



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:31 PM UTC

PDB ID : 4OB1 / pdb_00004ob1
Title : Crystal Structure of Nitrile Hydratase from *Pseudonocardia thermophila*
bound to Butaneboronic Acid via Co-crystallization
Authors : Rui, W.; Salette, M.; Ruslan, S.; Richard, H.; Dali, L.
Deposited on : 2014-01-06
Resolution : 1.63 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

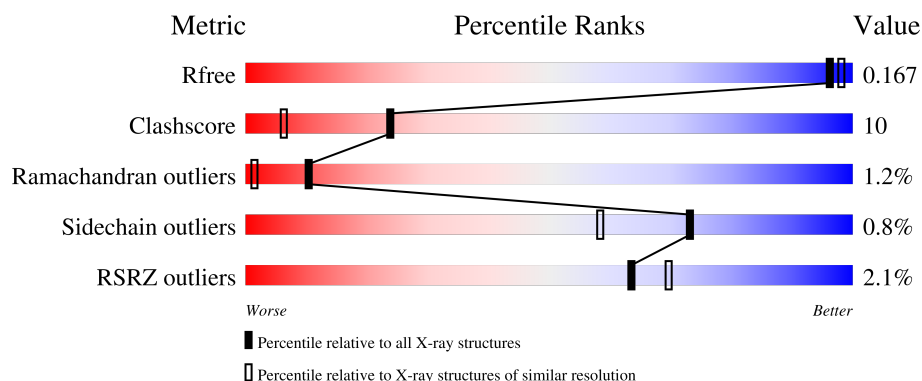
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	290	<div> <div>%</div> <div> <div></div> <div>62%</div> <div>7%</div> <div>30%</div> </div> </div>
2	B	233	<div> <div>3%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>...</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4310 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 Cobalt-containing nitrile hydratase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	7	0
			1646	1049	271	315	11			

There are 86 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-79	MET	-	initiating methionine	UNP Q7SID2
A	-78	THR	-	expression tag	UNP Q7SID2
A	-77	GLU	-	expression tag	UNP Q7SID2
A	-76	ASN	-	expression tag	UNP Q7SID2
A	-75	ILE	-	expression tag	UNP Q7SID2
A	-74	LEU	-	expression tag	UNP Q7SID2
A	-73	ARG	-	expression tag	UNP Q7SID2
A	-72	LYS	-	expression tag	UNP Q7SID2
A	-71	SER	-	expression tag	UNP Q7SID2
A	-70	ASP	-	expression tag	UNP Q7SID2
A	-69	GLU	-	expression tag	UNP Q7SID2
A	-68	GLU	-	expression tag	UNP Q7SID2
A	-67	ILE	-	expression tag	UNP Q7SID2
A	-66	GLN	-	expression tag	UNP Q7SID2
A	-65	LYS	-	expression tag	UNP Q7SID2
A	-64	GLU	-	expression tag	UNP Q7SID2
A	-63	ILE	-	expression tag	UNP Q7SID2
A	-62	THR	-	expression tag	UNP Q7SID2
A	-61	ALA	-	expression tag	UNP Q7SID2
A	-60	ARG	-	expression tag	UNP Q7SID2
A	-59	VAL	-	expression tag	UNP Q7SID2
A	-58	LYS	-	expression tag	UNP Q7SID2
A	-57	ALA	-	expression tag	UNP Q7SID2
A	-56	LEU	-	expression tag	UNP Q7SID2
A	-55	GLU	-	expression tag	UNP Q7SID2
A	-54	SER	-	expression tag	UNP Q7SID2
A	-53	MET	-	expression tag	UNP Q7SID2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-52	LEU	-	expression tag	UNP Q7SID2
A	-51	ILE	-	expression tag	UNP Q7SID2
A	-50	GLU	-	expression tag	UNP Q7SID2
A	-49	GLN	-	expression tag	UNP Q7SID2
A	-48	GLY	-	expression tag	UNP Q7SID2
A	-47	ILE	-	expression tag	UNP Q7SID2
A	-46	LEU	-	expression tag	UNP Q7SID2
A	-45	THR	-	expression tag	UNP Q7SID2
A	-44	THR	-	expression tag	UNP Q7SID2
A	-43	SER	-	expression tag	UNP Q7SID2
A	-42	MET	-	expression tag	UNP Q7SID2
A	-41	ILE	-	expression tag	UNP Q7SID2
A	-40	ASP	-	expression tag	UNP Q7SID2
A	-39	ARG	-	expression tag	UNP Q7SID2
A	-38	MET	-	expression tag	UNP Q7SID2
A	-37	ALA	-	expression tag	UNP Q7SID2
A	-36	GLU	-	expression tag	UNP Q7SID2
A	-35	ILE	-	expression tag	UNP Q7SID2
A	-34	TYR	-	expression tag	UNP Q7SID2
A	-33	GLU	-	expression tag	UNP Q7SID2
A	-32	ASN	-	expression tag	UNP Q7SID2
A	-31	GLU	-	expression tag	UNP Q7SID2
A	-30	VAL	-	expression tag	UNP Q7SID2
A	-29	GLY	-	expression tag	UNP Q7SID2
A	-28	PRO	-	expression tag	UNP Q7SID2
A	-27	HIS	-	expression tag	UNP Q7SID2
A	-26	LEU	-	expression tag	UNP Q7SID2
A	-25	GLY	-	expression tag	UNP Q7SID2
A	-24	ALA	-	expression tag	UNP Q7SID2
A	-23	LYS	-	expression tag	UNP Q7SID2
A	-22	VAL	-	expression tag	UNP Q7SID2
A	-21	VAL	-	expression tag	UNP Q7SID2
A	-20	VAL	-	expression tag	UNP Q7SID2
A	-19	LYS	-	expression tag	UNP Q7SID2
A	-18	ALA	-	expression tag	UNP Q7SID2
A	-17	TRP	-	expression tag	UNP Q7SID2
A	-16	THR	-	expression tag	UNP Q7SID2
A	-15	ASP	-	expression tag	UNP Q7SID2
A	-14	PRO	-	expression tag	UNP Q7SID2
A	-13	GLU	-	expression tag	UNP Q7SID2
A	-12	PHE	-	expression tag	UNP Q7SID2
A	-11	LYS	-	expression tag	UNP Q7SID2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	LYS	-	expression tag	UNP Q7SID2
A	-9	ARG	-	expression tag	UNP Q7SID2
A	-8	LEU	-	expression tag	UNP Q7SID2
A	-7	LEU	-	expression tag	UNP Q7SID2
A	-6	ALA	-	expression tag	UNP Q7SID2
A	-5	ASP	-	expression tag	UNP Q7SID2
A	-4	GLY	-	expression tag	UNP Q7SID2
A	-3	THR	-	expression tag	UNP Q7SID2
A	-2	GLU	-	expression tag	UNP Q7SID2
A	-1	ALA	-	expression tag	UNP Q7SID2
A	0	CYS	-	expression tag	UNP Q7SID2
A	205	HIS	-	expression tag	UNP Q7SID2
A	206	HIS	-	expression tag	UNP Q7SID2
A	207	HIS	-	expression tag	UNP Q7SID2
A	208	HIS	-	expression tag	UNP Q7SID2
A	209	HIS	-	expression tag	UNP Q7SID2
A	210	HIS	-	expression tag	UNP Q7SID2

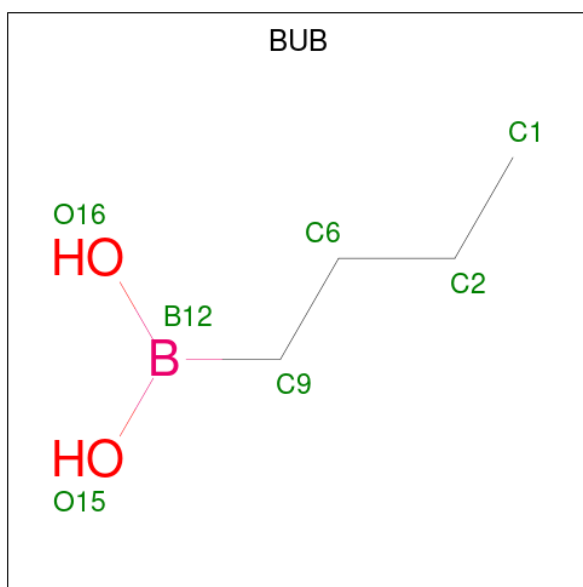
- Molecule 2 is a protein called Cobalt-containing nitrile hydratase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	228	Total	C	N	O	S	0	11	0
			1908	1225	329	347	7			

- Molecule 3 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Co	0	0
			1	1		

- Molecule 4 is 1-BUTANE BORONIC ACID (CCD ID: BUB) (formula: C₄H₁₁BO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	B	C	O	0	0
			7	1	4	2		

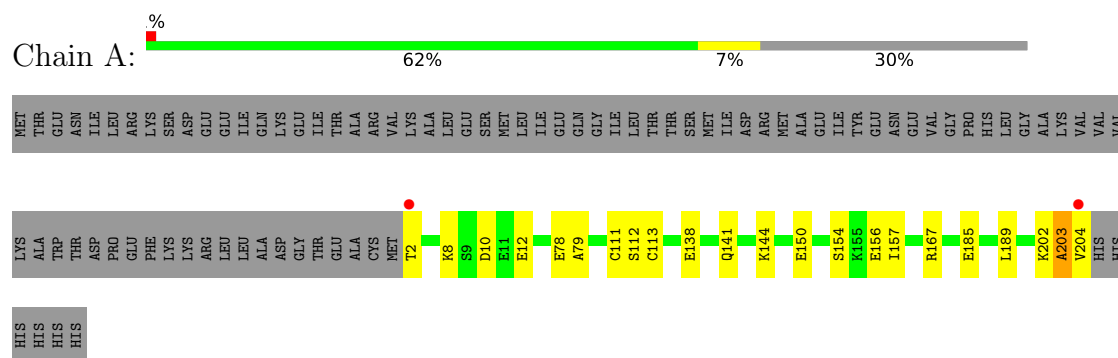
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	316	Total	O	0	0
			316	316		
5	B	432	Total	O	0	0
			432	432		

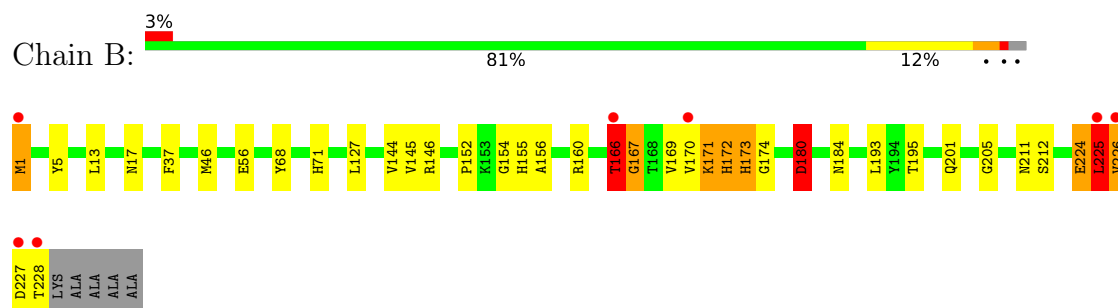
3 Residue-property plots [i](#)

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: Cobalt-containing nitrile hydratase subunit alpha



- Molecule 2: Cobalt-containing nitrile hydratase subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	65.87Å 65.87Å 186.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.06 – 1.63 36.06 – 1.63	Depositor EDS
% Data completeness (in resolution range)	100.0 (36.06-1.63) 99.9 (36.06-1.63)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.97 (at 1.63Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.3_1479)	Depositor
R, R_{free}	0.142 , 0.166 0.144 , 0.167	Depositor DCC
R_{free} test set	2998 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	11.8	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4310	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: CO, BUB, CSD

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	0.94	3/1694 (0.2%)	1.03	4/2298 (0.2%)
2	B	1.09	9/2003 (0.4%)	1.20	21/2722 (0.8%)
All	All	1.02	12/3697 (0.3%)	1.12	25/5020 (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
2	B	0	2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	226	VAL	CA-C	-12.83	1.36	1.52
2	B	225	LEU	CA-C	11.88	1.68	1.52
2	B	180	ASP	C-O	10.77	1.36	1.24
2	B	226	VAL	N-CA	7.76	1.56	1.46
2	B	225	LEU	C-N	7.31	1.43	1.33
1	A	79	ALA	C-O	6.51	1.31	1.24
1	A	185	GLU	C-O	6.47	1.31	1.24
2	B	37	PHE	CA-C	5.53	1.58	1.52
2	B	226	VAL	CA-CB	5.47	1.61	1.54
1	A	189	LEU	N-CA	5.33	1.53	1.46
2	B	68	TYR	C-O	5.17	1.30	1.24
2	B	172	HIS	CA-C	5.15	1.60	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	172	HIS	N-CA-C	9.57	124.51	113.02
2	B	225	LEU	CA-C-N	-9.37	105.10	121.97
2	B	225	LEU	C-N-CA	-9.37	105.10	121.97
2	B	226	VAL	CA-C-O	-9.08	109.43	120.78
2	B	226	VAL	CB-CA-C	-8.24	97.77	111.29
2	B	226	VAL	N-CA-CB	8.24	124.83	111.23
2	B	226	VAL	N-CA-C	-7.83	93.06	109.34
2	B	173	HIS	N-CA-C	7.70	127.19	110.80
2	B	226	VAL	O-C-N	7.68	132.17	122.57
2	B	226	VAL	CG1-CB-CG2	-7.50	94.31	110.80
2	B	169	VAL	CA-C-N	7.36	129.95	122.96
2	B	169	VAL	C-N-CA	7.36	129.95	122.96
2	B	172	HIS	O-C-N	-6.27	114.31	122.39
2	B	224	GLU	CA-C-N	-6.26	109.57	121.54
2	B	224	GLU	C-N-CA	-6.26	109.57	121.54
2	B	167	GLY	N-CA-C	6.14	117.78	111.95
1	A	203	ALA	CA-C-N	5.44	131.50	121.70
1	A	203	ALA	C-N-CA	5.44	131.50	121.70
1	A	189	LEU	N-CA-C	5.23	119.50	113.18
1	A	185	GLU	CA-C-O	5.15	126.23	120.82
2	B	205	GLY	CA-C-N	5.15	124.94	119.28
2	B	205	GLY	C-N-CA	5.15	124.94	119.28
2	B	166	THR	CB-CA-C	5.10	120.57	110.42
2	B	71	HIS	CA-C-N	-5.07	113.49	120.28
2	B	71	HIS	C-N-CA	-5.07	113.49	120.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	180	ASP	Mainchain
2	B	225	LEU	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	A	1646	0	1655	17	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1908	0	1830	55	0
3	A	1	0	0	0	0
4	A	7	0	10	0	0
5	A	316	0	0	7	0
5	B	432	0	0	15	2
All	All	4310	0	3495	71	2

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 (71) 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:146[A]:ARG:HB2	2:B:226:VAL:HG23	1.34	1.10
2:B:167:GLY:N	5:B:707:HOH:O	1.86	1.08
2:B:146[B]:ARG:HB2	2:B:226:VAL:HG23	1.38	1.02
1:A:8:LYS:NZ	5:A:536:HOH:O	1.98	0.94
2:B:17[A]:ASN:ND2	5:B:694:HOH:O	2.01	0.93
2:B:171:LYS:N	5:B:671:HOH:O	2.02	0.91
2:B:224:GLU:O	2:B:226:VAL:N	2.03	0.91
2:B:225:LEU:HB3	2:B:227:ASP:HA	1.53	0.90
2:B:172:HIS:O	2:B:174:GLY:N	2.12	0.82
1:A:2:THR:N	5:A:684:HOH:O	2.13	0.81
2:B:146[B]:ARG:H	2:B:226:VAL:H	1.29	0.80
2:B:146[A]:ARG:H	2:B:226:VAL:H	1.29	0.79
2:B:127[A]:LEU:HD23	5:B:697:HOH:O	1.86	0.76
2:B:172:HIS:O	2:B:172:HIS:CD2	2.39	0.76
2:B:224:GLU:O	2:B:225:LEU:C	2.28	0.74
2:B:166:THR:C	5:B:707:HOH:O	2.26	0.73
1:A:138:GLU:HB3	1:A:141:GLN:HG3	1.75	0.68
1:A:10[B]:ASP:OD1	5:A:472:HOH:O	2.12	0.67
2:B:170:VAL:O	2:B:195:THR:HB	1.94	0.66
2:B:144:VAL:O	2:B:226:VAL:O	2.14	0.65
2:B:1:MET:N	2:B:152:PRO:O	2.30	0.65
2:B:160:ARG:NE	5:B:464:HOH:O	2.30	0.65
2:B:144:VAL:H	2:B:228:THR:C	2.05	0.64
2:B:180:ASP:O	2:B:184:ASN:ND2	2.27	0.63
1:A:203:ALA:O	1:A:204:VAL:HG12	1.99	0.62
2:B:170:VAL:CA	5:B:671:HOH:O	2.48	0.62
2:B:1:MET:CE	2:B:156:ALA:H	2.14	0.60
2:B:166:THR:HG22	2:B:167:GLY:H	1.66	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:172:HIS:O	2:B:172:HIS:CG	2.55	0.60
2:B:1:MET:SD	2:B:5:TYR:OH	2.60	0.60
2:B:1:MET:HG3	2:B:154:GLY:O	2.02	0.59
2:B:1:MET:HE2	2:B:156:ALA:H	1.67	0.59
2:B:170:VAL:HA	5:B:671:HOH:O	2.02	0.59
1:A:78:GLU:OE1	5:A:629:HOH:O	2.17	0.59
1:A:138:GLU:CD	1:A:141:GLN:HE21	2.10	0.58
2:B:166:THR:HG23	2:B:226:VAL:HG21	1.85	0.57
2:B:170:VAL:C	5:B:671:HOH:O	2.40	0.56
2:B:1:MET:HE2	2:B:155:HIS:HA	1.88	0.56
1:A:138:GLU:HG2	5:B:650:HOH:O	2.07	0.54
1:A:2:THR:OG1	5:A:692:HOH:O	2.19	0.53
1:A:144:LYS:NZ	1:A:150:GLU:OE1	2.28	0.53
2:B:146[B]:ARG:CB	2:B:226:VAL:HG23	2.24	0.52
2:B:146[B]:ARG:HB2	2:B:226:VAL:CG2	2.27	0.52
2:B:225:LEU:CB	2:B:227:ASP:HA	2.31	0.51
2:B:201:GLN:NE2	5:B:730:HOH:O	2.23	0.51
1:A:138:GLU:OE2	1:A:141:GLN:NE2	2.44	0.51
2:B:166:THR:OG1	2:B:226:VAL:HG21	2.12	0.50
1:A:138:GLU:CD	1:A:141:GLN:NE2	2.70	0.49
2:B:1:MET:SD	2:B:56:GLU:OE1	2.70	0.49
2:B:166:THR:HG22	2:B:167:GLY:N	2.27	0.48
2:B:160:ARG:NH1	5:B:533:HOH:O	2.02	0.47
2:B:1:MET:HG3	2:B:155:HIS:HA	1.96	0.47
2:B:166:THR:CA	5:B:707:HOH:O	2.61	0.47
2:B:145:VAL:O	2:B:166:THR:C	2.59	0.46
2:B:212:SER:OG	5:B:580:HOH:O	2.20	0.46
2:B:1:MET:HB2	2:B:154:GLY:O	2.16	0.45
2:B:172:HIS:O	2:B:193:LEU:O	2.34	0.45
2:B:170:VAL:O	2:B:195:THR:O	2.36	0.44
2:B:166:THR:CG2	2:B:226:VAL:HG21	2.47	0.43
1:A:156:GLU:HG3	1:A:157:ILE:N	2.33	0.43
2:B:146[B]:ARG:H	2:B:226:VAL:N	2.07	0.42
2:B:146[A]:ARG:HB2	2:B:226:VAL:CG2	2.25	0.42
2:B:146[A]:ARG:H	2:B:226:VAL:N	2.08	0.42
1:A:154:SER:O	2:B:211[A]:ASN:ND2	2.53	0.41
1:A:202:LYS:HG2	5:A:620:HOH:O	2.20	0.41
2:B:1:MET:HB3	5:B:356:HOH:O	2.20	0.41
2:B:145:VAL:C	2:B:226:VAL:HB	2.45	0.41
2:B:1:MET:CG	2:B:154:GLY:O	2.68	0.41
1:A:12:GLU:HG3	5:A:680:HOH:O	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:CYS:HB3	1:A:167:ARG:CZ	2.52	0.40

All (2) 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)
5:B:676:HOH:O	5:B:687:HOH:O[4_555]	1.60	0.60
5:B:668:HOH:O	5:B:668:HOH:O[4_555]	2.13	0.07

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	207/290 (71%)	203 (98%)	3 (1%)	1 (0%)	24	9
2	B	237/233 (102%)	225 (95%)	8 (3%)	4 (2%)	7	0
All	All	444/523 (85%)	428 (96%)	11 (2%)	5 (1%)	10	1

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	173	HIS
2	B	225	LEU
2	B	166	THR
2	B	171	LYS
1	A	112	SER

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	185/254 (73%)	185 (100%)	0	100	100
2	B	199/191 (104%)	194 (98%)	5 (2%)	42	14
All	All	384/445 (86%)	379 (99%)	5 (1%)	73	38

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	1	MET
2	B	13[A]	LEU
2	B	13[B]	LEU
2	B	46[A]	MET
2	B	46[B]	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue 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	$\# Z > 2$
1	CSD	A	111	1	4,7,8	1.83	1 (25%)	1,8,10	3.16	1 (100%)

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
1	CSD	A	111	1	-	1/2/6/8	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	111	CSD	OD1-SG	3.36	1.50	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	111	CSD	OD1-SG-CB	-3.16	99.78	105.60

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	111	CSD	CA-CB-SG-OD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	BUB	A	302	3	6,6,6	2.35	1 (16%)	4,6,6	0.85	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
4	BUB	A	302	3	-	1/2/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	302	BUB	B12-C9	5.71	1.66	1.56

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	302	BUB	C2-C6-C9-B12

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	202/290 (69%)	-0.71	2 (0%) 79 84	5, 11, 24, 41	7 (3%)
2	B	228/233 (97%)	-0.46	7 (3%) 51 56	6, 10, 24, 127	11 (4%)
All	All	430/523 (82%)	-0.58	9 (2%) 63 70	5, 11, 24, 127	18 (4%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	226	VAL	17.7
2	B	227	ASP	7.2
2	B	228	THR	6.2
2	B	170	VAL	6.1
1	A	204	VAL	5.3
2	B	225	LEU	4.6
2	B	1	MET	4.2
2	B	166	THR	3.7
1	A	2	THR	2.4

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CSD	A	111	8/9	0.99	0.04	5,8,10,11	2

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(\AA^2)	Q<0.9
4	BUB	A	302	7/7	0.94	0.09	22,25,28,29	0
3	CO	A	301	1/1	1.00	0.01	6,6,6,6	1

6.5 Other polymers [i](#)

There are no such residues in this entry.