



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 06:00 pm BST

PDB ID : 1QDL
Title : THE CRYSTAL STRUCTURE OF ANTHRANILATE SYNTHASE FROM
SULFOLOBUS SOLFATARICUS
Authors : Knoechel, T.; Ivens, A.; Hester, G.; Gonzalez, A.; Bauerle, R.; Wilmanns, M.;
Kirschner, K.; Jansonius, J.N.
Deposited on : 1999-05-20
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

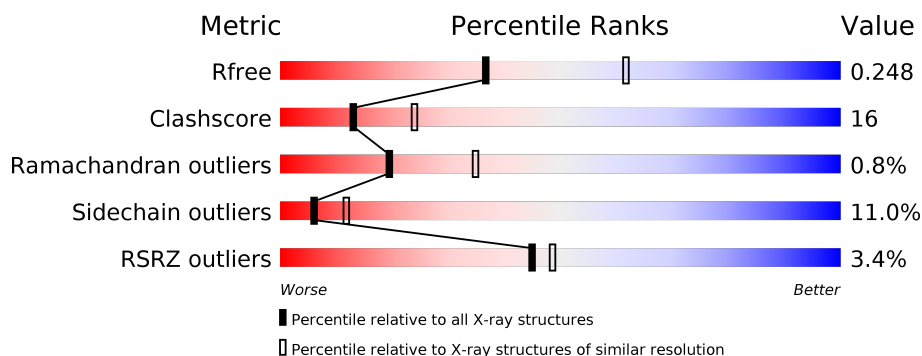
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	422	<div> <div>5%</div> <div> <div></div> <div>57%</div> <div>36%</div> <div>5%</div> </div> </div>
2	B	195	<div> <div>%</div> <div> <div></div> <div>66%</div> <div>32%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5003 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 (ANTHRANILATE SYNTHASE (TRPE-SUBUNIT)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	1
			3324	2150	544	619	11			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	337	ALA	ARG	CONFLICT	UNP Q06128

- Molecule 2 is a protein called PROTEIN (ANTHRANILATE SYNTHASE (TRPG-SUBUNIT)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	195	Total	C	N	O	S	0	0	0
			1549	999	264	283	3			

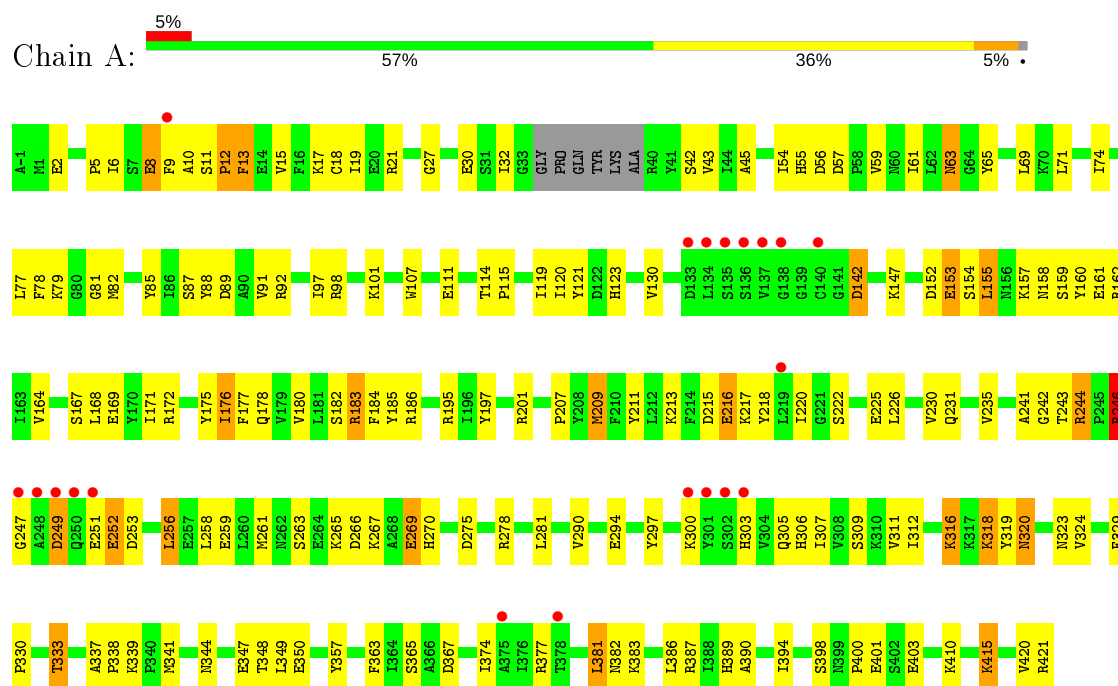
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	90	Total	O	0	0
			90	90		
3	B	40	Total	O	0	0
			40	40		

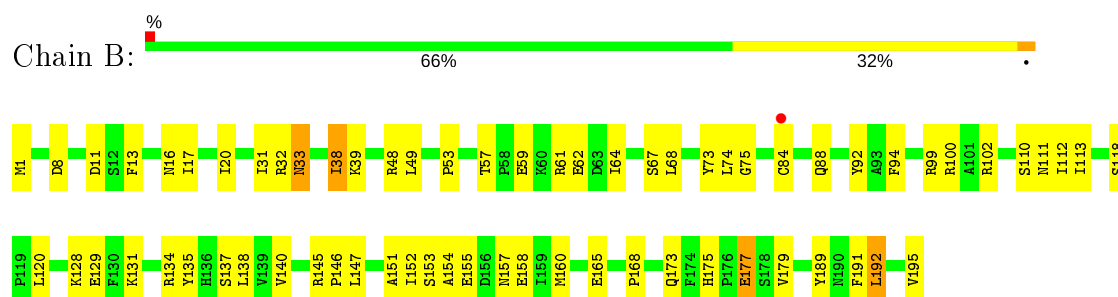
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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 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.

• Molecule 1: PROTEIN (ANTHRANILATE SYNTHASE (TRPE-SUBUNIT))



• Molecule 2: PROTEIN (ANTHRANILATE SYNTHASE (TRPG-SUBUNIT))



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	162.00Å 162.00Å 212.35Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.50 29.65 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.50) 99.8 (29.65-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.46 (at 2.51Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.226 , 0.254 0.216 , 0.248	Depositor DCC
R_{free} test set	2887 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 64.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5003	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.45	0/3404	0.67	0/4605
2	B	0.39	0/1583	0.65	0/2145
All	All	0.43	0/4987	0.67	0/6750

There are no bond length outliers.

There are no bond angle outliers.

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	3324	0	3299	116	0
2	B	1549	0	1574	45	0
3	A	90	0	0	2	0
3	B	40	0	0	1	0
All	All	5003	0	4873	160	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 (160) 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)
2:B:48:ARG:HD2	2:B:192:LEU:HD22	1.52	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:LEU:HD22	1:A:386:LEU:HG	1.55	0.85
1:A:249:ASP:OD2	1:A:251:GLU:HB2	1.78	0.82
1:A:42:SER:HB2	1:A:121:TYR:HB2	1.62	0.80
2:B:146:PRO:O	2:B:165:GLU:HG3	1.84	0.78
1:A:420:VAL:HG12	1:A:421:ARG:HG2	1.64	0.77
2:B:145:ARG:HD2	2:B:165:GLU:OE2	1.89	0.72
2:B:128:LYS:HG3	2:B:129:GLU:HG2	1.70	0.71
1:A:56:ASP:HB2	1:A:61:ILE:HD11	1.72	0.71
1:A:387:ARG:HH11	1:A:389:HIS:HD2	1.39	0.68
1:A:77:LEU:HD11	1:A:231:GLN:HG3	1.75	0.68
1:A:85:TYR:CE2	1:A:111:GLU:HB2	2.29	0.68
1:A:415:LYS:HG2	1:A:420:VAL:HG11	1.77	0.67
1:A:154:SER:O	1:A:155:LEU:HB2	1.96	0.66
1:A:222:SER:O	1:A:377:ARG:HD2	1.96	0.65
1:A:159:SER:O	1:A:162:ARG:HG2	1.98	0.63
1:A:45:ALA:HB1	1:A:115:PRO:HB3	1.82	0.62
1:A:10:ALA:HB3	1:A:15:VAL:HG22	1.82	0.61
1:A:381:LEU:N	1:A:381:LEU:HD23	2.14	0.61
1:A:81:GLY:O	1:A:114:THR:HG23	2.01	0.61
1:A:11:SER:O	1:A:15:VAL:HG23	2.00	0.61
1:A:270:HIS:HE1	1:A:309:SER:OG	1.84	0.60
2:B:75:GLY:O	2:B:168:PRO:HD2	2.02	0.60
1:A:263:SER:O	1:A:267:LYS:HG3	2.01	0.60
1:A:175:TYR:O	1:A:341:MET:HG2	2.01	0.59
2:B:110:SER:HB2	2:B:134:ARG:HD3	1.84	0.59
2:B:38:ILE:HG22	3:B:210:HOH:O	2.02	0.58
1:A:318:LYS:HD2	1:A:319:TYR:CZ	2.39	0.57
1:A:15:VAL:O	1:A:19:ILE:HG12	2.05	0.57
2:B:64:ILE:HB	2:B:67:SER:HB2	1.86	0.57
1:A:10:ALA:HB3	1:A:15:VAL:CG2	2.35	0.56
1:A:101:LYS:HD3	1:A:347:GLU:O	2.05	0.56
2:B:53:PRO:HB3	2:B:84:CYS:HB3	1.87	0.56
1:A:13:PHE:HD1	1:A:201:ARG:HH11	1.54	0.56
1:A:32:ILE:H	1:A:207:PRO:HA	1.71	0.56
1:A:98:ARG:HG3	1:A:344:ASN:OD1	2.07	0.55
2:B:153:SER:HB3	2:B:158:GLU:H	1.71	0.55
1:A:176:ILE:HA	1:A:338:PRO:HD2	1.89	0.55
2:B:146:PRO:HA	2:B:165:GLU:OE2	2.06	0.55
2:B:38:ILE:HD11	2:B:74:LEU:HD13	1.89	0.55
2:B:111:ASN:HD21	2:B:131:LYS:HE3	1.71	0.54
1:A:154:SER:HB2	1:A:184:PHE:CZ	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:LEU:HD12	1:A:259:GLU:OE2	2.08	0.54
1:A:77:LEU:CD1	1:A:231:GLN:HG3	2.38	0.53
1:A:241:ALA:HB1	1:A:269:GLU:HG2	1.90	0.53
1:A:294:GLU:HB3	1:A:297:TYR:CD2	2.43	0.53
1:A:365:SER:HB3	1:A:367:ASP:OD2	2.09	0.53
2:B:74:LEU:CD1	2:B:74:LEU:N	2.72	0.53
1:A:225:GLU:HB3	1:A:374:ILE:HD12	1.90	0.53
2:B:153:SER:O	2:B:157:ASN:HA	2.09	0.53
1:A:249:ASP:N	1:A:249:ASP:OD1	2.43	0.52
1:A:111:GLU:OE2	1:A:217:LYS:HE3	2.09	0.52
1:A:157:LYS:NZ	1:A:401:GLU:HG3	2.25	0.52
1:A:43:VAL:HG13	1:A:120:ILE:CD1	2.40	0.52
2:B:48:ARG:CD	2:B:192:LEU:HD22	2.31	0.52
2:B:59:GLU:HG2	2:B:92:TYR:CE1	2.45	0.52
2:B:64:ILE:CD1	2:B:68:LEU:HD12	2.40	0.52
1:A:218:TYR:HB2	1:A:381:LEU:HG	1.91	0.51
1:A:153:GLU:HB2	1:A:185:TYR:CD2	2.46	0.51
1:A:164:VAL:HG11	1:A:400:PRO:HA	1.93	0.51
2:B:74:LEU:HD12	2:B:74:LEU:N	2.26	0.50
1:A:247:GLY:HA3	1:A:253:ASP:CA	2.41	0.50
1:A:249:ASP:C	1:A:251:GLU:H	2.15	0.50
1:A:275:ASP:OD1	1:A:278:ARG:NH1	2.45	0.50
1:A:249:ASP:C	1:A:251:GLU:N	2.65	0.49
1:A:256:LEU:HB3	1:A:305:GLN:HE21	1.77	0.49
2:B:17:ILE:CG1	2:B:177:GLU:HG2	2.42	0.49
1:A:11:SER:HB2	1:A:12:PRO:HD2	1.94	0.49
1:A:318:LYS:HG2	3:A:432:HOH:O	2.10	0.49
1:A:17:LYS:O	1:A:21:ARG:HG3	2.12	0.49
1:A:119:ILE:HG12	1:A:130:VAL:HG13	1.94	0.49
1:A:230:VAL:HG22	1:A:235:VAL:HG22	1.96	0.48
1:A:316:LYS:HB2	1:A:319:TYR:CD1	2.49	0.48
1:A:180:VAL:HG21	1:A:333:THR:O	2.13	0.48
1:A:243:THR:HG23	1:A:306:HIS:CE1	2.48	0.48
1:A:5:PRO:HG2	1:A:8:GLU:HB2	1.95	0.48
1:A:263:SER:HB3	1:A:266:ASP:HB2	1.95	0.48
1:A:243:THR:O	1:A:244:ARG:HD3	2.14	0.48
1:A:74:ILE:HD11	1:A:79:LYS:HA	1.95	0.48
1:A:89:ASP:O	1:A:92:ARG:HG3	2.14	0.48
1:A:241:ALA:O	1:A:307:ILE:HB	2.14	0.48
1:A:330:PRO:HD2	1:A:339:LYS:HD3	1.96	0.47
1:A:142:ASP:N	1:A:142:ASP:OD1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:TYR:CD1	1:A:220:ILE:HG12	2.50	0.47
1:A:213:LYS:HG3	1:A:218:TYR:CE2	2.50	0.47
1:A:294:GLU:HB3	1:A:297:TYR:CE2	2.50	0.47
2:B:145:ARG:HB3	2:B:165:GLU:OE2	2.14	0.47
1:A:348:THR:HG22	1:A:349:LEU:HD23	1.96	0.47
1:A:88:TYR:HA	1:A:329:PHE:CZ	2.51	0.46
2:B:64:ILE:HD11	2:B:68:LEU:HD12	1.97	0.46
1:A:320:ASN:O	1:A:324:VAL:HG23	2.15	0.46
2:B:53:PRO:HB3	2:B:84:CYS:CB	2.46	0.46
1:A:258:LEU:HA	1:A:261:MET:HB2	1.97	0.46
2:B:134:ARG:HD2	2:B:160:MET:HE2	1.97	0.46
2:B:175:HIS:HB3	2:B:177:GLU:OE2	2.16	0.45
1:A:182:SER:HB2	1:A:390:ALA:O	2.16	0.45
1:A:175:TYR:O	1:A:176:ILE:HG23	2.17	0.45
1:A:337:ALA:HA	1:A:338:PRO:C	2.36	0.45
1:A:357:TYR:CD1	1:A:389:HIS:HB3	2.51	0.45
1:A:101:LYS:HE2	1:A:350:GLU:O	2.17	0.45
1:A:256:LEU:O	1:A:259:GLU:HB2	2.15	0.45
2:B:64:ILE:HG13	2:B:64:ILE:H	1.66	0.45
1:A:247:GLY:HA3	1:A:253:ASP:CB	2.47	0.45
1:A:186:ARG:HG2	1:A:387:ARG:HB2	1.99	0.45
2:B:111:ASN:ND2	2:B:131:LYS:HG3	2.32	0.45
1:A:177:PHE:CZ	2:B:135:TYR:HB3	2.52	0.44
1:A:247:GLY:HA3	1:A:253:ASP:HB2	1.99	0.44
1:A:91:VAL:HG13	1:A:97:ILE:HD13	2.00	0.44
2:B:112:ILE:O	2:B:129:GLU:HA	2.18	0.44
1:A:57:ASP:O	1:A:61:ILE:HG12	2.17	0.44
1:A:13:PHE:CE1	1:A:197:TYR:HD2	2.35	0.44
1:A:168:LEU:O	1:A:172:ARG:HD3	2.18	0.44
1:A:27:GLY:HA2	1:A:211:TYR:O	2.17	0.44
1:A:216:GLU:HG3	1:A:216:GLU:H	1.39	0.43
2:B:191:PHE:HA	2:B:195:VAL:CG2	2.48	0.43
1:A:54:ILE:HG23	1:A:61:ILE:HG13	1.99	0.43
2:B:16:ASN:O	2:B:20:ILE:HG13	2.18	0.43
1:A:160:TYR:CG	1:A:183:ARG:HD2	2.53	0.43
1:A:242:GLY:O	1:A:306:HIS:HA	2.18	0.43
2:B:73:TYR:N	2:B:73:TYR:CD1	2.85	0.43
1:A:121:TYR:HB3	1:A:123:HIS:CE1	2.53	0.43
2:B:94:PHE:HB2	2:B:147:LEU:CD1	2.49	0.43
1:A:30:GLU:HB2	1:A:209:MET:HB2	2.01	0.43
1:A:153:GLU:HB2	1:A:185:TYR:HD2	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:PHE:HD1	1:A:363:PHE:CD1	2.37	0.42
1:A:91:VAL:HG21	1:A:330:PRO:HD3	2.01	0.42
2:B:113:ILE:HD11	2:B:154:ALA:HB2	2.01	0.42
2:B:155:GLU:H	2:B:155:GLU:CD	2.22	0.42
1:A:19:ILE:H	1:A:19:ILE:HG12	1.72	0.42
1:A:246:ARG:HD2	1:A:246:ARG:HA	1.53	0.42
2:B:112:ILE:HG22	2:B:151:ALA:HB1	2.01	0.42
2:B:94:PHE:HB2	2:B:147:LEU:HD11	2.02	0.42
2:B:189:TYR:C	2:B:189:TYR:CD2	2.92	0.42
2:B:33:ASN:HB2	2:B:64:ILE:O	2.20	0.42
1:A:246:ARG:O	1:A:256:LEU:HD23	2.20	0.42
1:A:281:LEU:HD22	1:A:311:VAL:HG12	2.01	0.42
1:A:320:ASN:HB2	1:A:323:ASN:OD1	2.19	0.42
1:A:107:TRP:CZ2	1:A:382:ASN:HB2	2.55	0.42
1:A:59:VAL:O	1:A:63:ASN:HB2	2.20	0.42
1:A:65:TYR:O	1:A:69:LEU:HG	2.20	0.42
2:B:88:GLN:NE2	2:B:173:GLN:HE22	2.17	0.42
1:A:61:ILE:H	1:A:61:ILE:HG12	1.73	0.41
2:B:100:ARG:HG2	2:B:138:LEU:HD23	2.02	0.41
1:A:316:LYS:HB2	1:A:319:TYR:HD1	1.84	0.41
1:A:74:ILE:O	1:A:79:LYS:HD3	2.20	0.41
1:A:157:LYS:O	1:A:161:GLU:HG3	2.21	0.41
1:A:55:HIS:HB2	3:A:434:HOH:O	2.19	0.41
2:B:17:ILE:HG13	2:B:177:GLU:HG2	2.01	0.41
1:A:167:SER:O	1:A:171:ILE:HG13	2.20	0.41
1:A:152:ASP:OD2	1:A:186:ARG:HB2	2.21	0.41
1:A:290:VAL:HA	1:A:312:ILE:O	2.20	0.41
2:B:8:ASP:O	2:B:32:ARG:HA	2.21	0.41
1:A:381:LEU:HD22	1:A:386:LEU:CG	2.39	0.41
1:A:252:GLU:O	1:A:256:LEU:HD22	2.21	0.41
1:A:178:GLN:HA	1:A:394:ILE:O	2.21	0.41
2:B:17:ILE:HG12	2:B:177:GLU:HG2	2.02	0.41
2:B:11:ASP:OD1	2:B:13:PHE:HB2	2.21	0.40
1:A:182:SER:OG	1:A:389:HIS:HE1	2.04	0.40
1:A:43:VAL:HG13	1:A:120:ILE:HD13	2.03	0.40
1:A:381:LEU:N	1:A:381:LEU:CD2	2.84	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	412/422 (98%)	386 (94%)	22 (5%)	4 (1%)	15	28
2	B	193/195 (99%)	186 (96%)	6 (3%)	1 (0%)	29	48
All	All	605/617 (98%)	572 (94%)	28 (5%)	5 (1%)	19	35

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	246	ARG
1	A	155	LEU
2	B	33	ASN
1	A	215	ASP
1	A	12	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	356/360 (99%)	316 (89%)	40 (11%)	6	11
2	B	169/169 (100%)	151 (89%)	18 (11%)	6	13
All	All	525/529 (99%)	467 (89%)	58 (11%)	6	12

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU

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Mol	Chain	Res	Type
1	A	6	ILE
1	A	8	GLU
1	A	9	PHE
1	A	13	PHE
1	A	18	CYS
1	A	63	ASN
1	A	71	LEU
1	A	82	MET
1	A	87	SER
1	A	142	ASP
1	A	147	LYS
1	A	153	GLU
1	A	158	ASN
1	A	169	GLU
1	A	176	ILE
1	A	183	ARG
1	A	195	ARG
1	A	209	MET
1	A	216	GLU
1	A	226	LEU
1	A	244	ARG
1	A	246	ARG
1	A	249	ASP
1	A	252	GLU
1	A	256	LEU
1	A	265	LYS
1	A	269	GLU
1	A	300	LYS
1	A	303	HIS
1	A	316	LYS
1	A	318	LYS
1	A	320	ASN
1	A	333	THR
1	A	381	LEU
1	A	383	LYS
1	A	398	SER
1	A	403	GLU
1	A	410	LYS
1	A	415	LYS
2	B	1	MET
2	B	31	ILE
2	B	38	ILE

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Mol	Chain	Res	Type
2	B	39	LYS
2	B	49	LEU
2	B	57	THR
2	B	61	ARG
2	B	62	GLU
2	B	99	ARG
2	B	102	ARG
2	B	118	SER
2	B	120	LEU
2	B	137	SER
2	B	140	VAL
2	B	152	ILE
2	B	177	GLU
2	B	179	VAL
2	B	192	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	63	ASN
1	A	270	HIS
1	A	306	HIS
1	A	320	ASN
1	A	389	HIS
2	B	88	GLN
2	B	111	ASN

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 carbohydrates 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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	416/422 (98%)	0.02	20 (4%) 30 32	28, 47, 88, 100	0
2	B	195/195 (100%)	-0.20	1 (0%) 91 91	32, 48, 75, 90	0
All	All	611/617 (99%)	-0.05	21 (3%) 45 48	28, 47, 84, 100	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	VAL	5.2
1	A	133	ASP	4.1
1	A	138	GLY	4.1
1	A	135	SER	3.5
1	A	302	SER	3.5
1	A	303	HIS	3.2
1	A	134	LEU	3.2
1	A	248	ALA	3.2
1	A	301	TYR	3.1
1	A	136	SER	3.1
1	A	251	GLU	3.1
1	A	247	GLY	3.0
1	A	375	ALA	2.9
1	A	249	ASP	2.9
1	A	250	GLN	2.7
1	A	140	CYS	2.4
1	A	9	PHE	2.2
2	B	84	CYS	2.2
1	A	219	LEU	2.1
1	A	300	LYS	2.1
1	A	378	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.