



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:18 PM UTC

PDB ID : 1ACC / pdb_00001acc
Title : ANTHRAX PROTECTIVE ANTIGEN
Authors : Petosa, C.; Liddington, R.C.
Deposited on : 1997-02-05
Resolution : 2.10 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)	:	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

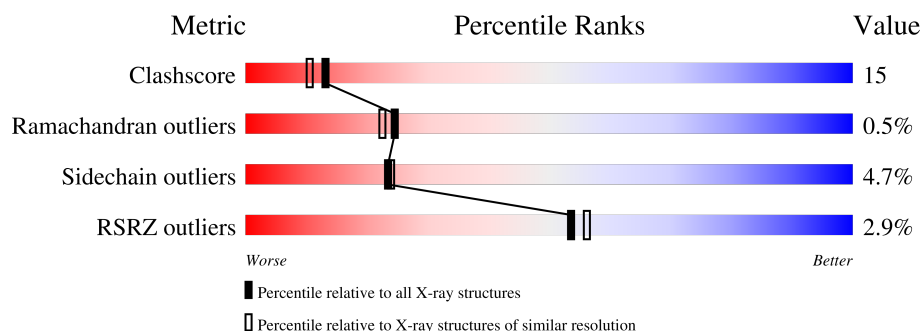
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	735	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5672 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 ANTHRAX PROTECTIVE ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	665	Total	C	N	O	S	0	0	0
			5282	3317	898	1060	7			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		

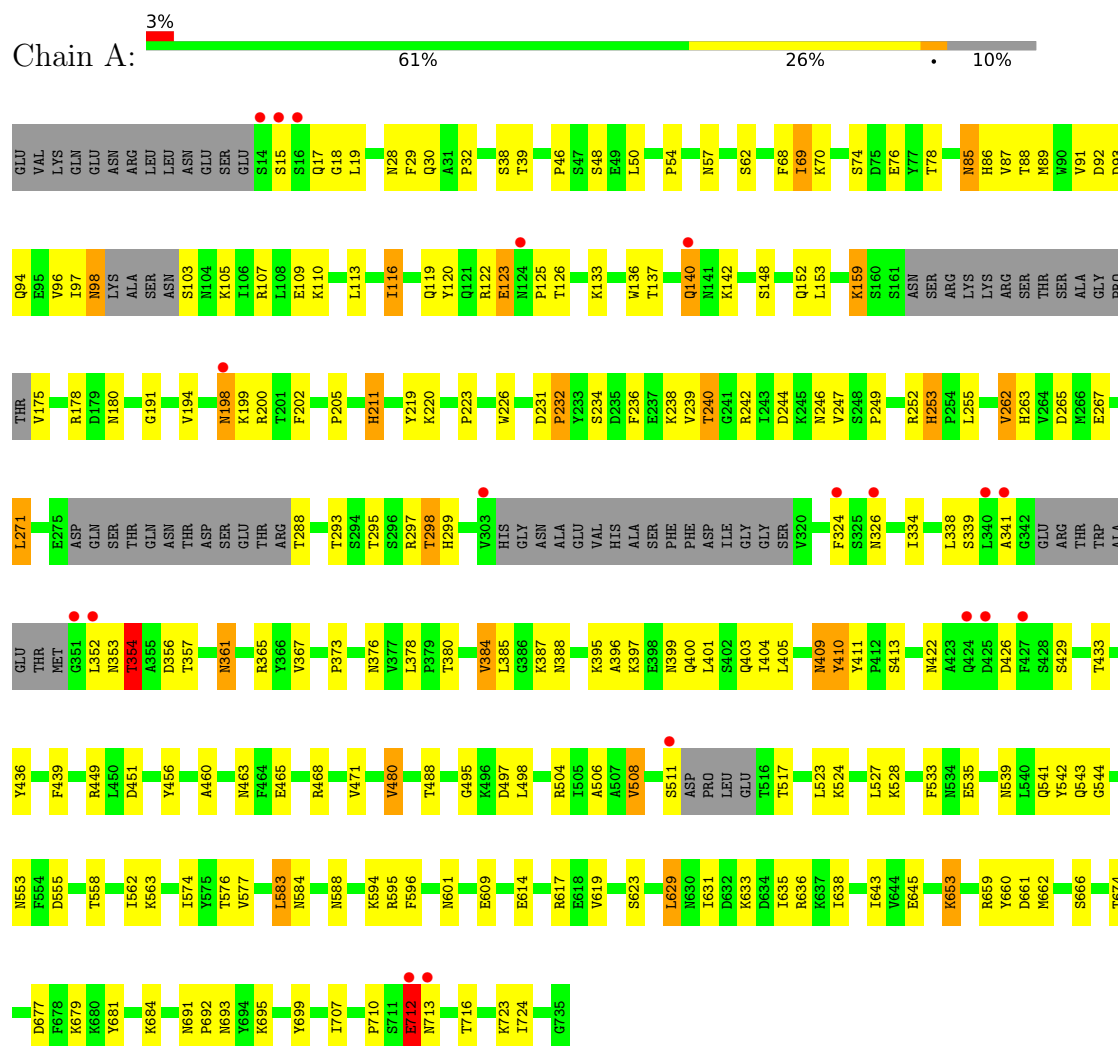
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	388	Total	O	0	0
			388	388		

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: ANTHRAX PROTECTIVE ANTIGEN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.30Å 93.70Å 82.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.10 6.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.4 (6.00-2.10) 92.1 (6.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.40 (at 1.89Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.223 , 0.296 0.223 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.411	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.69 , 128.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5672	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.48	0/5371	1.03	33/7271 (0.5%)

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
1	A	0	1

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	211	HIS	N-CA-C	9.11	120.82	111.07
1	A	126	THR	N-CA-C	-7.42	104.47	113.97
1	A	480	VAL	N-CA-C	6.85	117.82	111.45
1	A	543	GLN	N-CA-C	6.75	120.70	111.39
1	A	339	SER	N-CA-C	-6.51	105.97	114.04
1	A	123	GLU	N-CA-C	6.27	118.92	111.33
1	A	712	GLU	N-CA-C	6.09	119.26	108.24
1	A	677	ASP	N-CA-C	6.06	118.07	108.67
1	A	596	PHE	N-CA-C	5.91	119.19	109.85
1	A	232	PRO	N-CA-C	5.85	120.76	114.68
1	A	495	GLY	N-CA-C	5.80	120.84	113.24
1	A	140	GLN	N-CA-C	-5.74	102.96	110.53
1	A	253	HIS	N-CA-C	-5.67	102.63	109.72
1	A	555	ASP	N-CA-C	-5.64	103.25	110.53
1	A	384	VAL	N-CA-C	5.62	116.57	108.48
1	A	596	PHE	CA-C-O	-5.47	115.06	121.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	653	LYS	N-CA-C	5.47	118.49	109.85
1	A	297	ARG	N-CA-C	-5.31	102.53	110.23
1	A	376	ASN	N-CA-C	5.30	120.92	113.02
1	A	15	SER	N-CA-C	5.30	119.74	113.28
1	A	679	LYS	N-CA-C	5.29	117.12	111.36
1	A	410	TYR	N-CA-C	5.29	118.20	109.85
1	A	175	VAL	CA-C-N	5.23	126.37	119.84
1	A	175	VAL	C-N-CA	5.23	126.37	119.84
1	A	594	LYS	N-CA-C	5.22	118.91	112.54
1	A	116	ILE	CB-CA-C	-5.18	102.83	110.82
1	A	456	TYR	N-CA-C	-5.17	102.37	110.17
1	A	596	PHE	CA-C-N	5.16	129.47	122.19
1	A	596	PHE	C-N-CA	5.16	129.47	122.19
1	A	574	ILE	N-CA-C	5.14	116.60	111.00
1	A	712	GLU	CB-CA-C	-5.08	104.33	112.05
1	A	508	VAL	N-CA-C	5.07	116.10	109.21
1	A	262	VAL	N-CA-C	5.06	115.77	108.48

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	712	GLU	Mainchain

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	5282	0	5223	157	0
2	A	2	0	0	0	0
3	A	388	0	0	23	0
All	All	5672	0	5223	157	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 15.

All (157) 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:409:ASN:HD22	1:A:410:TYR:H	0.98	0.92
1:A:326:ASN:HD21	1:A:504:ARG:HH21	1.20	0.87
1:A:352:LEU:HD12	1:A:356:ASP:HB2	1.59	0.84
1:A:352:LEU:CD1	1:A:356:ASP:HB2	2.09	0.82
1:A:240:THR:HG21	1:A:242:ARG:HH11	1.44	0.82
1:A:298:THR:OG1	1:A:601:ASN:HB3	1.78	0.82
1:A:199:LYS:HG3	1:A:200:ARG:HG3	1.62	0.82
1:A:326:ASN:ND2	1:A:504:ARG:HH21	1.81	0.79
1:A:253:HIS:HD2	1:A:255:LEU:H	1.29	0.78
1:A:231:ASP:HB2	1:A:232:PRO:HD2	1.66	0.75
1:A:409:ASN:HD22	1:A:410:TYR:N	1.79	0.74
1:A:76:GLU:HB3	1:A:105:LYS:HD3	1.71	0.72
1:A:710:PRO:HG3	1:A:716:THR:HG22	1.72	0.71
1:A:17:GLN:O	1:A:39:THR:HA	1.90	0.71
1:A:409:ASN:ND2	1:A:410:TYR:H	1.82	0.70
1:A:629:LEU:HD12	1:A:631:ILE:HD11	1.72	0.70
1:A:404:ILE:HG12	3:A:1136:HOH:O	1.91	0.69
1:A:263:HIS:HE1	3:A:920:HOH:O	1.76	0.69
1:A:361:ASN:HD21	1:A:422:ASN:H	1.42	0.67
1:A:643:ILE:HD12	1:A:723:LYS:HD3	1.76	0.67
1:A:633:LYS:HE3	3:A:1162:HOH:O	1.94	0.67
1:A:357:THR:CG2	1:A:433:THR:HG23	2.24	0.66
1:A:326:ASN:HD21	1:A:504:ARG:NH2	1.91	0.66
1:A:263:HIS:HD2	3:A:867:HOH:O	1.78	0.65
1:A:691:ASN:ND2	1:A:692:PRO:HD2	2.13	0.64
1:A:69:ILE:HG12	1:A:116:ILE:HD11	1.80	0.63
1:A:352:LEU:HG	1:A:436:TYR:HD1	1.62	0.63
1:A:244:ASP:HB3	1:A:247:VAL:HG23	1.82	0.62
1:A:357:THR:HG23	1:A:433:THR:HG23	1.81	0.61
1:A:70:LYS:HB3	1:A:113:LEU:HD23	1.81	0.61
1:A:226:TRP:CE2	1:A:234:SER:HB3	2.36	0.60
1:A:341:ALA:HB2	3:A:878:HOH:O	2.02	0.60
1:A:18:GLY:O	1:A:153:LEU:HD12	2.01	0.60
1:A:159:LYS:HD2	1:A:180:ASN:HB2	1.84	0.60
1:A:271:LEU:HD23	1:A:439:PHE:HE1	1.69	0.58
1:A:449:ARG:HD2	1:A:451:ASP:OD1	2.04	0.57
1:A:98:ASN:O	1:A:98:ASN:ND2	2.35	0.57
1:A:123:GLU:C	1:A:125:PRO:HD3	2.30	0.57
1:A:660:TYR:HB2	1:A:707:ILE:HG13	1.86	0.56
1:A:29:PHE:CE1	1:A:62:SER:HB2	2.41	0.56
1:A:140:GLN:HG2	3:A:1092:HOH:O	2.05	0.56
1:A:232:PRO:HD3	1:A:480:VAL:HG11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:LEU:HD12	1:A:356:ASP:CB	2.34	0.55
1:A:265:ASP:OD1	1:A:295:THR:HB	2.07	0.55
1:A:352:LEU:HD11	1:A:356:ASP:HB2	1.87	0.55
1:A:249:PRO:HA	1:A:252:ARG:HD2	1.89	0.55
1:A:712:GLU:O	1:A:712:GLU:HG2	2.08	0.54
1:A:635:ILE:O	1:A:638:ILE:HG13	2.09	0.53
1:A:380:THR:HG23	1:A:395:LYS:HG2	1.89	0.53
1:A:583:LEU:C	1:A:584:ASN:HD22	2.15	0.53
1:A:198:ASN:HD22	1:A:199:LYS:N	2.07	0.53
1:A:636:ARG:NH1	1:A:674:THR:HG23	2.23	0.53
1:A:76:GLU:HB3	1:A:105:LYS:CD	2.39	0.52
1:A:541:GLN:NE2	1:A:544:GLY:H	2.07	0.52
1:A:684:LYS:HD2	1:A:684:LYS:N	2.25	0.52
1:A:96:VAL:O	1:A:103:SER:HA	2.10	0.51
1:A:18:GLY:HA3	1:A:38:SER:O	2.11	0.51
1:A:298:THR:HG23	3:A:951:HOH:O	2.10	0.51
1:A:238:LYS:HD2	3:A:945:HOH:O	2.11	0.51
1:A:122:ARG:HG2	3:A:1000:HOH:O	2.10	0.51
1:A:85:ASN:HD22	1:A:85:ASN:N	2.08	0.50
1:A:385:LEU:HD12	3:A:833:HOH:O	2.10	0.50
1:A:365:ARG:NH2	1:A:413:SER:O	2.44	0.50
1:A:54:PRO:HG2	1:A:57:ASN:OD1	2.11	0.50
1:A:240:THR:HG23	1:A:242:ARG:HG3	1.92	0.50
1:A:468:ARG:HG2	3:A:1074:HOH:O	2.11	0.50
1:A:645:GLU:OE2	1:A:653:LYS:HG2	2.11	0.50
1:A:125:PRO:HG3	3:A:1000:HOH:O	2.12	0.49
1:A:636:ARG:HG3	3:A:941:HOH:O	2.13	0.49
1:A:352:LEU:CD1	1:A:356:ASP:CB	2.87	0.49
1:A:231:ASP:HB2	1:A:232:PRO:CD	2.39	0.49
1:A:691:ASN:OD1	1:A:693:ASN:HB2	2.13	0.49
1:A:92:ASP:O	1:A:93:ASP:HB2	2.13	0.48
1:A:97:ILE:HG22	1:A:98:ASN:N	2.28	0.48
1:A:539:ASN:HB3	3:A:1110:HOH:O	2.12	0.48
1:A:397:LYS:HE3	1:A:400:GLN:OE1	2.13	0.48
1:A:76:GLU:HB3	1:A:105:LYS:CG	2.43	0.48
1:A:488:THR:HG22	1:A:506:ALA:HA	1.96	0.48
1:A:85:ASN:H	1:A:85:ASN:ND2	2.12	0.48
1:A:68:PHE:HB2	1:A:152:GLN:HG2	1.96	0.48
1:A:246:ASN:HB2	1:A:373:PRO:HG3	1.96	0.47
1:A:387:LYS:O	1:A:449:ARG:NH2	2.46	0.47
1:A:46:PRO:HB2	1:A:48:SER:OG	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:GLU:CB	1:A:105:LYS:HD3	2.42	0.47
1:A:91:VAL:O	1:A:94:GLN:HB3	2.15	0.47
1:A:523:LEU:O	1:A:527:LEU:HD13	2.15	0.47
1:A:85:ASN:N	1:A:85:ASN:ND2	2.61	0.47
1:A:159:LYS:CD	1:A:180:ASN:HB2	2.45	0.47
1:A:352:LEU:HG	1:A:436:TYR:CD1	2.47	0.47
1:A:354:THR:HA	1:A:436:TYR:CG	2.50	0.47
1:A:623:SER:HB2	3:A:803:HOH:O	2.14	0.47
1:A:659:ARG:HD3	1:A:662:MET:SD	2.55	0.47
1:A:137:THR:HA	1:A:142:LYS:O	2.16	0.46
1:A:533:PHE:CE2	1:A:542:TYR:HB2	2.50	0.46
1:A:463:ASN:OD1	1:A:465:GLU:HB2	2.15	0.46
1:A:74:SER:OG	1:A:109:GLU:HA	2.14	0.46
1:A:395:LYS:HA	3:A:885:HOH:O	2.15	0.46
1:A:78:THR:O	1:A:136:TRP:HA	2.16	0.46
1:A:399:ASN:ND2	3:A:1143:HOH:O	2.49	0.45
1:A:645:GLU:CD	1:A:653:LYS:HG2	2.41	0.45
1:A:324:PHE:HD1	1:A:588:ASN:ND2	2.15	0.45
1:A:62:SER:HA	1:A:120:TYR:O	2.17	0.45
1:A:636:ARG:HH12	1:A:666:SER:HB2	1.82	0.45
1:A:357:THR:HG21	1:A:433:THR:HG23	1.96	0.44
1:A:405:LEU:HD13	1:A:411:TYR:HB2	1.98	0.44
1:A:74:SER:HB3	1:A:107:ARG:NH2	2.32	0.44
1:A:361:ASN:HD21	1:A:422:ASN:N	2.13	0.44
1:A:123:GLU:O	1:A:125:PRO:HD3	2.17	0.44
1:A:240:THR:CG2	1:A:242:ARG:HH11	2.24	0.44
1:A:293:THR:HG22	1:A:334:ILE:HA	2.00	0.44
1:A:69:ILE:HG12	1:A:116:ILE:CG1	2.48	0.44
1:A:69:ILE:HG12	1:A:116:ILE:CD1	2.46	0.44
1:A:191:GLY:HA2	1:A:219:TYR:O	2.18	0.44
1:A:29:PHE:CZ	1:A:119:GLN:HG3	2.53	0.43
1:A:30:GLN:O	1:A:32:PRO:HD3	2.18	0.43
1:A:262:VAL:HA	1:A:367:VAL:O	2.18	0.43
1:A:614:GLU:OE2	1:A:617:ARG:HD2	2.18	0.43
1:A:576:THR:HG22	1:A:577:VAL:HG13	2.01	0.43
1:A:619:VAL:HG22	1:A:629:LEU:HD22	2.00	0.43
1:A:89:MET:HE3	1:A:116:ILE:HG23	2.00	0.43
1:A:636:ARG:HG2	3:A:939:HOH:O	2.17	0.43
1:A:70:LYS:HB3	1:A:113:LEU:CD2	2.45	0.43
1:A:78:THR:HA	3:A:1002:HOH:O	2.19	0.43
1:A:85:ASN:ND2	1:A:86:HIS:ND1	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:ARG:HH21	1:A:223:PRO:CG	2.32	0.43
1:A:384:VAL:HG11	1:A:449:ARG:NH2	2.34	0.43
1:A:460:ALA:HA	1:A:471:VAL:HA	2.00	0.43
1:A:403:GLN:HE22	1:A:413:SER:H	1.67	0.42
1:A:299:HIS:HB2	3:A:850:HOH:O	2.19	0.42
1:A:699:TYR:HB3	1:A:723:LYS:HB2	2.02	0.42
1:A:133:LYS:HE3	3:A:1010:HOH:O	2.19	0.42
1:A:267:GLU:OE2	1:A:365:ARG:HD2	2.19	0.42
1:A:558:THR:O	1:A:562:ILE:HG12	2.19	0.42
1:A:205:PRO:O	1:A:211:HIS:HE1	2.02	0.42
1:A:524:LYS:O	1:A:528:LYS:HG3	2.20	0.42
1:A:338:LEU:HG	1:A:661:ASP:OD2	2.19	0.41
1:A:19:LEU:HD21	1:A:116:ILE:HD12	2.01	0.41
1:A:271:LEU:HD23	1:A:439:PHE:CE1	2.52	0.41
1:A:70:LYS:HD3	1:A:152:GLN:NE2	2.34	0.41
1:A:87:VAL:CG1	1:A:88:THR:N	2.82	0.41
1:A:378:LEU:HB2	1:A:396:ALA:HB3	2.02	0.41
1:A:397:LYS:HB2	1:A:400:GLN:HB2	2.02	0.41
1:A:609:GLU:HG2	1:A:724:ILE:HG12	2.02	0.41
1:A:194:VAL:HA	1:A:202:PHE:O	2.21	0.41
1:A:236:PHE:O	1:A:240:THR:HB	2.21	0.41
1:A:50:LEU:HA	1:A:50:LEU:HD23	1.83	0.41
1:A:86:HIS:CD2	1:A:122:ARG:HD2	2.56	0.41
1:A:240:THR:HG21	1:A:242:ARG:NH1	2.24	0.41
1:A:595:ARG:HD2	3:A:998:HOH:O	2.19	0.41
1:A:253:HIS:CD2	1:A:255:LEU:H	2.20	0.40
1:A:662:MET:HE1	3:A:962:HOH:O	2.20	0.40
1:A:240:THR:HG21	1:A:242:ARG:HD2	2.04	0.40
1:A:497:ASP:O	1:A:498:LEU:HB2	2.20	0.40
1:A:601:ASN:HD22	1:A:601:ASN:HA	1.72	0.40
1:A:662:MET:HE2	1:A:681:TYR:CD1	2.57	0.40
1:A:220:LYS:HB3	1:A:220:LYS:NZ	2.36	0.40
1:A:194:VAL:HG21	1:A:239:VAL:O	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	651/735 (89%)	617 (95%)	31 (5%)	3 (0%)	24	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	354	THR
1	A	426	ASP
1	A	713	ASN

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	600/661 (91%)	572 (95%)	28 (5%)	23	24

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	69	ILE
1	A	85	ASN
1	A	98	ASN
1	A	110	LYS
1	A	148	SER
1	A	159	LYS
1	A	198	ASN

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Mol	Chain	Res	Type
1	A	240	THR
1	A	271	LEU
1	A	288	THR
1	A	298	THR
1	A	353	ASN
1	A	354	THR
1	A	361	ASN
1	A	388	ASN
1	A	401	LEU
1	A	409	ASN
1	A	429	SER
1	A	508	VAL
1	A	511	SER
1	A	517	THR
1	A	535	GLU
1	A	553	ASN
1	A	563	LYS
1	A	583	LEU
1	A	629	LEU
1	A	695	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	85	ASN
1	A	94	GLN
1	A	124	ASN
1	A	150	ASN
1	A	152	GLN
1	A	180	ASN
1	A	198	ASN
1	A	246	ASN
1	A	253	HIS
1	A	263	HIS
1	A	326	ASN
1	A	361	ASN
1	A	388	ASN
1	A	403	GLN
1	A	409	ASN
1	A	424	GLN
1	A	541	GLN

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Mol	Chain	Res	Type
1	A	584	ASN
1	A	601	ASN
1	A	697	ASN
1	A	705	ASN
1	A	709	ASN
1	A	713	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	665/735 (90%)	-0.31	19 (2%) 53 56	6, 24, 57, 87	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	713	ASN	6.6
1	A	351	GLY	4.5
1	A	15	SER	3.6
1	A	16	SER	3.1
1	A	124	ASN	2.9
1	A	140	GLN	2.9
1	A	14	SER	2.8
1	A	712	GLU	2.7
1	A	341	ALA	2.6
1	A	427	PHE	2.6
1	A	511	SER	2.5
1	A	340	LEU	2.5
1	A	303	VAL	2.5
1	A	198	ASN	2.4
1	A	425	ASP	2.3
1	A	352	LEU	2.2
1	A	326	ASN	2.1
1	A	324	PHE	2.1
1	A	424	GLN	2.1

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
2	CA	A	801	1/1	0.97	0.02	13,13,13,13	0
2	CA	A	800	1/1	1.00	0.02	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.