A:54:ARG:NH1	2.30	0.46	
1:A:254:LEU:HD22	1:A:261:MET:HE1	1.98	0.46	
1:C:98:TYR:HB2	1:C:99:VAL:HG13	1.98	0.46	
1:C:106:HIS:ND1	1:C:107:PRO:HD2	2.30	0.46	
1:C:231:GLN:NE2	1:C:231:GLN:CA	2.79	0.45	
1:A:187:ILE:HG22	1:A:247:PHE:HE2	1.82	0.45	
1:B:132:ASN:HA	1:B:167:ARG:HE	1.82	0.45	
1:B:48:PHE:HB2	1:B:74:SER:HA	1.97	0.45	
1:B:36:PRO:HD2	3:B:486:HOH:O	2.16	0.45	
1:B:293:ILE:O	1:B:297:GLN:HB2	2.17	0.45	
1:B:50:GLU:OE2	1:B:105:ARG:NE	2.49	0.45	
1:B:31:LYS:NZ	1:B:291:ASN:HD21	2.16	0.44	
1:A:219:MET:HE2	1:A:257:ALA:CB	2.47	0.44	
1:A:257:ALA:HB1	1:A:261:MET:HE3	1.97	0.44	
1:A:265:HIS:ND1	1:A:266:PRO:HD2	2.32	0.44	
1:C:256:ASN:HD22	1:C:256:ASN:N	1.97	0.44	
1:A:54:ARG:HH11	1:A:54:ARG:CB	2.30	0.44	
1:A:218:VAL:O	1:A:219:MET:C	2.55	0.44	
1:A:219:MET:HE3	1:A:222:VAL:CG2	2.46	0.44	
1:A:1:ALA:HA	1:A:307:ASP:OD1	2.18	0.43	
1:A:164:LYS:HE2	1:A:165:TYR:CZ	2.52	0.43	
1:A:214:SER:O	1:A:217:GLU:HG3	2.17	0.43	
1:A:16:SER:O	1:A:19:ASP:HB2	2.17	0.43	
1:A:154:ASN:HA	1:A:181:GLY:O	2.18	0.43	
1:B:35:GLN:HB3	1:B:38:LEU:HB2	1.99	0.43	
1:A:121:ASN:OD1	1:A:121:ASN:N	2.29	0.43	
1:A:293:ILE:HG22	1:A:297:GLN:HE21	1.82	0.43	
1:A:84:LYS:HB3	1:A:86:GLU:OE1	2.19	0.43	
1:A:106:HIS:CE1	1:A:107:PRO:HD2	2.53	0.43	
1:A:17:ARG:NH2	1:A:178:LYS:O	2.52	0.43	
1:A:160:VAL:HG22	1:A:187:ILE:HB	2.01	0.43	
1:C:12:ILE:HD13	1:C:15:LEU:HD12	2.00	0.43	
1:A:172:LEU:HG	1:A:176:LEU:HD12	2.00	0.43	
1:C:125:LEU:HD23	1:C:299:LEU:CD2	2.49	0.42	
1:A:189:PRO:HG2	1:A:192:LEU:HD12	2.00	0.42	



Interatomic Clash					
Atom-1	Atom-2	distance (Å)	overlap (Å)		
1:B:199:LEU:HD13	1:B:209:TRP:CH2	2.54	0.42		
1:B:254:LEU:HD13	1:B:282:HIS:HD2	1.84	0.42		
1:B:270:VAL:CG1	1:B:271:ASP:H	2.21	0.42		
1:C:142:LEU:HD11	1:C:175:ALA:HB1	2.00	0.42		
1:B:250:ARG:O	1:B:253:ASP:HB2	2.19	0.42		
1:C:112:ALA:O	1:C:116:THR:HG23	2.20	0.42		
1:A:113:ARG:HD3	1:A:113:ARG:HA	1.92	0.42		
1:C:51:ALA:HB2	1:C:75:ASP:HB2	2.01	0.42		
1:B:56:ARG:O	1:B:60:GLU:HG3	2.20	0.42		
1:A:60:GLU:O	1:A:63:MET:HB2	2.19	0.42		
1:B:39:LEU:CD2	1:B:305:ASN:ND2	2.83	0.42		
1:C:59:PHE:O	1:C:62:SER:HB2	2.19	0.42		
1:C:20:LEU:O	1:C:24:LEU:HG	2.19	0.42		
1:C:54:ARG:NE	3:C:578:HOH:O	2.53	0.42		
1:A:33:ASN:O	1:A:35:GLN:NE2	2.47	0.41		
1:C:279:LYS:N	1:C:279:LYS:CE	2.79	0.41		
1:A:291:ASN:HD22	1:A:291:ASN:HA	1.59	0.41		
1:A:4:LEU:HD22	1:A:9:ILE:HD11	2.03	0.41		
1:A:54:ARG:NH1	1:A:54:ARG:CB	2.83	0.41		
1:A:164:LYS:CE	1:A:165:TYR:CZ	3.03	0.41		
1:A:258:LYS:NZ	3:A:424:HOH:O	2.53	0.41		
1:A:156:HIS:HB3	1:A:185:TYR:HE1	1.85	0.41		
1:C:121:ASN:H	1:C:121:ASN:ND2	2.12	0.41		
1:C:167:ARG:HA	1:C:167:ARG:HE	1.85	0.41		
1:B:8:HIS:CE1	1:B:123:PRO:HA	2.55	0.41		
1:A:4:LEU:HD22	1:A:9:ILE:CD1	2.50	0.41		
1:A:105:ARG:HD3	1:A:128:GLY:HA3	2.02	0.41		
1:A:216:GLU:H	1:A:216:GLU:HG3	1.57	0.41		
1:A:105:ARG:NH2	3:A:525:HOH:O	2.30	0.41		
1:A:219:MET:CE	1:A:222:VAL:CG2	2.99	0.41		
1:B:267:LEU:HB3	1:B:268:PRO:HA	2.03	0.41		
1:C:236:ASP:OD1	1:C:237:PRO:N	2.53	0.41		
1:A:267:LEU:CB	1:A:268:PRO:HA	2.42	0.40		
1:B:12:ILE:HD12	1:B:12:ILE:HA	1.77	0.40		
1:B:132:ASN:CG	1:B:133:GLN:H	2.25	0.40		
1:B:309:VAL:O	1:B:309:VAL:HG12	2.20	0.40		
1:C:56:ARG:HD2	3:C:415:HOH:O	2.22	0.40		
1:A:12:ILE:HG22	3:A:400:HOH:O	2.22	0.40		
1:A:84:LYS:HA	1:A:84:LYS:HD3	1.92	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	Perce	entiles
1	А	282/310~(91%)	268~(95%)	14~(5%)	0	100	100
1	В	283/310~(91%)	269~(95%)	12 (4%)	2(1%)	22	11
1	С	300/310~(97%)	290~(97%)	9~(3%)	1 (0%)	41	30
All	All	865/930~(93%)	827 (96%)	35~(4%)	3~(0%)	41	30

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	132	ASN
1	В	270	VAL
1	С	270	VAL

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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percei	ntiles
1	А	238/261~(91%)	219~(92%)	19 (8%)	12	4
1	В	239/261~(92%)	227~(95%)	12 (5%)	24	13
1	С	$256/261 \ (98\%)$	241 (94%)	15 (6%)	19	9
All	All	733/783~(94%)	687 (94%)	46 (6%)	18	7

All (46) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	А	3	PRO
1	А	11	SER
1	А	16	SER
1	А	29	LYS
1	А	50	GLU
1	А	59	PHE
1	А	86	GLU
1	А	134	HIS
1	А	171	SER
1	А	183	ARG
1	А	190	ASP
1	А	201	MET
1	А	217	GLU
1	A	247	PHE
1	А	250	ARG
1	А	269	ARG
1	А	271	ASP
1	А	285	TYR
1	А	306	ARG
1	В	59	PHE
1	В	132	ASN
1	В	134	HIS
1	В	147	GLU
1	В	167	ARG
1	В	258	LYS
1	В	272	GLU
1	В	279	LYS
1	В	285	TYR
1	В	297	GLN
1	В	306	ARG
1	В	310	LEU
1	С	59	PHE
1	С	78	ASN
1	С	86	GLU
1	С	98	TYR
1	С	121	ASN
1	С	131	SER
1	С	167	ARG
1	C	171	SER
1	C	174	GLN
1	C	234	ARG
1	C	256	ASN
1	С	271	ASP



Continued from previous page...

Mol	Chain	Res	Type
1	С	279	LYS
1	С	285	TYR
1	С	307	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	8	HIS
1	А	108	GLN
1	А	196	GLN
1	А	282	HIS
1	А	291	ASN
1	В	33	ASN
1	В	291	ASN
1	В	305	ASN
1	С	8	HIS
1	С	13	ASN
1	С	64	HIS
1	С	121	ASN
1	С	137	GLN
1	С	146	GLN
1	С	174	GLN
1	С	182	ASN
1	С	231	GLN
1	С	255	HIS
1	С	256	ASN
1	С	282	HIS

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, 2 are monoatomic - leaving 0 for Mogul analysis. 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. No monomer is involved in short contacts.

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.

