



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:26 PM UTC

PDB ID : 1TET / pdb_00001tet
Title : CRYSTAL STRUCTURE OF AN ANTICHOLOERA TOXIN PEPTIDE COM-
PLEX AT 2.3 ANGSTROMS
Authors : Shoham, M.
Deposited on : 1993-06-21
Resolution : 2.30 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)	:	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

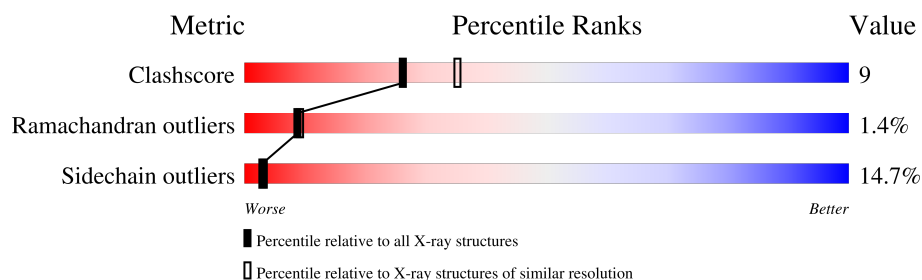
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

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	L	216	 63% 27% 8% •
2	H	210	 58% 30% 10% •
3	P	15	 33% 47% 20%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3541 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 IGG1 TE33 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1681	1059	278	338	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	24	LYS	ARG	conflict	PIR PC4203
L	31A	SER	THR	conflict	PIR PC4203
L	31B	SER	ASN	conflict	PIR PC4203
L	33	PHE	LEU	conflict	PIR PC4203
L	94	ILE	VAL	conflict	PIR PC4203
L	96	PHE	ARG	conflict	PIR PC4203
L	100	SER	GLY	conflict	PIR PC4203
L	188	TRP	ARG	conflict	PIR PC4203

- Molecule 2 is a protein called IGG1 TE33 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	210	Total	C	N	O	S	0	0	1
			1604	1026	262	309	7			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	13	THR	LYS	conflict	GB 2072131
H	26	GLY	ASP	conflict	GB 2072131
H	28	THR	SER	conflict	GB 2072131
H	30	THR	MET	conflict	GB 2072131
H	32	TYR	SER	conflict	GB 2072131
H	35	SER	GLN	conflict	GB 2072131
H	38	LYS	GLN	conflict	GB 2072131
H	40	THR	MET	conflict	GB 2072131

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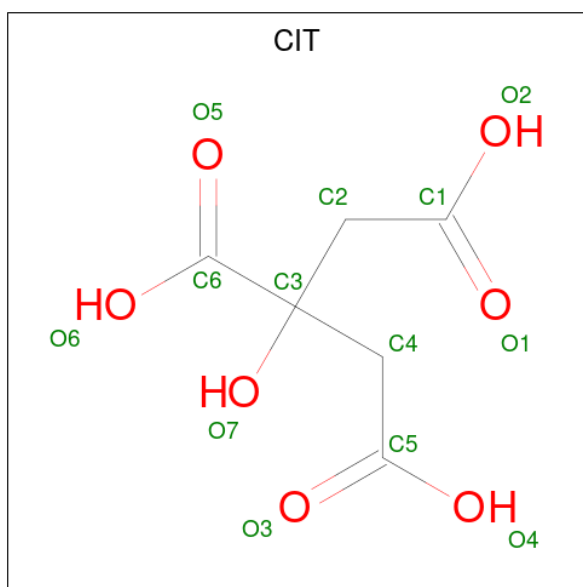
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Chain	Residue	Modelled	Actual	Comment	Reference
H	45	PHE	LEU	conflict	GB 2072131
H	48	MET	ILE	conflict	GB 2072131
H	51	ILE	LEU	conflict	GB 2072131
H	53	TYR	GLN	conflict	GB 2072131
H	58	THR	GLU	conflict	GB 2072131
H	61	ASP	GLU	conflict	GB 2072131
H	76	SER	THR	conflict	GB 2072131
H	?	-	THR	deletion	GB 2072131
H	?	-	TRP	deletion	GB 2072131
H	?	-	GLY	deletion	GB 2072131
H	94	ARG	GLY	conflict	GB 2072131
H	95	ARG	ASN	conflict	GB 2072131
H	100A	TRP	ALA	conflict	GB 2072131
H	100C	PHE	-	insertion	GB 2072131
H	101	ASP	-	insertion	GB 2072131
H	102	VAL	-	insertion	GB 2072131
H	105	THR	GLN	conflict	GB 2072131
H	109	VAL	LEU	conflict	GB 2072131
H	?	-	SER	deletion	GB 2072131
H	?	-	ALA	deletion	GB 2072131
H	?	-	ALA	deletion	GB 2072131
H	?	-	GLN	deletion	GB 2072131
H	?	-	THR	deletion	GB 2072131
H	?	-	ASN	deletion	GB 2072131
H	187	PRO	THR	conflict	GB 2072131
H	188	ARG	TRP	conflict	GB 2072131

- Molecule 3 is a protein called CHOLERA TOXIN PEPTIDE 3 (CTP3).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	12	Total	C	N	O	0	0	0
			92	56	18	18			

- Molecule 4 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	C	O	0	0
			13	6	7		

- Molecule 5 is water.

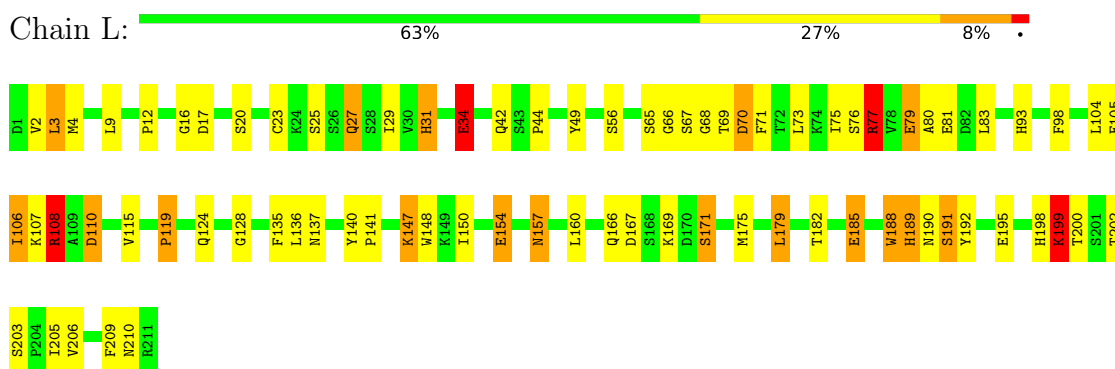
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	66	Total	O	0	0
			66	66		
5	H	78	Total	O	0	0
			78	78		
5	P	7	Total	O	0	0
			7	7		

3 Residue-property plots

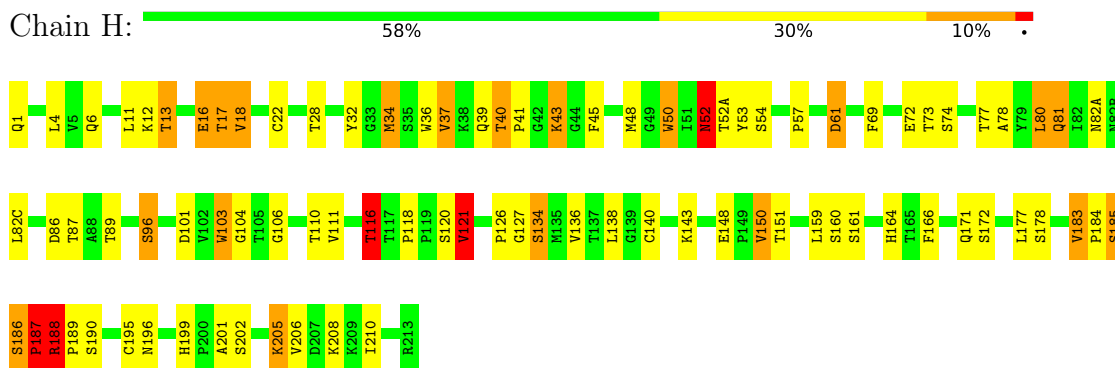
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: IGG1 TE33 FAB (LIGHT CHAIN)



• Molecule 2: IGG1 TE33 FAB (HEAVY CHAIN)



• Molecule 3: CHOLERA TOXIN PEPTIDE 3 (CTP3)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	104.15Å 110.61Å 40.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (5.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.148 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3541	wwPDB-VP
Average B, all atoms (Å ²)	13.0	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: CIT

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	L	1.07	6/1723 (0.3%)	1.89	47/2337 (2.0%)
2	H	1.13	6/1651 (0.4%)	2.00	56/2258 (2.5%)
3	P	1.09	1/93 (1.1%)	2.69	6/123 (4.9%)
All	All	1.10	13/3467 (0.4%)	1.97	109/4718 (2.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	H	0	1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	189	HIS	CD2-NE2	-7.24	1.29	1.37
1	L	198	HIS	CD2-NE2	-6.96	1.30	1.37
1	L	31	HIS	CD2-NE2	-6.88	1.30	1.37
2	H	116	THR	CA-CB	6.75	1.64	1.53
2	H	126	PRO	C-N	-6.69	1.24	1.33
3	P	8	HIS	CD2-NE2	-6.53	1.30	1.37
2	H	199	HIS	CD2-NE2	-6.41	1.30	1.37
2	H	164	HIS	CD2-NE2	-6.35	1.30	1.37
2	H	37	VAL	CA-CB	6.22	1.63	1.54
1	L	31	HIS	CG-ND1	-6.17	1.31	1.38
1	L	93	HIS	CD2-NE2	-5.42	1.31	1.37
1	L	189	HIS	CG-ND1	-5.34	1.32	1.38
2	H	121	VAL	CA-CB	5.07	1.60	1.54

All (109) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	6	SER	N-CA-C	13.27	131.47	114.75
1	L	188	TRP	CG-CD2-CE3	12.18	146.08	133.90
2	H	187	PRO	N-CA-C	11.09	129.32	113.47
2	H	183	VAL	N-CA-CB	-11.00	102.32	111.67
1	L	188	TRP	CB-CG-CD1	-9.79	112.21	126.90
1	L	137	ASN	OD1-CG-ND2	-9.68	112.92	122.60
1	L	42	GLN	OE1-CD-NE2	-9.56	113.04	122.60
2	H	196	ASN	OD1-CG-ND2	-9.09	113.51	122.60
2	H	166	PHE	CA-CB-CG	-8.40	105.40	113.80
3	P	7	GLN	CG-CD-NE2	8.07	128.50	116.40
3	P	8	HIS	CA-CB-CG	-8.07	105.73	113.80
1	L	157	ASN	OD1-CG-ND2	-7.90	114.70	122.60
1	L	29	ILE	CB-CG1-CD1	-7.67	97.70	113.80
2	H	186	SER	CA-C-N	7.40	126.67	118.97
2	H	186	SER	C-N-CA	7.40	126.67	118.97
2	H	52	ASN	OD1-CG-ND2	-7.36	115.25	122.60
1	L	2	VAL	N-CA-C	-7.33	97.18	107.80
1	L	175	MET	CG-SD-CE	-7.11	85.25	100.90
2	H	34	MET	CG-SD-CE	-7.09	85.30	100.90
2	H	196	ASN	CB-CG-ND2	7.07	127.01	116.40
2	H	86	ASP	CA-CB-CG	6.99	119.59	112.60
2	H	164	HIS	CB-CG-CD2	-6.96	122.15	131.20
1	L	110	ASP	CA-CB-CG	6.93	119.53	112.60
2	H	164	HIS	CA-CB-CG	6.76	120.56	113.80
1	L	34	GLU	N-CA-C	6.75	119.70	109.23
3	P	7	GLN	OE1-CD-NE2	-6.73	115.87	122.60
2	H	81	GLN	CA-CB-CG	-6.61	100.87	114.10
1	L	27	GLN	OE1-CD-NE2	-6.56	116.04	122.60
1	L	76	SER	CA-C-O	-6.48	114.02	121.47
1	L	182	THR	N-CA-C	6.45	122.27	113.37
1	L	171	SER	N-CA-C	6.45	120.41	111.90
3	P	8	HIS	CB-CG-CD2	-6.40	122.88	131.20
2	H	187	PRO	O-C-N	-6.23	115.13	122.17
1	L	42	GLN	CG-CD-NE2	6.17	125.65	116.40
1	L	195	GLU	N-CA-C	6.17	119.18	108.76
1	L	93	HIS	CB-CG-CD2	-6.15	123.20	131.20
1	L	188	TRP	CB-CG-CD2	6.04	135.26	126.80
2	H	36	TRP	CG-CD2-CE3	6.01	139.91	133.90
2	H	11	LEU	N-CA-C	-6.00	99.13	108.90
2	H	40	THR	CA-C-N	5.99	127.33	119.84
2	H	40	THR	C-N-CA	5.99	127.33	119.84
2	H	120	SER	CA-C-N	-5.92	115.85	123.19
2	H	120	SER	C-N-CA	-5.92	115.85	123.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	128	GLY	O-C-N	-5.91	116.51	122.54
2	H	111	VAL	O-C-N	-5.89	116.50	123.03
1	L	188	TRP	CE2-CD2-CE3	-5.87	112.93	118.80
2	H	18	VAL	CB-CA-C	-5.84	101.35	110.69
2	H	148	GLU	CA-C-O	-5.83	115.36	120.60
1	L	157	ASN	CB-CG-ND2	5.79	125.09	116.40
2	H	151	THR	CA-CB-OG1	-5.74	100.99	109.60
2	H	120	SER	N-CA-C	-5.73	99.84	109.07
1	L	160	LEU	CA-C-O	5.73	126.31	120.24
2	H	101	ASP	CA-CB-CG	5.73	118.33	112.60
2	H	77	THR	CA-CB-OG1	-5.72	101.02	109.60
2	H	45	PHE	O-C-N	-5.69	116.73	123.22
1	L	198	HIS	CB-CG-CD2	-5.68	123.81	131.20
2	H	150	VAL	N-CA-C	-5.68	100.05	108.45
1	L	98	PHE	CA-CB-CG	5.67	119.47	113.80
1	L	70	ASP	CA-CB-CG	-5.67	106.93	112.60
2	H	45	PHE	N-CA-C	5.67	118.87	109.46
2	H	69	PHE	N-CA-C	-5.64	100.51	109.59
1	L	210	ASN	CA-CB-CG	5.60	118.20	112.60
2	H	202	SER	N-CA-C	-5.58	105.69	112.88
1	L	202	THR	N-CA-C	-5.54	106.66	113.41
3	P	9	ILE	N-CA-CB	-5.51	102.29	112.43
2	H	161	SER	N-CA-C	5.49	117.82	110.35
1	L	9	LEU	N-CA-CB	-5.43	102.18	110.33
2	H	13	THR	N-CA-CB	-5.43	103.46	110.03
1	L	195	GLU	CA-CB-CG	5.41	124.93	114.10
2	H	96	SER	O-C-N	-5.40	115.41	122.59
1	L	136	LEU	N-CA-C	-5.40	96.20	107.37
1	L	191	SER	N-CA-C	5.38	117.67	108.90
1	L	137	ASN	CB-CG-ND2	5.38	124.47	116.40
1	L	106	ILE	N-CA-C	5.38	115.84	107.99
2	H	183	VAL	CA-C-N	5.37	125.63	119.83
2	H	183	VAL	C-N-CA	5.37	125.63	119.83
1	L	93	HIS	CB-CG-ND1	5.37	130.75	122.70
1	L	108	ARG	NE-CZ-NH2	-5.36	114.37	119.20
2	H	188	ARG	CB-CG-CD	5.36	123.64	111.30
1	L	148	TRP	CG-CD2-CE3	5.36	139.26	133.90
2	H	206	VAL	O-C-N	-5.36	117.68	123.14
1	L	79	GLU	CB-CG-CD	5.35	121.69	112.60
2	H	195	CYS	N-CA-CB	-5.34	101.34	110.16
2	H	39	GLN	OE1-CD-NE2	-5.31	117.29	122.60
2	H	28	THR	CA-CB-OG1	-5.30	101.64	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	57	PRO	CA-C-O	-5.29	115.52	122.19
2	H	61	ASP	CA-CB-CG	5.24	117.84	112.60
1	L	3	LEU	N-CA-C	-5.24	101.16	109.96
2	H	199	HIS	CA-C-N	5.23	124.85	119.05
2	H	199	HIS	C-N-CA	5.23	124.85	119.05
2	H	45	PHE	CA-CB-CG	5.23	119.03	113.80
1	L	124	GLN	OE1-CD-NE2	-5.21	117.39	122.60
1	L	166	GLN	OE1-CD-NE2	-5.20	117.40	122.60
1	L	188	TRP	CE2-CD2-CG	-5.19	100.98	107.20
1	L	188	TRP	CA-CB-CG	5.14	123.37	113.60
2	H	164	HIS	CB-CG-ND1	5.13	130.39	122.70
2	H	183	VAL	CB-CA-C	5.07	118.16	110.97
2	H	43	LYS	N-CA-C	5.06	121.58	110.80
2	H	89	THR	CA-CB-OG1	-5.06	102.01	109.60
2	H	201	ALA	N-CA-C	5.06	117.53	111.71
2	H	148	GLU	N-CA-C	-5.05	103.31	110.29
2	H	103	TRP	CG-CD2-CE3	5.05	138.95	133.90
2	H	118	PRO	CA-C-N	5.05	124.98	119.78
2	H	118	PRO	C-N-CA	5.05	124.98	119.78
1	L	206	VAL	O-C-N	-5.04	117.82	123.26
1	L	73	LEU	N-CA-C	-5.04	101.48	109.59
1	L	137	ASN	CA-C-N	5.01	129.73	122.36
1	L	137	ASN	C-N-CA	5.01	129.73	122.36
1	L	137	ASN	CA-CB-CG	5.01	117.61	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	186	SER	Peptide

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	L	1681	0	1617	28	0
2	H	1604	0	1570	35	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	P	92	0	92	3	0
4	L	13	0	5	1	0
5	H	78	0	0	7	0
5	L	66	0	0	2	1
5	P	7	0	0	0	0
All	All	3541	0	3284	63	1

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 9.

All (63) 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:H:116:THR:CG2	5:H:268:HOH:O	1.82	1.26
2:H:116:THR:HG21	5:H:268:HOH:O	1.43	1.07
2:H:143:LYS:NZ	5:H:267:HOH:O	1.80	1.04
2:H:116:THR:HG22	5:H:268:HOH:O	1.57	0.88
1:L:141:PRO:HD2	1:L:199:LYS:HD3	1.63	0.80
2:H:188:ARG:HH11	2:H:188:ARG:HB3	1.48	0.79
2:H:127:GLY:N	5:H:234:HOH:O	2.18	0.76
1:L:107:LYS:HA	1:L:140:TYR:OH	1.88	0.72
1:L:205:ILE:HD12	5:L:214:HOH:O	1.88	0.72
2:H:138:LEU:HD11	2:H:188:ARG:HG3	1.74	0.69
1:L:80:ALA:HA	1:L:106:ILE:HD13	1.76	0.68
1:L:185:GLU:HA	1:L:188:TRP:CD1	2.30	0.67
2:H:134:SER:N	2:H:185:SER:HG	1.94	0.65
2:H:121:VAL:HG23	2:H:208:LYS:HG3	1.79	0.65
1:L:190:ASN:HB3	5:L:274:HOH:O	1.97	0.64
2:H:171:GLN:NE2	5:H:267:HOH:O	2.29	0.64
1:L:4:MET:HE3	1:L:23:CYS:SG	2.40	0.61
1:L:185:GLU:HA	1:L:188:TRP:HD1	1.66	0.60
1:L:199:LYS:HD2	1:L:199:LYS:H	1.67	0.60
2:H:52:ASN:HD21	2:H:53:TYR:HB3	1.67	0.60
2:H:87:THR:HG23	2:H:110:THR:HA	1.85	0.58
1:L:147:LYS:HE3	1:L:154:GLU:HB2	1.84	0.58
1:L:34:GLU:HG3	1:L:49:TYR:HA	1.86	0.57
2:H:32:TYR:CE2	2:H:96:SER:HB3	2.40	0.57
1:L:115:VAL:HA	1:L:135:PHE:O	2.04	0.57
1:L:12:PRO:HA	1:L:105:GLU:O	2.06	0.56
2:H:50:TRP:CD2	3:P:5:GLY:HA2	2.41	0.55
2:H:6:GLN:NE2	2:H:106:GLY:H	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:188:ARG:HD2	2:H:189:PRO:HA	1.90	0.53
1:L:189:HIS:HB2	1:L:192:TYR:OH	2.10	0.52
2:H:184:PRO:HB2	2:H:187:PRO:HD2	1.92	0.52
2:H:6:GLN:HE21	2:H:104:GLY:HA3	1.76	0.51
2:H:81:GLN:NE2	2:H:82(A):ASN:HD21	2.08	0.51
1:L:16:GLY:O	1:L:77:ARG:HA	2.11	0.51
1:L:147:LYS:HZ3	1:L:147:LYS:HA	1.76	0.51
2:H:52:ASN:ND2	2:H:54:SER:H	2.09	0.51
1:L:150:ILE:HD11	1:L:179:LEU:HD11	1.93	0.50
2:H:134:SER:O	2:H:184:PRO:HA	2.11	0.50
2:H:16:GLU:O	2:H:82(C):LEU:HB2	2.12	0.50
1:L:108:ARG:HD2	1:L:171:SER:HB2	1.95	0.49
1:L:199:LYS:HD2	1:L:199:LYS:N	2.28	0.48
1:L:31:HIS:NE2	4:L:212:CIT:H22	2.28	0.47
1:L:108:ARG:CD	1:L:171:SER:HB2	2.45	0.47
2:H:17:THR:HG23	2:H:81:GLN:HE21	1.80	0.46
2:H:138:LEU:HD22	2:H:210:ILE:HG21	1.97	0.46
2:H:52:ASN:ND2	2:H:53:TYR:HB3	2.30	0.46
2:H:52:ASN:HD22	2:H:53:TYR:H	1.64	0.46
1:L:119:PRO:HG3	1:L:209:PHE:CD2	2.51	0.45
1:L:44:PRO:HB2	2:H:103:TRP:CE2	2.52	0.45
3:P:13:LYS:HB3	3:P:13:LYS:NZ	2.32	0.45
2:H:4:LEU:HD22	2:H:34:MET:HE1	2.00	0.43
2:H:48:MET:HE1	2:H:80:LEU:HD11	2.00	0.43
1:L:167:ASP:O	1:L:171:SER:HA	2.19	0.43
2:H:50:TRP:CE2	3:P:5:GLY:HA2	2.54	0.43
1:L:25:SER:OG	1:L:69:THR:HA	2.19	0.42
2:H:171:GLN:O	2:H:171:GLN:HG3	2.19	0.42
2:H:205:LYS:HB3	5:H:241:HOH:O	2.18	0.42
2:H:184:PRO:O	2:H:187:PRO:HB2	2.19	0.42
2:H:121:VAL:CG2	2:H:208:LYS:HG3	2.47	0.42
1:L:66:GLY:HA3	1:L:71:PHE:HA	2.01	0.41
1:L:110:ASP:HB3	1:L:200:THR:HG22	2.02	0.41
1:L:75:ILE:HD12	1:L:75:ILE:N	2.36	0.41
2:H:22:CYS:HB3	2:H:78:ALA:HB3	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:159:LEU:CD1	5:L:256:HOH:O[1_556]	2.14	0.06

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	L	214/216 (99%)	202 (94%)	8 (4%)	4 (2%)	6	5
2	H	206/210 (98%)	194 (94%)	10 (5%)	2 (1%)	12	15
3	P	10/15 (67%)	8 (80%)	2 (20%)	0	100	100
All	All	430/441 (98%)	404 (94%)	20 (5%)	6 (1%)	9	9

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	68	GLY
2	H	172	SER
1	L	77	ARG
1	L	199	LYS
2	H	41	PRO
1	L	119	PRO

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	L	194/194 (100%)	170 (88%)	24 (12%)	4	5
2	H	182/182 (100%)	150 (82%)	32 (18%)	2	2
3	P	11/13 (85%)	10 (91%)	1 (9%)	9	11
All	All	387/389 (100%)	330 (85%)	57 (15%)	3	3

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	3	LEU
1	L	17	ASP
1	L	20	SER
1	L	27	GLN
1	L	34	GLU
1	L	56	SER
1	L	65	SER
1	L	67	SER
1	L	70	ASP
1	L	77	ARG
1	L	79	GLU
1	L	81	GLU
1	L	83	LEU
1	L	104	LEU
1	L	108	ARG
1	L	147	LYS
1	L	154	GLU
1	L	157	ASN
1	L	169	LYS
1	L	179	LEU
1	L	185	GLU
1	L	191	SER
1	L	199	LYS
1	L	203	SER
2	H	1	GLN
2	H	12	LYS
2	H	13	THR
2	H	16	GLU
2	H	17	THR
2	H	18	VAL
2	H	37	VAL
2	H	40	THR
2	H	43	LYS
2	H	50	TRP
2	H	52	ASN
2	H	52(A)	THR
2	H	61	ASP
2	H	72	GLU
2	H	73	THR
2	H	74	SER
2	H	80	LEU
2	H	116	THR

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Mol	Chain	Res	Type
2	H	121	VAL
2	H	134	SER
2	H	136	VAL
2	H	140	CYS
2	H	150	VAL
2	H	160	SER
2	H	177	LEU
2	H	178	SER
2	H	183	VAL
2	H	185	SER
2	H	187	PRO
2	H	188	ARG
2	H	190	SER
2	H	205	LYS
3	P	14	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	42	GLN
2	H	6	GLN
2	H	52	ASN
2	H	81	GLN
2	H	82(A)	ASN
2	H	164	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

1 ligand is 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
4	CIT	L	212	-	12,12,12	1.45	2 (16%)	17,17,17	4.16	10 (58%)

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
4	CIT	L	212	-	-	4/16/16/16	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	212	CIT	O7-C3	-2.13	1.39	1.43
4	L	212	CIT	O4-C5	-2.10	1.23	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	212	CIT	C4-C3-C2	-10.23	83.07	109.31
4	L	212	CIT	O7-C3-C2	7.17	125.73	109.38
4	L	212	CIT	O7-C3-C4	-7.17	93.03	109.38
4	L	212	CIT	O7-C3-C6	4.49	115.33	108.96
4	L	212	CIT	C2-C3-C6	3.58	117.94	110.03
4	L	212	CIT	C3-C4-C5	3.37	123.12	113.92
4	L	212	CIT	O6-C6-C3	3.30	119.47	113.14
4	L	212	CIT	O4-C5-C4	2.39	121.92	114.35
4	L	212	CIT	C4-C3-C6	-2.32	104.91	110.03
4	L	212	CIT	O1-C1-C2	-2.07	117.09	122.95

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	L	212	CIT	C2-C3-C6-O5
4	L	212	CIT	C2-C3-C6-O6
4	L	212	CIT	C1-C2-C3-O7
4	L	212	CIT	C3-C4-C5-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	212	CIT	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

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.