



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:31 PM UTC

PDB ID : 1H3F / pdb_00001h3f
Title : Tyrosyl-tRNA synthetase from *Thermus thermophilus* complexed with tyrosinol
Authors : Cusack, S.; Yaremchuk, A.; Kriklivyi, I.; Tukalo, M.
Deposited on : 2002-08-28
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 ⓘ 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

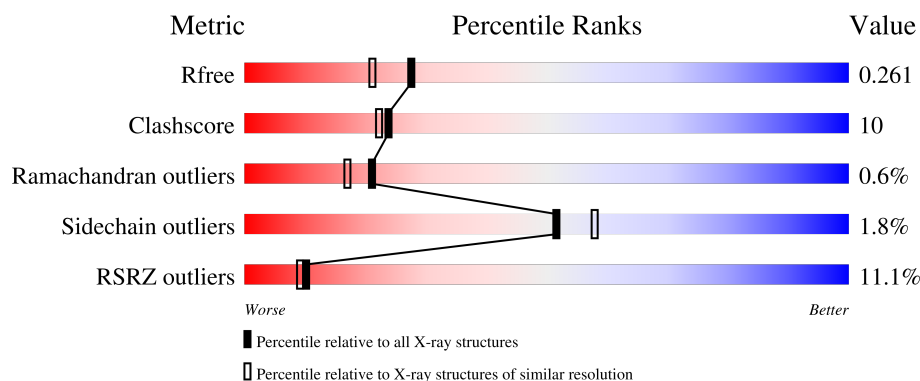
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 $\leq 5\%$. 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	432	8% 75% 17% • 7%
1	B	432	12% 73% 19% • 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	TYE	A	1433	-	-	X	-

2 Entry composition [i](#)

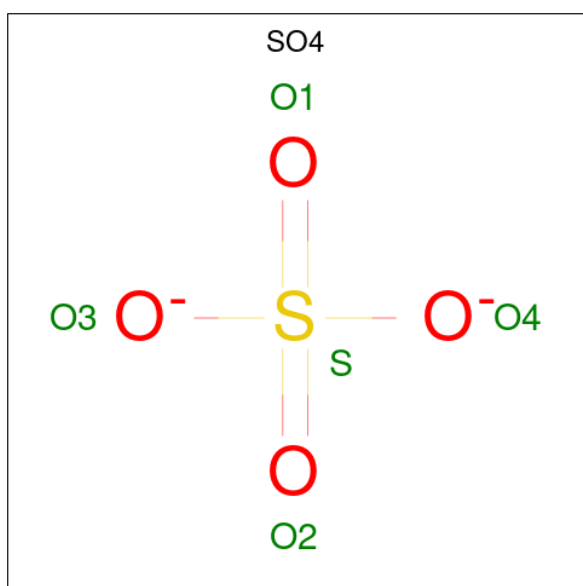
There are 4 unique types of molecules in this entry. The entry contains 6777 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 TYROSYL-TRNA SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3229	2056	574	589	10			
1	B	406	Total	C	N	O	S	0	0	0
			3245	2066	576	592	11			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



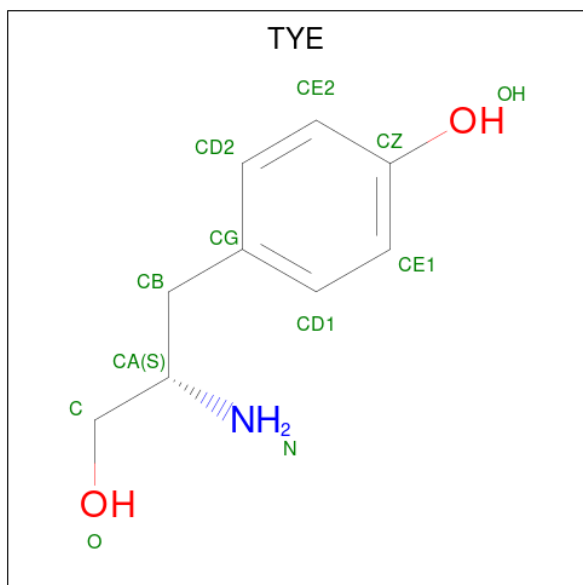
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 4-[(2S)-2-amino-3-hydroxypropyl]phenol (CCD ID: TYE) (formula: $C_9H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			12	9	1	2		
3	B	1	Total	C	N	O	0	0
			12	9	1	2		

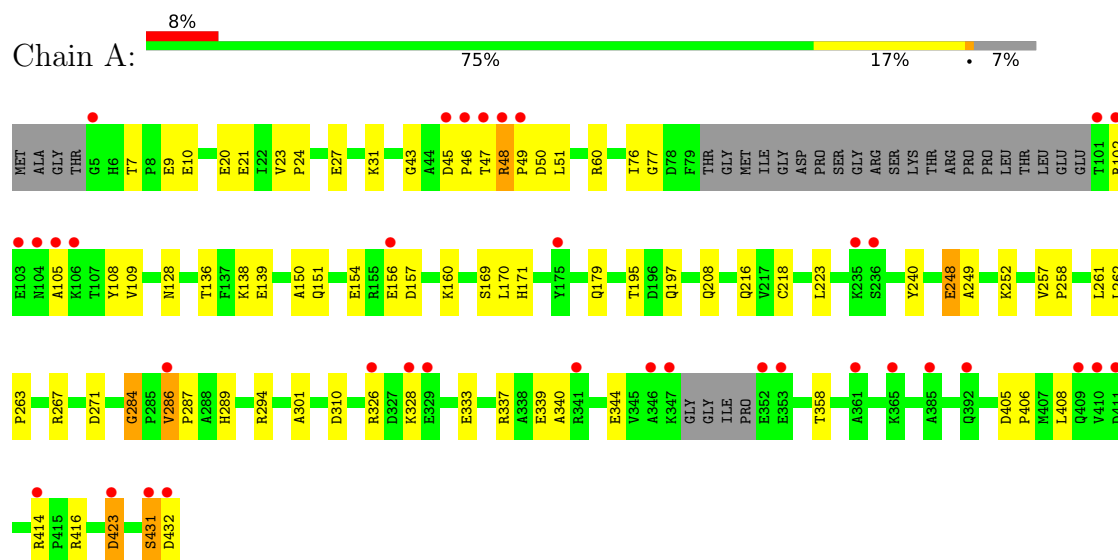
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	146	Total	O	0	0
			146	146		
4	B	108	Total	O	0	0
			108	108		

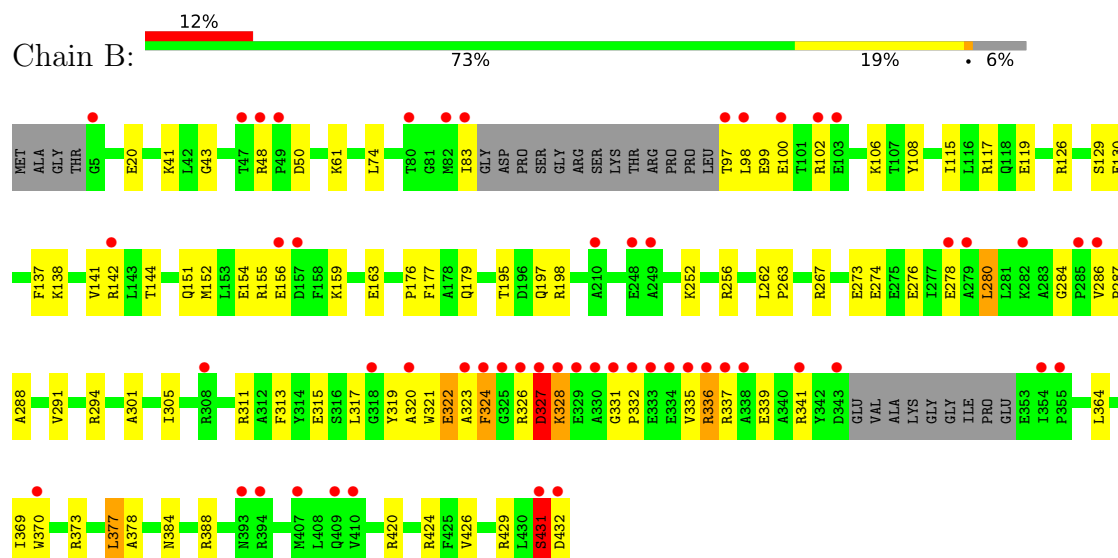
3 Residue-property plots [i](#)

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 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: TYROSYL-TRNA SYNTHETASE



• Molecule 1: TYROSYL-TRNA SYNTHETASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.79Å 111.05Å 141.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.00 25.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.5 (25.00-2.00) 97.5 (25.00-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 1.99Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.234 , 0.261 0.233 , 0.261	Depositor DCC
R_{free} test set	3559 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtrriage
Anisotropy	0.484	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 45.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6777	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.58 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.5682e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

Bond lengths and bond angles in the following residue types are not validated in this section: TYE, SO4

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	0.40	0/3291	0.85	5/4445 (0.1%)
1	B	0.39	0/3307	0.86	9/4467 (0.2%)
All	All	0.40	0/6598	0.86	14/8912 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	328	LYS	N-CA-C	6.68	120.89	112.87
1	A	284	GLY	CA-C-N	6.02	125.23	118.97
1	A	284	GLY	C-N-CA	6.02	125.23	118.97
1	B	117	ARG	N-CA-C	-5.93	102.14	110.50
1	B	322	GLU	N-CA-C	5.91	119.67	112.23
1	A	240	TYR	N-CA-C	5.82	118.33	109.95
1	A	301	ALA	N-CA-C	-5.72	105.80	112.89
1	A	423	ASP	N-CA-C	5.60	119.91	112.30
1	B	61	LYS	N-CA-C	-5.39	105.48	111.36
1	B	48	ARG	CA-C-N	5.19	124.69	119.19
1	B	48	ARG	C-N-CA	5.19	124.69	119.19
1	B	129	SER	N-CA-C	-5.18	106.22	112.54
1	B	41	LYS	N-CA-C	5.08	118.49	109.96
1	B	301	ALA	N-CA-C	-5.02	107.00	113.23

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	3229	0	3295	71	0
1	B	3245	0	3311	64	0
2	A	25	0	0	1	0
3	A	12	0	12	7	0
3	B	12	0	12	3	0
4	A	146	0	0	9	0
4	B	108	0	0	1	0
All	All	6777	0	6630	133	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 10.

All (133) 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:286:VAL:HG23	1:A:287:PRO:HD3	1.18	1.13
1:A:405:ASP:HA	4:A:2134:HOH:O	1.55	1.07
1:A:197:GLN:NE2	3:A:1433:TYE:HNA	1.64	0.94
1:A:358:THR:HG21	1:A:432:ASP:HB2	1.49	0.91
1:A:286:VAL:CG2	1:A:287:PRO:HD3	2.04	0.86
1:A:208:GLN:NE2	1:A:216:GLN:H	1.79	0.80
1:B:294:ARG:NH2	1:B:328:LYS:HA	1.97	0.79
1:A:333:GLU:O	1:A:337:ARG:HG3	1.84	0.78
1:A:197:GLN:HE22	3:A:1433:TYE:HNA	0.83	0.77
1:B:326:ARG:C	1:B:328:LYS:H	1.93	0.76
1:B:323:ALA:HB3	1:B:327:ASP:OD2	1.85	0.76
1:B:364:LEU:HD13	1:B:369:ILE:HB	1.69	0.74
1:A:406:PRO:HD2	4:A:2134:HOH:O	1.86	0.73
1:A:431:SER:HA	4:A:2142:HOH:O	1.87	0.72
1:A:208:GLN:HE22	1:A:216:GLN:H	1.38	0.72
1:B:267:ARG:HD3	4:B:2084:HOH:O	1.92	0.69
1:B:431:SER:O	1:B:432:ASP:HB2	1.93	0.68
1:B:431:SER:O	1:B:432:ASP:CB	2.43	0.66
1:A:340:ALA:O	1:A:344:GLU:HG3	1.96	0.65
1:B:273:GLU:HB2	1:B:276:GLU:HG3	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:THR:C	1:A:48:ARG:HD2	2.23	0.64
1:B:332:PRO:HG2	1:B:335:VAL:HB	1.78	0.64
1:B:20:GLU:HG2	1:B:195:THR:CG2	2.27	0.64
1:B:320:ALA:O	1:B:327:ASP:OD2	2.14	0.64
1:B:280:LEU:HD22	1:B:288:ALA:HA	1.80	0.63
1:B:328:LYS:HG3	1:B:339:GLU:OE1	1.98	0.63
1:B:20:GLU:HG2	1:B:195:THR:HG22	1.81	0.63
1:A:20:GLU:HG2	1:A:195:THR:HG22	1.80	0.62
1:B:370:TRP:HZ2	1:B:388:ARG:HD3	1.63	0.62
1:B:332:PRO:O	1:B:336:ARG:HB3	2.00	0.62
1:A:24:PRO:HG2	1:A:27:GLU:HB2	1.80	0.61
1:B:98:LEU:HD23	1:B:98:LEU:O	2.00	0.61
1:B:280:LEU:HD21	1:B:291:VAL:HG21	1.81	0.61
1:B:323:ALA:HB3	1:B:327:ASP:CG	2.24	0.61
3:A:1433:TYE:H	3:A:1433:TYE:HD1	1.81	0.61
1:A:136:THR:OG1	1:A:139:GLU:HG3	2.01	0.61
1:B:159:LYS:O	1:B:163:GLU:HG3	2.01	0.60
1:B:317:LEU:O	1:B:331:GLY:HA2	2.00	0.60
1:A:286:VAL:HG23	1:A:287:PRO:CD	2.12	0.60
1:A:150:ALA:O	1:A:154:GLU:HG2	2.02	0.60
3:A:1433:TYE:H	3:A:1433:TYE:CD1	2.32	0.59
1:A:160:LYS:HA	1:A:160:LYS:HE2	1.85	0.58
1:B:155:ARG:NH1	1:B:176:PRO:HG3	2.19	0.57
1:A:294:ARG:HH21	1:A:328:LYS:HG2	1.70	0.57
1:A:20:GLU:HG2	1:A:195:THR:CG2	2.35	0.57
1:B:43:GLY:HA3	3:B:1433:TYE:CE1	2.35	0.56
1:B:326:ARG:O	1:B:328:LYS:N	2.31	0.56
1:B:197:GLN:OE1	3:B:1433:TYE:N	2.38	0.56
1:B:286:VAL:HB	1:B:287:PRO:HD3	1.89	0.55
1:A:51:LEU:HD21	1:A:108:TYR:CE2	2.42	0.55
1:B:262:LEU:HB2	1:B:263:PRO:HD3	1.89	0.54
1:A:294:ARG:NH2	1:A:328:LYS:HG2	2.23	0.54
1:B:370:TRP:CZ2	1:B:388:ARG:HD3	2.43	0.54
1:A:77:GLY:HA2	1:A:128:ASN:ND2	2.24	0.53
1:B:274:GLU:O	1:B:278:GLU:HG3	2.08	0.53
1:B:252:LYS:HE3	1:B:256:ARG:NH2	2.23	0.53
1:A:7:THR:OG1	1:A:10:GLU:HG3	2.10	0.52
1:A:257:VAL:O	1:A:289:HIS:HE1	1.91	0.52
1:A:263:PRO:O	1:A:267:ARG:HG3	2.09	0.52
1:A:169:SER:OG	1:A:171:HIS:HD2	1.93	0.52
1:B:326:ARG:C	1:B:328:LYS:N	2.62	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:LEU:HD21	1:B:108:TYR:HE1	1.75	0.52
1:A:258:PRO:HD2	1:A:261:LEU:HD12	1.92	0.52
1:B:138:LYS:HE2	1:B:142:ARG:NH2	2.25	0.52
1:A:179:GLN:OE1	3:A:1433:TYE:N	2.43	0.52
1:A:208:GLN:HE22	1:A:216:GLN:CB	2.23	0.51
1:A:151:GLN:O	1:A:154:GLU:HB2	2.12	0.50
1:A:138:LYS:HG3	1:B:83:ILE:HD13	1.92	0.49
1:B:74:LEU:HD21	1:B:108:TYR:CE1	2.47	0.49
1:A:208:GLN:NE2	1:A:216:GLN:HG2	2.28	0.49
1:A:43:GLY:HA3	3:A:1433:TYE:CE2	2.43	0.49
1:A:23:VAL:HG23	1:A:218:CYS:HB2	1.95	0.49
1:A:60:ARG:HD3	4:A:2099:HOH:O	2.13	0.49
1:A:262:LEU:HB2	1:A:263:PRO:HD3	1.94	0.49
1:A:48:ARG:HD2	1:A:48:ARG:N	2.28	0.48
1:B:378:ALA:HB1	1:B:426:VAL:HG11	1.95	0.48
1:A:326:ARG:HD3	4:A:2115:HOH:O	2.12	0.48
1:B:294:ARG:NH2	1:B:328:LYS:CA	2.74	0.48
1:A:24:PRO:HG2	1:A:27:GLU:CB	2.42	0.48
1:A:102:ARG:HD3	4:A:2039:HOH:O	2.13	0.48
1:B:284:GLY:C	1:B:287:PRO:HD2	2.39	0.48
1:A:170:LEU:HD23	1:B:177:PHE:CE2	2.49	0.47
1:A:208:GLN:HE22	1:A:216:GLN:N	2.08	0.47
1:B:294:ARG:HD3	1:B:319:TYR:CZ	2.50	0.47
1:A:60:ARG:NH2	1:A:271:ASP:OD2	2.47	0.47
1:A:138:LYS:HG3	1:B:83:ILE:CD1	2.45	0.47
1:A:45:ASP:C	1:A:45:ASP:OD2	2.57	0.47
1:B:198:ARG:HH11	1:B:198:ARG:HG3	1.80	0.47
1:A:208:GLN:HE22	1:A:216:GLN:HB3	1.79	0.47
1:B:322:GLU:O	1:B:324:PHE:HD1	1.98	0.47
1:A:60:ARG:HH22	1:A:271:ASP:CG	2.23	0.46
1:B:373:ARG:NE	1:B:377:LEU:HD23	2.30	0.46
1:B:97:THR:HG22	1:B:98:LEU:N	2.29	0.46
1:A:46:PRO:HG3	1:A:76:ILE:CG2	2.45	0.46
1:A:284:GLY:C	1:A:287:PRO:HD2	2.41	0.45
1:B:373:ARG:HA	1:B:384:ASN:OD1	2.16	0.45
1:A:105:ALA:O	1:A:109:VAL:HG23	2.16	0.45
1:B:126:ARG:HH21	1:B:130:GLU:CD	2.25	0.45
1:B:429:ARG:HH11	1:B:429:ARG:HG3	1.82	0.45
1:A:31:LYS:HE3	4:A:2076:HOH:O	2.16	0.44
1:A:328:LYS:HG2	1:A:339:GLU:OE2	2.17	0.44
1:B:337:ARG:O	1:B:341:ARG:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:ARG:HH11	1:A:48:ARG:HG2	1.83	0.44
1:B:328:LYS:O	1:B:328:LYS:HG2	2.17	0.44
1:B:179:GLN:OE1	3:B:1433:TYE:N	2.48	0.44
1:A:170:LEU:HB3	1:B:144:THR:O	2.18	0.44
1:A:156:GLU:HG3	1:A:157:ASP:H	1.81	0.44
1:A:195:THR:OG1	1:A:223:LEU:CD2	2.66	0.44
1:B:311:ARG:O	1:B:315:GLU:HG3	2.18	0.44
1:B:274:GLU:HG3	1:B:278:GLU:OE2	2.18	0.43
1:A:252:LYS:HD2	4:A:2094:HOH:O	2.17	0.43
1:B:276:GLU:HB3	1:B:321:TRP:CZ2	2.54	0.43
1:B:420:ARG:O	1:B:424:ARG:HB2	2.18	0.43
1:A:248:GLU:HG3	1:A:249:ALA:N	2.33	0.43
1:B:364:LEU:HD22	1:B:364:LEU:N	2.34	0.43
1:A:49:PRO:HG2	1:A:50:ASP:H	1.84	0.43
1:B:100:GLU:H	1:B:100:GLU:HG2	1.60	0.43
1:A:310:ASP:HB2	4:A:2100:HOH:O	2.20	0.42
1:B:102:ARG:O	1:B:106:LYS:HG3	2.20	0.42
1:A:197:GLN:NE2	3:A:1433:TYE:N	2.45	0.42
1:B:137:PHE:O	1:B:141:VAL:HG23	2.20	0.41
1:A:284:GLY:O	1:A:287:PRO:HD2	2.20	0.41
1:B:313:PHE:O	1:B:317:LEU:HG	2.20	0.41
1:A:21:GLU:HG2	1:A:23:VAL:HG13	2.02	0.41
1:A:9:GLU:H	1:A:9:GLU:CD	2.29	0.41
1:A:48:ARG:HG2	2:A:502:SO4:O4	2.20	0.41
1:A:77:GLY:HA2	1:A:128:ASN:HD21	1.84	0.41
1:A:23:VAL:HA	1:A:24:PRO:HA	1.77	0.41
1:B:152:MET:C	1:B:154:GLU:H	2.29	0.41
1:B:115:ILE:HG23	1:B:305:ILE:HD11	2.03	0.40
1:A:405:ASP:HB3	1:A:408:LEU:HB2	2.03	0.40
1:A:414:ARG:CZ	1:A:416:ARG:HH12	2.34	0.40
1:B:151:GLN:O	1:B:154:GLU:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	397/432 (92%)	388 (98%)	8 (2%)	1 (0%)	36	35
1	B	400/432 (93%)	386 (96%)	10 (2%)	4 (1%)	12	8
All	All	797/864 (92%)	774 (97%)	18 (2%)	5 (1%)	21	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	431	SER
1	B	327	ASP
1	B	156	GLU
1	B	431	SER
1	B	324	PHE

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	341/363 (94%)	337 (99%)	4 (1%)	63	70
1	B	343/363 (94%)	335 (98%)	8 (2%)	44	49
All	All	684/726 (94%)	672 (98%)	12 (2%)	51	58

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	ARG
1	A	248	GLU
1	A	286	VAL
1	A	423	ASP
1	B	50	ASP
1	B	99	GLU
1	B	119	GLU
1	B	280	LEU

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Mol	Chain	Res	Type
1	B	327	ASP
1	B	336	ARG
1	B	377	LEU
1	B	431	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	HIS
1	A	171	HIS
1	A	197	GLN
1	A	200	ASN
1	A	208	GLN
1	A	289	HIS
1	A	392	GLN
1	A	409	GLN
1	B	64	GLN
1	B	392	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	503	-	4,4,4	0.35	0	6,6,6	0.10	0
2	SO4	A	501	-	4,4,4	0.37	0	6,6,6	0.11	0
3	TYE	B	1433	-	12,12,12	1.26	1 (8%)	13,15,15	0.69	0
2	SO4	A	504	-	4,4,4	0.39	0	6,6,6	0.10	0
3	TYE	A	1433	-	12,12,12	1.24	1 (8%)	13,15,15	0.55	0
2	SO4	A	502	-	4,4,4	0.36	0	6,6,6	0.09	0
2	SO4	A	505	-	4,4,4	0.38	0	6,6,6	0.06	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TYE	B	1433	-	-	2/6/6/6	0/1/1/1
3	TYE	A	1433	-	-	0/6/6/6	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1433	TYE	O-C	-3.86	1.26	1.42
3	A	1433	TYE	O-C	-3.85	1.26	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1433	TYE	O-C-CA-CB
3	B	1433	TYE	O-C-CA-N

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1433	TYE	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1433	TYE	7	0
2	A	502	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	403/432 (93%)	0.36	36 (8%) 15 14	14, 28, 49, 66	0
1	B	406/432 (93%)	0.64	54 (13%) 7 6	15, 31, 59, 81	0
All	All	809/864 (93%)	0.50	90 (11%) 10 9	14, 30, 55, 81	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49	PRO	6.6
1	B	328	LYS	5.6
1	B	327	ASP	5.2
1	B	329	GLU	5.2
1	B	330	ALA	5.1
1	B	324	PHE	4.9
1	B	432	ASP	4.8
1	B	370	TRP	4.6
1	A	347	LYS	4.3
1	A	48	ARG	4.3
1	B	83	ILE	4.1
1	A	46	PRO	4.1
1	B	331	GLY	4.0
1	B	326	ARG	3.9
1	A	235	LYS	3.9
1	B	98	LEU	3.6
1	B	103	GLU	3.6
1	B	318	GLY	3.6
1	B	48	ARG	3.6
1	B	335	VAL	3.5
1	A	352	GLU	3.5
1	B	337	ARG	3.5
1	A	105	ALA	3.4
1	A	101	THR	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	97	THR	3.3
1	A	47	THR	3.3
1	B	5	GLY	3.3
1	A	45	ASP	3.3
1	B	407	MET	3.2
1	B	49	PRO	3.2
1	A	392	GLN	3.2
1	A	346	ALA	3.1
1	B	410	VAL	3.1
1	A	104	ASN	3.1
1	B	278	GLU	3.1
1	A	103	GLU	3.0
1	B	325	GLY	3.0
1	B	341	ARG	2.9
1	B	286	VAL	2.9
1	B	336	ARG	2.9
1	A	365	LYS	2.9
1	A	5	GLY	2.9
1	B	156	GLU	2.9
1	B	343	ASP	2.8
1	B	354	ILE	2.8
1	A	432	ASP	2.7
1	A	102	ARG	2.7
1	A	156	GLU	2.7
1	A	329	GLU	2.7
1	A	431	SER	2.7
1	A	328	LYS	2.7
1	B	409	GLN	2.7
1	A	353	GLU	2.7
1	A	286	VAL	2.7
1	A	361	ALA	2.6
1	B	338	ALA	2.6
1	B	332	PRO	2.6
1	A	411	ASP	2.5
1	B	47	THR	2.5
1	B	248	GLU	2.5
1	A	106	LYS	2.5
1	B	334	GLU	2.4
1	B	80	THR	2.4
1	B	323	ALA	2.4
1	B	282	LYS	2.4
1	B	393	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	341	ARG	2.4
1	A	175	TYR	2.4
1	B	333	GLU	2.4
1	A	423	ASP	2.4
1	B	157	ASP	2.4
1	B	431	SER	2.4
1	B	210	ALA	2.3
1	A	414	ARG	2.3
1	B	142	ARG	2.3
1	B	100	GLU	2.3
1	B	249	ALA	2.3
1	A	409	GLN	2.2
1	A	326	ARG	2.1
1	B	394	ARG	2.1
1	B	308	ARG	2.1
1	A	410	VAL	2.1
1	A	385	ALA	2.1
1	B	285	PRO	2.1
1	B	320	ALA	2.1
1	B	355	PRO	2.0
1	A	236	SER	2.0
1	B	82	MET	2.0
1	B	279	ALA	2.0
1	B	102	ARG	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	TYE	A	1433	12/12	0.67	0.23	35,53,64,66	0
3	TYE	B	1433	12/12	0.71	0.21	37,50,61,64	0
2	SO4	A	501	5/5	0.86	0.10	84,84,85,85	0
2	SO4	A	505	5/5	0.89	0.10	64,65,67,67	0
2	SO4	A	504	5/5	0.90	0.10	49,50,52,53	0
2	SO4	A	502	5/5	0.94	0.10	57,57,61,61	0
2	SO4	A	503	5/5	0.95	0.08	45,51,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.