



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:37 PM UTC

PDB ID : 1BKO / pdb_00001bko
Title : THERMOSTABLE THYMIDYLATE SYNTHASE A FROM BACILLUS SUBTILIS
Authors : Stout, T.J.; Schellenberger, U.; Santi, D.V.; Stroud, R.M.
Deposited on : 1998-07-09
Resolution : 2.75 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

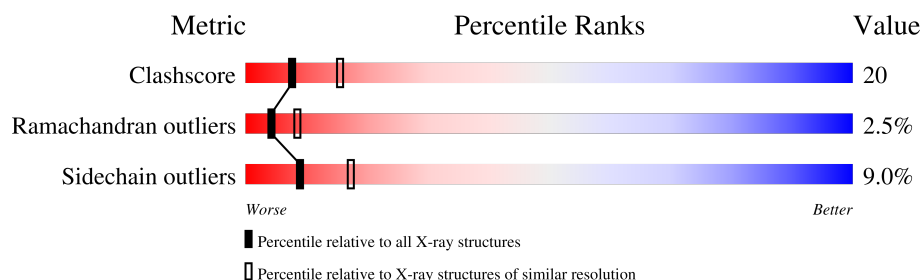
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	278	53% 37% 8% .
1	B	278	54% 38% 6% .
1	C	278	56% 39% . .
1	D	278	52% 41% 5% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9239 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 THYMIDYLATE SYNTHASE A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	0	0
			2301	1471	390	430	10			
1	B	278	Total	C	N	O	S	0	0	0
			2301	1471	390	430	10			
1	C	278	Total	C	N	O	S	0	0	0
			2301	1471	390	430	10			
1	D	278	Total	C	N	O	S	0	0	0
			2301	1471	390	430	10			

- Molecule 2 is water.

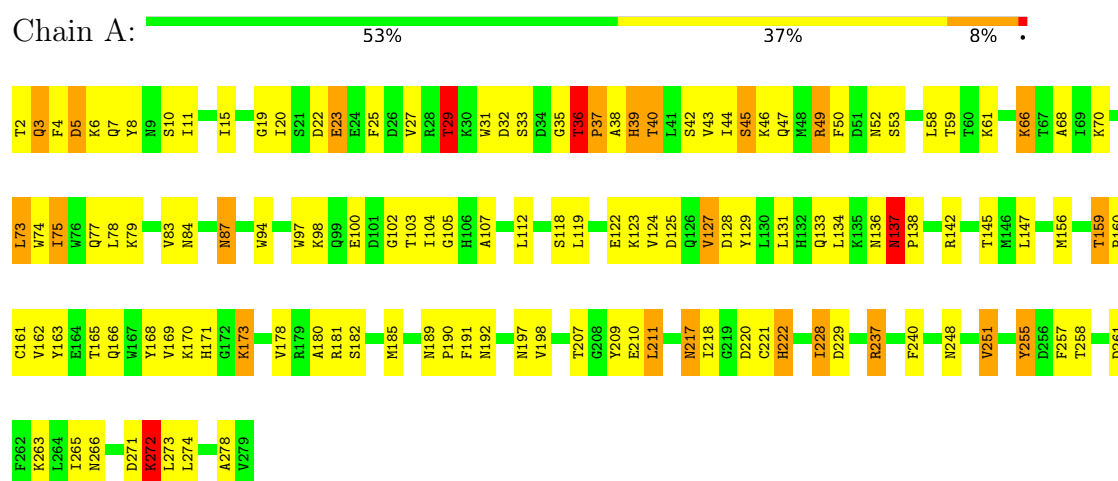
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	35	Total	O	0	0
			35	35		

3 Residue-property plots

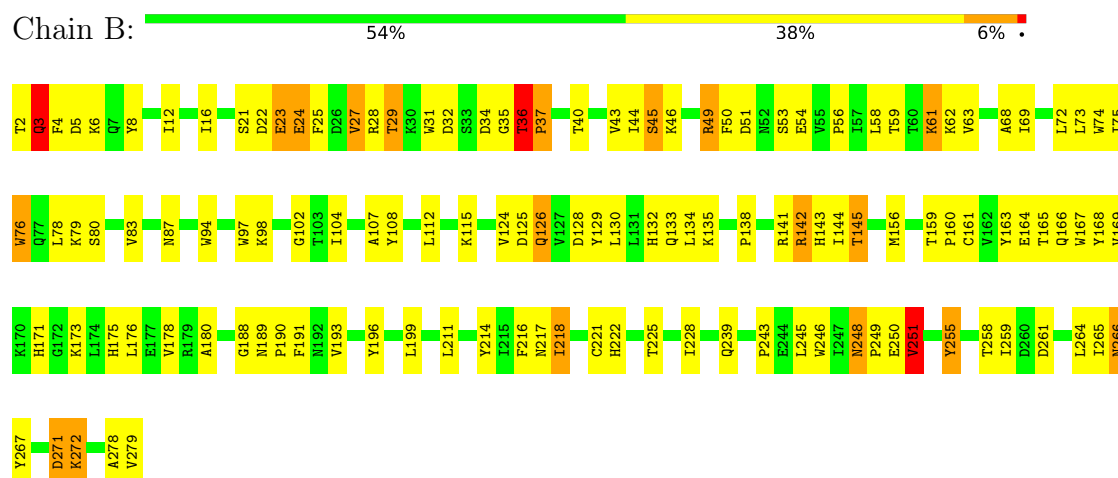
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

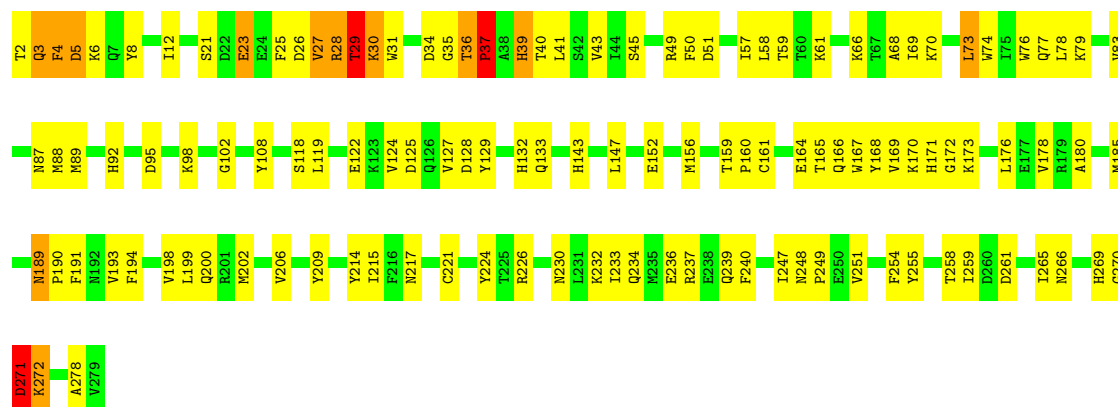


• Molecule 1: THYMIDYLATE SYNTHASE A



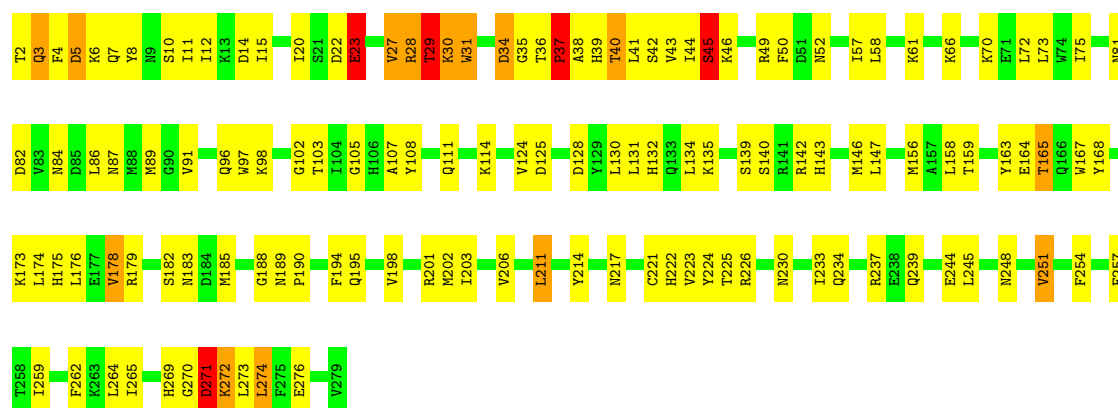
• Molecule 1: THYMIDYLATE SYNTHASE A





• Molecule 1: THYMIDYLATE SYNTHASE A

Chain D: 52% 41% 5%



4 Data and refinement statistics

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	99.36Å 52.83Å 129.39Å 90.00° 101.70° 90.00°	Depositor
Resolution (Å)	50.00 – 2.75	Depositor
% Data completeness (in resolution range)	76.1 (50.00-2.75)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
Refinement program	X-PLOR 3.854	Depositor
R, R_{free}	0.213 , 0.320	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9239	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	0.42	0/2359	1.00	17/3196 (0.5%)
1	B	0.44	0/2359	1.04	18/3196 (0.6%)
1	C	0.44	0/2359	0.96	10/3196 (0.3%)
1	D	0.45	0/2359	0.99	14/3196 (0.4%)
All	All	0.44	0/9436	1.00	59/12784 (0.5%)

There are no bond length outliers.

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	188	GLY	N-CA-C	10.14	126.75	113.27
1	B	188	GLY	N-CA-C	9.71	124.38	112.73
1	B	29	THR	N-CA-C	8.04	120.70	110.24
1	C	28	ARG	N-CA-C	7.91	120.33	108.31
1	B	76	TRP	N-CA-C	7.76	119.52	111.14
1	B	36	THR	CA-C-N	7.38	129.07	119.84
1	B	36	THR	C-N-CA	7.38	129.07	119.84
1	A	5	ASP	N-CA-C	-7.36	102.42	111.40
1	B	222	HIS	N-CA-C	7.16	119.81	109.14
1	B	59	THR	N-CA-C	-6.79	105.93	114.56
1	A	36	THR	CA-C-N	6.77	128.30	119.84
1	A	36	THR	C-N-CA	6.77	128.30	119.84
1	B	104	ILE	N-CA-C	-6.74	104.74	111.88
1	A	40	THR	N-CA-C	6.66	121.13	109.80
1	D	75	ILE	N-CA-C	6.53	117.15	110.82
1	C	40	THR	N-CA-C	6.51	120.86	109.80
1	A	127	VAL	N-CA-C	6.49	116.52	110.42
1	C	76	TRP	N-CA-C	6.42	117.97	110.97
1	A	19	GLY	N-CA-C	6.29	120.78	110.97
1	C	29	THR	N-CA-C	6.28	124.18	110.80
1	A	53	SER	N-CA-C	-6.23	104.40	111.07
1	B	246	TRP	N-CA-C	6.13	118.61	109.59
1	B	164	GLU	N-CA-C	6.13	118.17	108.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	ASN	N-CA-C	-6.10	105.33	112.89
1	B	248	ASN	CA-C-N	6.09	125.98	119.28
1	B	248	ASN	C-N-CA	6.09	125.98	119.28
1	D	37	PRO	CA-N-CD	-6.08	103.49	112.00
1	A	44	ILE	N-CA-C	5.96	117.94	108.89
1	A	229	ASP	N-CA-C	5.94	117.45	110.97
1	D	159	THR	CA-C-N	5.84	125.61	119.76
1	D	159	THR	C-N-CA	5.84	125.61	119.76
1	D	40	THR	N-CA-C	5.83	115.98	108.34
1	B	32	ASP	N-CA-C	-5.83	106.51	113.97
1	A	198	VAL	N-CA-C	-5.80	104.68	111.00
1	B	266	ASN	N-CA-C	5.79	117.68	110.91
1	B	141	ARG	N-CA-C	-5.75	105.57	112.59
1	D	164	GLU	N-CA-C	5.74	118.13	109.23
1	D	244	GLU	N-CA-C	-5.74	102.80	110.55
1	D	146	MET	N-CA-C	5.71	117.72	108.41
1	D	29	THR	N-CA-C	5.70	122.94	110.80
1	A	192	ASN	N-CA-C	5.68	117.48	111.28
1	A	222	HIS	N-CA-C	5.67	119.20	110.42
1	A	75	ILE	N-CA-C	5.61	116.67	111.45
1	C	37	PRO	CA-N-CD	-5.58	104.19	112.00
1	D	5	ASP	N-CA-C	-5.56	104.85	111.69
1	D	37	PRO	N-CA-C	5.45	123.70	112.47
1	B	251	VAL	N-CA-C	5.41	116.68	109.37
1	A	272	LYS	N-CA-C	5.39	118.27	109.59
1	B	126	GLN	N-CA-C	5.37	116.94	111.14
1	C	28	ARG	CB-CA-C	-5.37	110.36	116.54
1	A	32	ASP	N-CA-C	-5.35	107.30	113.88
1	A	29	THR	N-CA-C	5.32	122.12	110.80
1	B	3	GLN	N-CA-C	5.26	118.86	112.23
1	D	274	LEU	N-CA-C	5.24	117.44	108.90
1	C	164	GLU	N-CA-C	5.19	116.96	109.07
1	C	34	ASP	N-CA-C	-5.13	107.57	113.88
1	D	31	TRP	N-CA-C	5.12	116.93	110.33
1	C	266	ASN	N-CA-C	5.03	116.80	110.91
1	C	152	GLU	N-CA-C	5.00	118.69	112.59

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	2301	0	2226	86	0
1	B	2301	0	2226	88	0
1	C	2301	0	2226	107	0
1	D	2301	0	2226	95	0
2	A	35	0	0	0	0
All	All	9239	0	8904	364	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 20.

All (364) 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:B:272:LYS:HE3	1:B:272:LYS:HA	1.48	0.96
1:D:30:LYS:HB2	1:D:36:THR:HA	1.56	0.87
1:C:43:VAL:HG23	1:C:221:CYS:HB3	1.56	0.87
1:C:36:THR:OG1	1:C:37:PRO:HD2	1.75	0.86
1:A:49:ARG:HH22	1:B:45:SER:HB2	1.39	0.85
1:B:74:TRP:HA	1:B:78:LEU:HD12	1.60	0.83
1:B:62:LYS:HB3	1:B:272:LYS:HZ3	1.43	0.82
1:C:165:THR:HG22	1:C:178:VAL:HG22	1.61	0.82
1:C:230:ASN:HA	1:C:233:ILE:HD12	1.61	0.81
1:A:272:LYS:NZ	1:A:272:LYS:HA	1.97	0.79
1:D:183:ASN:HD21	1:D:189:ASN:HA	1.48	0.79
1:C:77:GLN:HE21	1:C:255:TYR:HA	1.50	0.76
1:C:199:LEU:HA	1:C:202:MET:HE3	1.69	0.75
1:D:201:ARG:HG2	1:D:211:LEU:HD21	1.68	0.75
1:A:3:GLN:HB2	1:A:6:LYS:HD3	1.69	0.74
1:C:27:VAL:HG12	1:C:28:ARG:H	1.52	0.74
1:D:2:THR:HB	1:D:50:PHE:HD1	1.54	0.72
1:A:23:GLU:HA	1:A:39:HIS:NE2	2.05	0.72
1:A:11:ILE:O	1:A:15:ILE:HG13	1.90	0.72
1:D:36:THR:OG1	1:D:37:PRO:HD2	1.89	0.71
1:B:62:LYS:HB3	1:B:272:LYS:NZ	2.06	0.70
1:A:147:LEU:HD12	1:A:163:TYR:HA	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:THR:OG1	1:A:37:PRO:HD2	1.92	0.70
1:C:74:TRP:HA	1:C:78:LEU:HD12	1.74	0.69
1:A:49:ARG:NH2	1:B:45:SER:HB2	2.06	0.69
1:C:3:GLN:C	1:C:5:ASP:H	2.01	0.69
1:C:254:PHE:HD2	1:C:255:TYR:CD1	2.11	0.68
1:A:20:ILE:HG23	1:B:171:HIS:NE2	2.09	0.68
1:C:83:VAL:HG22	1:C:102:GLY:O	1.94	0.68
1:D:82:ASP:HA	1:D:103:THR:HG22	1.76	0.68
1:A:237:ARG:NH2	1:A:273:LEU:HG	2.10	0.67
1:C:30:LYS:HB2	1:C:36:THR:HA	1.75	0.67
1:A:207:THR:HG1	1:A:209:TYR:HD2	1.40	0.67
1:A:166:GLN:HG2	1:A:168:TYR:CZ	2.30	0.67
1:A:79:LYS:HZ3	1:A:127:VAL:HB	1.58	0.66
1:C:124:VAL:HG22	1:C:125:ASP:H	1.60	0.66
1:D:189:ASN:HB3	1:D:190:PRO:HD3	1.76	0.66
1:C:2:THR:HG21	1:C:51:ASP:H	1.59	0.66
1:A:31:TRP:O	1:A:35:GLY:HA2	1.96	0.66
1:D:36:THR:CB	1:D:37:PRO:HD2	2.25	0.66
1:C:3:GLN:HG3	1:C:6:LYS:NZ	2.09	0.66
1:B:58:LEU:CD2	1:B:190:PRO:HB3	2.25	0.65
1:D:36:THR:HG23	1:D:37:PRO:HD2	1.77	0.65
1:C:5:ASP:HB3	1:C:239:GLN:OE1	1.97	0.65
1:D:131:LEU:HA	1:D:134:LEU:HD12	1.77	0.65
1:B:3:GLN:HG3	1:B:6:LYS:NZ	2.12	0.65
1:D:70:LYS:HE2	1:D:89:MET:HB3	1.79	0.65
1:D:165:THR:HG22	1:D:176:LEU:HD11	1.79	0.64
1:C:198:VAL:O	1:C:202:MET:HG3	1.96	0.64
1:A:165:THR:HG22	1:A:178:VAL:HG22	1.80	0.64
1:C:234:GLN:HA	1:C:237:ARG:HD3	1.80	0.64
1:B:98:LYS:HG3	1:B:102:GLY:HA2	1.80	0.64
1:C:272:LYS:NZ	1:C:272:LYS:HA	2.13	0.64
1:A:272:LYS:HA	1:A:272:LYS:HZ2	1.61	0.63
1:D:36:THR:CG2	1:D:37:PRO:HD2	2.28	0.63
1:D:43:VAL:HG23	1:D:221:CYS:HB2	1.80	0.63
1:D:23:GLU:HG3	1:D:39:HIS:CG	2.34	0.63
1:B:225:THR:O	1:B:228:ILE:HG12	1.99	0.63
1:C:29:THR:HG21	1:C:224:TYR:HE1	1.64	0.63
1:D:147:LEU:HD12	1:D:163:TYR:HA	1.80	0.63
1:A:124:VAL:CG2	1:A:128:ASP:HB2	2.29	0.62
1:C:36:THR:CB	1:C:37:PRO:HD2	2.30	0.62
1:D:3:GLN:HB2	1:D:6:LYS:HD3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:254:PHE:O	1:D:257:PHE:HD1	1.83	0.62
1:D:183:ASN:ND2	1:D:189:ASN:HA	2.15	0.61
1:B:218:ILE:N	1:B:218:ILE:HD12	2.15	0.61
1:B:248:ASN:HD21	1:B:250:GLU:HB2	1.64	0.61
1:A:22:ASP:OD1	1:A:39:HIS:HB2	2.00	0.61
1:D:11:ILE:O	1:D:15:ILE:HG13	2.01	0.60
1:B:189:ASN:O	1:B:193:VAL:HG23	2.01	0.60
1:A:79:LYS:NZ	1:A:127:VAL:HB	2.17	0.60
1:B:58:LEU:HD21	1:B:190:PRO:HB3	1.83	0.60
1:A:31:TRP:CE3	1:A:278:ALA:HB2	2.37	0.60
1:B:31:TRP:O	1:B:35:GLY:HA2	2.00	0.60
1:C:2:THR:HB	1:C:50:PHE:HD1	1.67	0.60
1:C:199:LEU:HA	1:C:202:MET:CE	2.32	0.60
1:A:23:GLU:HA	1:A:39:HIS:CD2	2.37	0.60
1:C:98:LYS:CG	1:C:102:GLY:HA2	2.32	0.59
1:B:51:ASP:OD1	1:B:53:SER:HB2	2.02	0.59
1:B:98:LYS:CG	1:B:102:GLY:HA2	2.33	0.59
1:C:269:HIS:HD2	1:C:270:GLY:O	1.85	0.59
1:A:43:VAL:HG23	1:A:221:CYS:HB3	1.85	0.58
1:C:27:VAL:HG12	1:C:28:ARG:N	2.16	0.58
1:C:167:TRP:CH2	1:C:176:LEU:HD22	2.39	0.58
1:C:69:ILE:O	1:C:73:LEU:HD22	2.03	0.57
1:D:66:LYS:O	1:D:70:LYS:HG3	2.04	0.57
1:D:82:ASP:OD2	1:D:84:ASN:HB3	2.04	0.57
1:D:140:SER:HB3	1:D:143:HIS:CE1	2.39	0.57
1:A:128:ASP:HA	1:A:131:LEU:HD12	1.86	0.57
1:B:2:THR:O	1:B:54:GLU:HB2	2.05	0.57
1:D:179:ARG:HG3	1:D:217:ASN:HB2	1.87	0.57
1:A:2:THR:HB	1:A:50:PHE:HD2	1.70	0.57
1:C:23:GLU:HG3	1:C:39:HIS:CD2	2.40	0.57
1:D:72:LEU:HB2	1:D:195:GLN:HG3	1.87	0.57
1:A:3:GLN:O	1:A:4:PHE:HB3	2.06	0.56
1:A:182:SER:CB	1:A:220:ASP:HB3	2.35	0.56
1:C:272:LYS:HA	1:C:272:LYS:HZ2	1.70	0.56
1:C:156:MET:HE1	1:C:160:PRO:HD3	1.88	0.55
1:D:165:THR:HG23	1:D:178:VAL:HG13	1.88	0.55
1:A:29:THR:HB	1:A:38:ALA:O	2.06	0.55
1:B:83:VAL:HG21	1:B:98:LYS:HE2	1.88	0.55
1:D:3:GLN:C	1:D:5:ASP:H	2.14	0.55
1:C:59:THR:HB	1:C:240:PHE:H	1.71	0.55
1:C:29:THR:HG21	1:C:224:TYR:CE1	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:ILE:N	1:B:265:ILE:HD12	2.22	0.54
1:D:4:PHE:HE2	1:D:58:LEU:HD22	1.71	0.54
1:B:36:THR:OG1	1:B:37:PRO:HD2	2.07	0.54
1:D:45:SER:O	1:D:46:LYS:HD2	2.08	0.54
1:B:2:THR:HB	1:B:50:PHE:HD1	1.73	0.54
1:D:131:LEU:HD21	1:D:203:ILE:HD12	1.90	0.54
1:A:98:LYS:HG3	1:A:102:GLY:HA2	1.89	0.54
1:D:29:THR:HG21	1:D:224:TYR:HE2	1.72	0.53
1:B:63:VAL:HG23	1:B:267:TYR:OH	2.09	0.53
1:C:166:GLN:HG2	1:C:168:TYR:CE1	2.44	0.53
1:B:3:GLN:HG3	1:B:6:LYS:HZ2	1.74	0.53
1:C:3:GLN:HG3	1:C:6:LYS:HZ3	1.73	0.53
1:B:72:LEU:HD12	1:B:76:TRP:CD1	2.44	0.53
1:A:3:GLN:C	1:A:5:ASP:H	2.17	0.52
1:B:83:VAL:HG22	1:B:102:GLY:O	2.09	0.52
1:D:198:VAL:HG12	1:D:202:MET:HE3	1.91	0.52
1:B:255:TYR:CD1	1:B:255:TYR:N	2.77	0.52
1:A:8:TYR:HA	1:A:11:ILE:HD12	1.91	0.52
1:C:43:VAL:CG2	1:C:221:CYS:HB3	2.34	0.52
1:C:172:GLY:O	1:C:209:TYR:HD1	1.93	0.52
1:C:98:LYS:HG2	1:C:102:GLY:HA2	1.91	0.52
1:C:248:ASN:O	1:C:251:VAL:HG23	2.08	0.52
1:B:74:TRP:NE1	1:B:80:SER:HB3	2.25	0.52
1:B:255:TYR:N	1:B:255:TYR:HD1	2.06	0.52
1:D:57:ILE:HD12	1:D:194:PHE:CE1	2.45	0.52
1:D:98:LYS:HG3	1:D:102:GLY:HA2	1.91	0.52
1:A:272:LYS:HA	1:A:272:LYS:HZ1	1.74	0.51
1:D:245:LEU:HD13	1:D:264:LEU:HD12	1.93	0.51
1:B:168:TYR:HB2	1:B:175:HIS:HB2	1.90	0.51
1:C:98:LYS:HG3	1:C:102:GLY:HA2	1.92	0.51
1:C:124:VAL:HG22	1:C:128:ASP:HB2	1.93	0.51
1:C:170:LYS:HE3	1:D:42:SER:HB2	1.92	0.51
1:B:5:ASP:HB2	1:B:239:GLN:OE1	2.11	0.51
1:D:36:THR:HG23	1:D:37:PRO:CD	2.41	0.51
1:A:161:CYS:O	1:A:180:ALA:HA	2.11	0.51
1:D:234:GLN:HA	1:D:237:ARG:HD3	1.92	0.51
1:D:97:TRP:NE1	1:D:158:LEU:HD13	2.26	0.51
1:B:43:VAL:CG2	1:B:221:CYS:HB3	2.40	0.51
1:D:124:VAL:HG22	1:D:125:ASP:N	2.25	0.51
1:C:108:TYR:HA	1:C:147:LEU:HD22	1.92	0.51
1:C:23:GLU:CD	1:C:23:GLU:H	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:ARG:HA	1:D:214:TYR:O	2.11	0.50
1:D:132:HIS:HA	1:D:135:LYS:NZ	2.25	0.50
1:C:200:GLN:HE22	1:C:214:TYR:HB2	1.77	0.50
1:A:129:TYR:CE1	1:A:133:GLN:HG3	2.46	0.50
1:C:215:ILE:HD12	1:C:215:ILE:N	2.27	0.50
1:C:79:LYS:NZ	1:C:127:VAL:HB	2.27	0.50
1:C:247:ILE:O	1:C:249:PRO:HD3	2.11	0.50
1:B:75:ILE:O	1:B:79:LYS:HA	2.11	0.50
1:C:124:VAL:HG22	1:C:125:ASP:N	2.25	0.50
1:C:265:ILE:N	1:C:265:ILE:HD12	2.27	0.50
1:D:20:ILE:O	1:D:41:LEU:HA	2.12	0.50
1:A:52:ASN:HD22	1:A:211:LEU:HG	1.76	0.49
1:B:189:ASN:HB3	1:B:190:PRO:HD3	1.93	0.49
1:D:182:SER:HB2	1:D:222:HIS:HE1	1.77	0.49
1:D:230:ASN:HA	1:D:233:ILE:HD12	1.94	0.49
1:D:5:ASP:HB3	1:D:239:GLN:NE2	2.26	0.49
1:A:83:VAL:HG22	1:A:102:GLY:O	2.12	0.49
1:D:29:THR:HB	1:D:38:ALA:HB3	1.95	0.49
1:D:86:LEU:HG	1:D:91:VAL:HB	1.95	0.49
1:D:272:LYS:HG3	1:D:273:LEU:N	2.26	0.49
1:B:27:VAL:HG12	1:B:28:ARG:N	2.28	0.49
1:C:49:ARG:HA	1:C:214:TYR:O	2.13	0.49
1:A:107:ALA:HB2	1:A:156:MET:HB3	1.94	0.49
1:B:132:HIS:HA	1:B:135:LYS:HE3	1.95	0.49
1:D:31:TRP:CZ3	1:D:276:GLU:HB2	2.48	0.49
1:C:3:GLN:HG3	1:C:6:LYS:HZ2	1.75	0.49
1:C:8:TYR:O	1:C:12:ILE:HG12	2.12	0.49
1:A:255:TYR:CD1	1:A:255:TYR:N	2.81	0.49
1:C:166:GLN:HG2	1:C:168:TYR:CZ	2.48	0.49
1:D:5:ASP:HB3	1:D:239:GLN:HE21	1.78	0.49
1:D:269:HIS:HD2	1:D:270:GLY:O	1.94	0.49
1:B:248:ASN:O	1:B:251:VAL:HG23	2.13	0.48
1:C:189:ASN:HB3	1:C:190:PRO:HD3	1.94	0.48
1:C:230:ASN:O	1:C:233:ILE:HB	2.13	0.48
1:A:138:PRO:O	1:A:169:VAL:HB	2.13	0.48
1:B:8:TYR:O	1:B:12:ILE:HG12	2.14	0.48
1:B:49:ARG:HA	1:B:214:TYR:O	2.12	0.48
1:D:202:MET:HE2	1:D:262:PHE:CE2	2.47	0.48
1:C:3:GLN:C	1:C:5:ASP:N	2.67	0.48
1:C:66:LYS:O	1:C:70:LYS:HG3	2.13	0.48
1:C:133:GLN:OE1	1:C:133:GLN:HA	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:111:GLN:HA	1:D:114:LYS:HG3	1.95	0.48
1:B:163:TYR:HD1	1:B:163:TYR:H	1.62	0.48
1:C:129:TYR:O	1:C:133:GLN:HG2	2.13	0.48
1:D:265:ILE:HD12	1:D:265:ILE:N	2.28	0.48
1:A:3:GLN:HG3	1:A:6:LYS:NZ	2.29	0.48
1:A:103:THR:C	1:A:105:GLY:H	2.22	0.48
1:A:3:GLN:CB	1:A:6:LYS:HD3	2.42	0.48
1:A:207:THR:OG1	1:A:209:TYR:HD2	1.96	0.48
1:D:43:VAL:CG2	1:D:221:CYS:HB2	2.42	0.47
1:A:77:GLN:HE22	1:A:257:PHE:HB2	1.79	0.47
1:A:181:ARG:NH1	1:B:142:ARG:HG3	2.28	0.47
1:B:29:THR:HA	1:B:279:VAL:OXT	2.14	0.47
1:D:124:VAL:CG2	1:D:128:ASP:HB2	2.45	0.47
1:B:23:GLU:C	1:B:25:PHE:H	2.21	0.47
1:C:45:SER:HB3	1:D:49:ARG:HH22	1.79	0.47
1:B:258:THR:HG22	1:B:261:ASP:OD2	2.15	0.47
1:A:66:LYS:O	1:A:70:LYS:HG3	2.14	0.47
1:C:4:PHE:HE2	1:C:58:LEU:HD22	1.80	0.47
1:C:161:CYS:O	1:C:180:ALA:HA	2.15	0.47
1:A:159:THR:HA	1:A:160:PRO:HD3	1.77	0.47
1:C:119:LEU:HD21	1:C:132:HIS:CG	2.50	0.47
1:A:47:GLN:HB2	1:A:217:ASN:HB3	1.97	0.47
1:A:145:THR:HB	1:A:165:THR:OG1	2.15	0.47
1:D:70:LYS:HA	1:D:89:MET:HE1	1.97	0.47
1:B:83:VAL:CG2	1:B:98:LYS:HE2	2.45	0.46
1:C:3:GLN:O	1:C:5:ASP:N	2.48	0.46
1:C:258:THR:HG22	1:C:261:ASP:CG	2.40	0.46
1:D:3:GLN:O	1:D:4:PHE:HB3	2.14	0.46
1:D:7:GLN:O	1:D:8:TYR:C	2.58	0.46
1:A:84:ASN:O	1:A:87:ASN:HB2	2.16	0.46
1:B:69:ILE:O	1:B:73:LEU:HD13	2.15	0.46
1:C:27:VAL:CG1	1:C:28:ARG:H	2.26	0.46
1:A:263:LYS:HE3	1:A:265:ILE:HD11	1.96	0.46
1:A:258:THR:HG22	1:A:261:ASP:OD2	2.15	0.46
1:B:124:VAL:CG2	1:B:128:ASP:HB3	2.46	0.46
1:A:59:THR:HB	1:A:240:PHE:HD1	1.81	0.46
1:C:124:VAL:HG11	1:C:129:TYR:HB2	1.97	0.46
1:C:185:MET:HE3	1:C:221:CYS:SG	2.55	0.46
1:D:2:THR:HB	1:D:50:PHE:CD1	2.43	0.46
1:D:105:GLY:C	1:D:107:ALA:H	2.23	0.46
1:B:161:CYS:O	1:B:180:ALA:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:59:THR:CB	1:C:240:PHE:H	2.29	0.46
1:B:68:ALA:HA	1:B:191:PHE:CZ	2.51	0.46
1:D:107:ALA:HA	1:D:156:MET:HE3	1.98	0.46
1:A:74:TRP:O	1:A:78:LEU:HB2	2.16	0.46
1:A:98:LYS:CG	1:A:102:GLY:HA2	2.46	0.46
1:B:159:THR:HA	1:B:160:PRO:HD3	1.85	0.46
1:D:274:LEU:HD12	1:D:274:LEU:HA	1.79	0.46
1:A:75:ILE:O	1:A:79:LYS:HA	2.16	0.45
1:A:248:ASN:O	1:A:251:VAL:HG23	2.15	0.45
1:A:255:TYR:N	1:A:255:TYR:HD1	2.13	0.45
1:C:124:VAL:CG2	1:C:128:ASP:HB2	2.46	0.45
1:C:36:THR:HG1	1:C:37:PRO:HD2	1.79	0.45
1:A:173:LYS:HD2	1:A:210:GLU:HB2	1.99	0.45
1:B:43:VAL:HG22	1:B:221:CYS:HB3	1.98	0.45
1:C:31:TRP:CE3	1:C:278:ALA:HB2	2.51	0.45
1:C:237:ARG:NH2	1:C:271:ASP:N	2.64	0.45
1:C:49:ARG:HB3	1:C:215:ILE:HG13	1.97	0.45
1:D:23:GLU:HG3	1:D:39:HIS:CD2	2.51	0.45
1:C:165:THR:CG2	1:C:178:VAL:HG22	2.40	0.45
1:D:28:ARG:O	1:D:30:LYS:HD3	2.17	0.45
1:D:248:ASN:O	1:D:251:VAL:HG23	2.16	0.45
1:A:94:TRP:HA	1:A:97:TRP:HZ3	1.82	0.45
1:A:118:SER:HA	1:A:122:GLU:O	2.16	0.45
1:C:68:ALA:HA	1:C:191:PHE:CZ	2.52	0.45
1:C:70:LYS:HB3	1:C:89:MET:HE2	1.99	0.45
1:A:45:SER:HB3	1:B:49:ARG:HH12	1.82	0.45
1:B:3:GLN:HG3	1:B:6:LYS:HZ3	1.79	0.45
1:D:34:ASP:CG	1:D:35:GLY:N	2.74	0.44
1:D:185:MET:HB2	1:D:223:VAL:HG22	2.00	0.44
1:B:167:TRP:CD2	1:B:176:LEU:HD13	2.52	0.44
1:C:23:GLU:HG3	1:C:39:HIS:CG	2.51	0.44
1:D:165:THR:CG2	1:D:176:LEU:HD11	2.44	0.44
1:B:34:ASP:OD2	1:B:36:THR:HG22	2.17	0.44
1:A:83:VAL:CG2	1:A:98:LYS:HE2	2.48	0.44
1:A:136:ASN:O	1:A:137:ASN:HB2	2.18	0.44
1:D:96:GLN:HB3	1:D:97:TRP:CE3	2.53	0.44
1:B:126:GLN:OE1	1:B:145:THR:HG22	2.17	0.44
1:A:68:ALA:HA	1:A:191:PHE:CZ	2.53	0.44
1:A:94:TRP:HA	1:A:97:TRP:CZ3	2.53	0.44
1:C:143:HIS:HB2	1:C:167:TRP:O	2.18	0.44
1:D:31:TRP:HB2	1:D:34:ASP:OD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:271:ASP:HB3	1:D:272:LYS:H	1.52	0.44
1:B:258:THR:O	1:B:261:ASP:HB2	2.17	0.43
1:C:21:SER:HB2	1:C:23:GLU:OE1	2.18	0.43
1:C:57:ILE:HD12	1:C:194:PHE:CE2	2.52	0.43
1:C:118:SER:HA	1:C:122:GLU:O	2.17	0.43
1:A:20:ILE:HG23	1:B:171:HIS:CE1	2.53	0.43
1:A:138:PRO:CB	1:A:169:VAL:HG11	2.48	0.43
1:B:107:ALA:HB2	1:B:156:MET:HG2	2.00	0.43
1:B:130:LEU:O	1:B:134:LEU:HG	2.18	0.43
1:A:3:GLN:HG3	1:A:6:LYS:HZ2	1.84	0.43
1:C:79:LYS:HE2	1:C:127:VAL:HG21	2.00	0.43
1:D:10:SER:O	1:D:14:ASP:HB2	2.18	0.43
1:A:40:THR:HB	1:A:222:HIS:HB2	2.01	0.43
1:A:168:TYR:CD2	1:B:44:ILE:HD13	2.54	0.43
1:B:129:TYR:CE1	1:B:133:GLN:HG3	2.54	0.43
1:D:27:VAL:HG12	1:D:28:ARG:N	2.33	0.43
1:D:29:THR:HG21	1:D:224:TYR:CE2	2.53	0.43
1:B:22:ASP:HB2	1:B:25:PHE:HD2	1.84	0.43
1:B:112:LEU:HD23	1:B:112:LEU:HA	1.79	0.43
1:B:3:GLN:C	1:B:5:ASP:H	2.27	0.43
1:B:21:SER:HA	1:B:40:THR:O	2.19	0.43
1:A:23:GLU:HG3	1:A:39:HIS:CE1	2.54	0.43
1:A:25:PHE:CE1	1:B:138:PRO:HG2	2.54	0.43
1:C:124:VAL:CG1	1:C:129:TYR:HB2	2.49	0.43
1:B:143:HIS:O	1:B:166:GLN:HA	2.19	0.42
1:B:248:ASN:HD22	1:B:251:VAL:CG2	2.32	0.42
1:B:124:VAL:HG22	1:B:125:ASP:N	2.34	0.42
1:C:70:LYS:HE2	1:C:89:MET:HB3	2.01	0.42
1:A:258:THR:O	1:A:261:ASP:HB2	2.19	0.42
1:C:23:GLU:OE2	1:C:39:HIS:HB2	2.19	0.42
1:C:59:THR:HG21	1:C:240:PHE:O	2.20	0.42
1:D:12:ILE:HD11	1:D:185:MET:HB3	2.02	0.42
1:D:132:HIS:HA	1:D:135:LYS:HZ2	1.83	0.42
1:B:115:LYS:HA	1:B:124:VAL:O	2.19	0.42
1:B:143:HIS:C	1:B:144:ILE:HG13	2.45	0.42
1:C:79:LYS:HZ3	1:C:127:VAL:HB	1.82	0.42
1:D:194:PHE:O	1:D:198:VAL:HG23	2.19	0.42
1:A:112:LEU:O	1:A:125:ASP:HB2	2.20	0.42
1:D:130:LEU:O	1:D:134:LEU:HG	2.19	0.42
1:B:94:TRP:HA	1:B:97:TRP:CZ3	2.55	0.42
1:B:138:PRO:O	1:B:169:VAL:HB	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:GLY:O	1:C:36:THR:HB	2.19	0.42
1:D:81:ASN:HD22	1:D:103:THR:HB	1.84	0.42
1:A:189:ASN:HB3	1:A:190:PRO:HD3	2.01	0.42
1:C:129:TYR:CE1	1:C:133:GLN:HG3	2.55	0.42
1:A:228:ILE:H	1:A:228:ILE:HG12	1.62	0.42
1:B:27:VAL:CG1	1:B:28:ARG:N	2.82	0.42
1:B:248:ASN:HD22	1:B:251:VAL:HG22	1.85	0.42
1:C:168:TYR:OH	1:D:179:ARG:NH2	2.53	0.42
1:C:170:LYS:HE3	1:D:42:SER:CB	2.50	0.42
1:A:31:TRP:CD1	1:A:38:ALA:HB2	2.55	0.41
1:A:123:LYS:HD3	1:A:123:LYS:HA	1.93	0.41
1:B:3:GLN:O	1:B:4:PHE:HB3	2.20	0.41
1:B:138:PRO:HB2	1:B:169:VAL:HG11	2.02	0.41
1:D:96:GLN:HB3	1:D:97:TRP:CZ3	2.55	0.41
1:D:131:LEU:O	1:D:135:LYS:HE3	2.20	0.41
1:B:248:ASN:HA	1:B:249:PRO:HD3	1.85	0.41
1:C:88:MET:HE2	1:C:88:MET:HA	2.01	0.41
1:C:271:ASP:HB3	1:C:272:LYS:H	1.58	0.41
1:C:189:ASN:O	1:C:193:VAL:HG23	2.19	0.41
1:D:84:ASN:O	1:D:87:ASN:HB2	2.20	0.41
1:A:185:MET:O	1:A:190:PRO:HD3	2.21	0.41
1:B:61:LYS:HD3	1:B:62:LYS:O	2.21	0.41
1:B:171:HIS:C	1:B:173:LYS:H	2.28	0.41
1:B:196:TYR:O	1:B:199:LEU:HB3	2.20	0.41
1:C:255:TYR:CD1	1:C:255:TYR:N	2.88	0.41
1:D:44:ILE:HG22	1:D:45:SER:N	2.36	0.41
1:D:174:LEU:HD12	1:D:174:LEU:HA	1.89	0.41
1:C:169:VAL:HA	1:C:173:LYS:O	2.21	0.41
1:A:162:VAL:HG13	1:A:178:VAL:CG1	2.50	0.41
1:D:131:LEU:O	1:D:135:LYS:HG2	2.21	0.41
1:A:170:LYS:O	1:A:171:HIS:HB2	2.20	0.41
1:A:134:LEU:HD23	1:A:134:LEU:HA	1.78	0.41
1:A:182:SER:HB3	1:A:220:ASP:HB3	2.03	0.41
1:B:16:ILE:O	1:B:16:ILE:HG22	2.20	0.41
1:B:166:GLN:NE2	1:B:168:TYR:CE1	2.89	0.41
1:C:170:LYS:O	1:C:171:HIS:HB2	2.21	0.41
1:C:232:LYS:O	1:C:236:GLU:HG2	2.21	0.41
1:D:168:TYR:HB2	1:D:175:HIS:HB2	2.02	0.41
1:A:73:LEU:HA	1:A:73:LEU:HD12	1.77	0.41
1:B:243:PRO:HG2	1:B:264:LEU:HD11	2.03	0.41
1:C:4:PHE:CE1	1:C:193:VAL:HG21	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:ASN:HD22	1:C:92:HIS:CE1	2.40	0.41
1:C:255:TYR:N	1:C:255:TYR:HD1	2.19	0.41
1:D:46:LYS:HD2	1:D:46:LYS:HA	1.87	0.40
1:D:130:LEU:HD11	1:D:167:TRP:CD1	2.55	0.40
1:B:178:VAL:HB	1:B:216:PHE:HD1	1.87	0.40
1:C:31:TRP:O	1:C:35:GLY:HA2	2.20	0.40
1:C:41:LEU:HD23	1:C:41:LEU:HA	1.97	0.40
1:A:124:VAL:HG22	1:A:125:ASP:N	2.36	0.40
1:C:23:GLU:C	1:C:25:PHE:H	2.29	0.40
1:C:247:ILE:HA	1:C:261:ASP:O	2.20	0.40
1:A:7:GLN:O	1:A:10:SER:HB2	2.21	0.40
1:B:31:TRP:CE3	1:B:278:ALA:HB2	2.56	0.40
1:D:40:THR:HB	1:D:41:LEU:H	1.78	0.40
1:D:206:VAL:HG21	1:D:254:PHE:HB2	2.04	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	276/278 (99%)	241 (87%)	30 (11%)	5 (2%)	6	13
1	B	276/278 (99%)	242 (88%)	27 (10%)	7 (2%)	4	8
1	C	276/278 (99%)	234 (85%)	34 (12%)	8 (3%)	3	6
1	D	276/278 (99%)	239 (87%)	29 (10%)	8 (3%)	3	6
All	All	1104/1112 (99%)	956 (87%)	120 (11%)	28 (2%)	4	8

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	VAL

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Mol	Chain	Res	Type
1	A	37	PRO
1	B	27	VAL
1	B	37	PRO
1	B	271	ASP
1	C	27	VAL
1	C	37	PRO
1	C	271	ASP
1	D	27	VAL
1	D	37	PRO
1	D	271	ASP
1	A	271	ASP
1	D	29	THR
1	C	23	GLU
1	D	28	ARG
1	D	45	SER
1	A	104	ILE
1	D	23	GLU
1	B	24	GLU
1	B	45	SER
1	B	108	TYR
1	C	4	PHE
1	C	26	ASP
1	C	189	ASN
1	D	108	TYR
1	A	137	ASN
1	B	56	PRO
1	C	206	VAL

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	252/253 (100%)	221 (88%)	31 (12%)	4	8
1	B	252/253 (100%)	231 (92%)	21 (8%)	10	19
1	C	252/253 (100%)	236 (94%)	16 (6%)	16	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	252/253 (100%)	229 (91%)	23 (9%)	9	17
All	All	1008/1012 (100%)	917 (91%)	91 (9%)	9	17

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	23	GLU
1	A	29	THR
1	A	33	SER
1	A	36	THR
1	A	39	HIS
1	A	42	SER
1	A	45	SER
1	A	46	LYS
1	A	49	ARG
1	A	58	LEU
1	A	61	LYS
1	A	66	LYS
1	A	73	LEU
1	A	87	ASN
1	A	100	GLU
1	A	119	LEU
1	A	137	ASN
1	A	142	ARG
1	A	159	THR
1	A	173	LYS
1	A	211	LEU
1	A	217	ASN
1	A	218	ILE
1	A	228	ILE
1	A	237	ARG
1	A	251	VAL
1	A	255	TYR
1	A	266	ASN
1	A	272	LYS
1	A	274	LEU
1	B	3	GLN
1	B	23	GLU
1	B	24	GLU
1	B	36	THR
1	B	46	LYS

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Mol	Chain	Res	Type
1	B	49	ARG
1	B	61	LYS
1	B	87	ASN
1	B	142	ARG
1	B	145	THR
1	B	165	THR
1	B	211	LEU
1	B	217	ASN
1	B	218	ILE
1	B	245	LEU
1	B	251	VAL
1	B	255	TYR
1	B	259	ILE
1	B	266	ASN
1	B	271	ASP
1	B	272	LYS
1	C	3	GLN
1	C	5	ASP
1	C	29	THR
1	C	30	LYS
1	C	36	THR
1	C	37	PRO
1	C	39	HIS
1	C	61	LYS
1	C	73	LEU
1	C	95	ASP
1	C	159	THR
1	C	217	ASN
1	C	226	ARG
1	C	259	ILE
1	C	271	ASP
1	C	272	LYS
1	D	3	GLN
1	D	22	ASP
1	D	23	GLU
1	D	29	THR
1	D	30	LYS
1	D	34	ASP
1	D	37	PRO
1	D	45	SER
1	D	52	ASN
1	D	61	LYS

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Mol	Chain	Res	Type
1	D	73	LEU
1	D	139	SER
1	D	142	ARG
1	D	165	THR
1	D	173	LYS
1	D	178	VAL
1	D	211	LEU
1	D	225	THR
1	D	226	ARG
1	D	251	VAL
1	D	259	ILE
1	D	271	ASP
1	D	272	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	77	GLN
1	A	87	ASN
1	A	99	GLN
1	A	120	ASN
1	A	136	ASN
1	A	137	ASN
1	A	195	GLN
1	A	200	GLN
1	A	217	ASN
1	A	266	ASN
1	A	269	HIS
1	B	7	GLN
1	B	9	ASN
1	B	87	ASN
1	B	99	GLN
1	B	106	HIS
1	B	120	ASN
1	B	137	ASN
1	B	166	GLN
1	B	183	ASN
1	B	189	ASN
1	B	195	GLN
1	B	200	GLN
1	B	234	GLN

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Mol	Chain	Res	Type
1	B	248	ASN
1	B	269	HIS
1	C	7	GLN
1	C	9	ASN
1	C	47	GLN
1	C	77	GLN
1	C	87	ASN
1	C	99	GLN
1	C	120	ASN
1	C	137	ASN
1	C	195	GLN
1	C	269	HIS
1	D	9	ASN
1	D	84	ASN
1	D	87	ASN
1	D	96	GLN
1	D	99	GLN
1	D	111	GLN
1	D	120	ASN
1	D	132	HIS
1	D	136	ASN
1	D	137	ASN
1	D	143	HIS
1	D	183	ASN
1	D	192	ASN
1	D	195	GLN
1	D	200	GLN
1	D	222	HIS
1	D	269	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.