



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 07:02 AM EDT

PDB ID : 2GVH / pdb_00002gvh
Title : Crystal structure of Acyl-CoA hydrolase (15159470) from AGROBACTERIUM TUMEFACIENS at 2.65 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2006-05-02
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

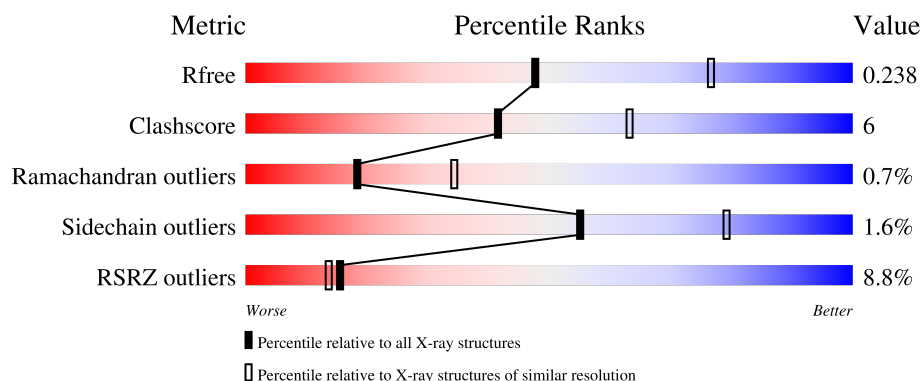
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	288	
1	B	288	
1	C	288	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5444 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 AGR_L_2016p.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	Se	0	0	0
			1785	1136	312	326	3	8			
1	B	250	Total	C	N	O	S	Se	0	0	0
			1828	1159	327	331	3	8			
1	C	250	Total	C	N	O	S	Se	0	0	0
			1830	1158	326	335	3	8			

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MSE	-	expression tag	GB 15159470
A	-10	GLY	-	expression tag	GB 15159470
A	-9	SER	-	expression tag	GB 15159470
A	-8	ASP	-	expression tag	GB 15159470
A	-7	LYS	-	expression tag	GB 15159470
A	-6	ILE	-	expression tag	GB 15159470
A	-5	HIS	-	expression tag	GB 15159470
A	-4	HIS	-	expression tag	GB 15159470
A	-3	HIS	-	expression tag	GB 15159470
A	-2	HIS	-	expression tag	GB 15159470
A	-1	HIS	-	expression tag	GB 15159470
A	0	HIS	-	expression tag	GB 15159470
A	1	MSE	MET	modified residue	GB 15159470
A	43	MSE	MET	modified residue	GB 15159470
A	99	MSE	MET	modified residue	GB 15159470
A	119	MSE	MET	modified residue	GB 15159470
A	150	MSE	MET	modified residue	GB 15159470
A	165	MSE	MET	modified residue	GB 15159470
A	174	MSE	MET	modified residue	GB 15159470
A	224	MSE	MET	modified residue	GB 15159470
A	250	MSE	MET	modified residue	GB 15159470
B	-11	MSE	-	expression tag	GB 15159470
B	-10	GLY	-	expression tag	GB 15159470

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-9	SER	-	expression tag	GB 15159470
B	-8	ASP	-	expression tag	GB 15159470
B	-7	LYS	-	expression tag	GB 15159470
B	-6	ILE	-	expression tag	GB 15159470
B	-5	HIS	-	expression tag	GB 15159470
B	-4	HIS	-	expression tag	GB 15159470
B	-3	HIS	-	expression tag	GB 15159470
B	-2	HIS	-	expression tag	GB 15159470
B	-1	HIS	-	expression tag	GB 15159470
B	0	HIS	-	expression tag	GB 15159470
B	1	MSE	MET	modified residue	GB 15159470
B	43	MSE	MET	modified residue	GB 15159470
B	99	MSE	MET	modified residue	GB 15159470
B	119	MSE	MET	modified residue	GB 15159470
B	150	MSE	MET	modified residue	GB 15159470
B	165	MSE	MET	modified residue	GB 15159470
B	174	MSE	MET	modified residue	GB 15159470
B	224	MSE	MET	modified residue	GB 15159470
B	250	MSE	MET	modified residue	GB 15159470
C	-11	MSE	-	expression tag	GB 15159470
C	-10	GLY	-	expression tag	GB 15159470
C	-9	SER	-	expression tag	GB 15159470
C	-8	ASP	-	expression tag	GB 15159470
C	-7	LYS	-	expression tag	GB 15159470
C	-6	ILE	-	expression tag	GB 15159470
C	-5	HIS	-	expression tag	GB 15159470
C	-4	HIS	-	expression tag	GB 15159470
C	-3	HIS	-	expression tag	GB 15159470
C	-2	HIS	-	expression tag	GB 15159470
C	-1	HIS	-	expression tag	GB 15159470
C	0	HIS	-	expression tag	GB 15159470
C	1	MSE	MET	modified residue	GB 15159470
C	43	MSE	MET	modified residue	GB 15159470
C	99	MSE	MET	modified residue	GB 15159470
C	119	MSE	MET	modified residue	GB 15159470
C	150	MSE	MET	modified residue	GB 15159470
C	165	MSE	MET	modified residue	GB 15159470
C	174	MSE	MET	modified residue	GB 15159470
C	224	MSE	MET	modified residue	GB 15159470
C	250	MSE	MET	modified residue	GB 15159470

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		

- Molecule 1: AGR L 2016p



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	121.21Å 121.21Å 81.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.71 – 2.50 48.71 – 2.50	Depositor EDS
% Data completeness (in resolution range)	86.1 (48.71-2.50) 86.1 (48.71-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.226 , 0.244 0.221 , 0.238	Depositor DCC
R_{free} test set	1773 reflections (4.32%)	wwPDB-VP
Wilson B-factor (Å ²)	58.9	Xtriage
Anisotropy	0.725	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5444	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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: NA

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	0.79	1/1810 (0.1%)	0.90	2/2448 (0.1%)
1	B	1.01	2/1856 (0.1%)	1.02	3/2508 (0.1%)
1	C	0.92	0/1855	1.03	6/2502 (0.2%)
All	All	0.91	3/5521 (0.1%)	0.99	11/7458 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	150	MSE	SE-CE	-10.88	1.62	1.95
1	A	150	MSE	SE-CE	-7.03	1.74	1.95
1	B	250	MSE	SE-CE	-6.84	1.75	1.95

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	258	ARG	CA-C-N	8.91	128.99	119.90
1	C	258	ARG	C-N-CA	8.91	128.99	119.90
1	C	104	ILE	N-CA-C	-6.42	104.50	110.53
1	C	9	LYS	CA-C-N	-6.23	113.53	120.14
1	C	9	LYS	C-N-CA	-6.23	113.53	120.14
1	B	258	ARG	CA-C-N	6.19	126.21	119.89
1	B	258	ARG	C-N-CA	6.19	126.21	119.89
1	A	258	ARG	CA-C-N	6.14	127.51	119.84
1	A	258	ARG	C-N-CA	6.14	127.51	119.84
1	B	146	ASP	N-CA-C	-5.99	105.94	113.18
1	C	148	VAL	N-CA-C	5.41	117.09	108.87

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	1785	0	1701	23	0
1	B	1828	0	1765	27	0
1	C	1830	0	1760	21	0
2	B	1	0	0	0	0
All	All	5444	0	5226	68	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 6.

All (68) 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:B:213:GLN:HG3	1:B:231:TRP:HZ3	1.36	0.91
1:C:190:VAL:HG23	1:C:250:MSE:CE	2.10	0.82
1:A:20:ILE:HD12	1:A:150:MSE:CE	2.14	0.76
1:B:150:MSE:HG3	1:B:173:TYR:OH	1.86	0.76
1:B:213:GLN:HG3	1:B:231:TRP:CZ3	2.21	0.75
1:B:213:GLN:CG	1:B:231:TRP:HZ3	2.00	0.73
1:A:190:VAL:HG23	1:A:250:MSE:CE	2.21	0.70
1:B:190:VAL:HG23	1:B:250:MSE:CE	2.22	0.69
1:A:20:ILE:HB	1:A:150:MSE:HE2	1.78	0.65
1:C:190:VAL:HG23	1:C:250:MSE:HE3	1.78	0.63
1:A:20:ILE:HD12	1:A:150:MSE:HE2	1.78	0.63
1:A:148:VAL:HG11	1:A:150:MSE:HE3	1.87	0.56
1:B:190:VAL:HG23	1:B:250:MSE:HE2	1.86	0.56
1:A:20:ILE:HB	1:A:150:MSE:CE	2.36	0.56
1:A:158:GLN:HB3	1:A:169:GLU:HG2	1.89	0.54
1:C:190:VAL:CG2	1:C:250:MSE:CE	2.83	0.54
1:B:222:SER:HB2	1:B:260:ALA:O	2.06	0.53
1:A:190:VAL:HG23	1:A:250:MSE:HE2	1.91	0.53
1:B:213:GLN:CD	1:B:231:TRP:CZ3	2.88	0.52
1:B:204:ILE:HD11	1:B:243:THR:HG21	1.92	0.51
1:A:20:ILE:CD1	1:A:150:MSE:CE	2.89	0.51
1:C:190:VAL:HG23	1:C:250:MSE:HE2	1.90	0.51
1:C:190:VAL:CG2	1:C:250:MSE:HE2	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:158:GLN:HB3	1:C:169:GLU:HG2	1.95	0.49
1:B:148:VAL:HG11	1:B:150:MSE:CE	2.42	0.49
1:B:190:VAL:CG2	1:B:250:MSE:HE2	2.43	0.48
1:B:179:PHE:C	1:B:179:PHE:CD1	2.90	0.48
1:C:19:LEU:C	1:C:19:LEU:HD23	2.38	0.48
1:A:204:ILE:HD11	1:A:243:THR:HG21	1.97	0.47
1:B:213:GLN:CG	1:B:231:TRP:CZ3	2.88	0.46
1:B:70:ARG:HG3	1:B:111:THR:HG22	1.97	0.45
1:A:179:PHE:C	1:A:179:PHE:CD1	2.94	0.45
1:A:148:VAL:CG1	1:A:150:MSE:HE3	2.45	0.45
1:A:190:VAL:HG23	1:A:250:MSE:HE3	1.98	0.45
1:A:40:LEU:HD23	1:A:40:LEU:HA	1.75	0.45
1:C:204:ILE:HD11	1:C:243:THR:HG21	1.98	0.45
1:C:179:PHE:HA	1:C:250:MSE:HE2	1.97	0.45
1:A:148:VAL:O	1:A:213:GLN:HA	2.18	0.44
1:B:72:PRO:HB3	1:B:110:HIS:HB3	2.00	0.44
1:B:148:VAL:O	1:B:213:GLN:HA	2.17	0.44
1:A:27:ASP:OD2	1:A:30:HIS:HB2	2.17	0.43
1:C:131:TYR:C	1:C:131:TYR:CD1	2.96	0.43
1:A:190:VAL:CG2	1:A:250:MSE:HE2	2.48	0.43
1:C:213:GLN:HG3	1:C:231:TRP:HZ3	1.82	0.43
1:B:132:VAL:HG23	1:B:133:LEU:O	2.17	0.43
1:A:44:ASP:HB3	1:C:168:GLY:HA3	2.01	0.42
1:C:179:PHE:CD1	1:C:179:PHE:C	2.97	0.42
1:B:54:PHE:CE2	1:B:134:PRO:HG2	2.55	0.42
1:C:18:ARG:HG2	1:C:82:THR:HG22	2.01	0.42
1:B:190:VAL:CG2	1:B:250:MSE:CE	2.96	0.42
1:B:213:GLN:O	1:B:228:THR:HA	2.20	0.42
1:B:168:GLY:HA3	1:C:44:ASP:HB3	2.02	0.42
1:A:29:ASN:CB	1:C:27:ASP:OD2	2.68	0.41
1:B:42:LEU:O	1:B:46:VAL:HG12	2.20	0.41
1:C:235:LEU:HD23	1:C:235:LEU:HA	1.92	0.41
1:A:141:THR:HA	1:A:142:PRO:HD3	1.79	0.41
1:B:158:GLN:HB3	1:B:169:GLU:HG2	2.03	0.41
1:C:190:VAL:CG2	1:C:250:MSE:HE3	2.47	0.41
1:A:150:MSE:HE2	1:A:150:MSE:HB3	1.90	0.41
1:A:190:VAL:CG2	1:A:250:MSE:CE	2.96	0.41
1:B:56:ARG:HB3	1:B:236:LEU:O	2.21	0.41
1:C:40:LEU:HD23	1:C:40:LEU:HA	1.92	0.41
1:C:150:MSE:HG3	1:C:173:TYR:OH	2.21	0.41
1:C:148:VAL:CG1	1:C:150:MSE:HE2	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:ALA:C	1:B:130:SER:H	2.29	0.40
1:A:68:ASP:O	1:A:113:THR:HG23	2.21	0.40
1:B:46:VAL:HG11	1:B:81:PHE:HB3	2.03	0.40
1:B:226:ILE:HD13	1:B:250:MSE:CG	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	242/288 (84%)	236 (98%)	5 (2%)	1 (0%)	30	49
1	B	246/288 (85%)	237 (96%)	7 (3%)	2 (1%)	16	31
1	C	244/288 (85%)	239 (98%)	3 (1%)	2 (1%)	16	31
All	All	732/864 (85%)	712 (97%)	15 (2%)	5 (1%)	18	34

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	ILE
1	C	75	ILE
1	B	74	ARG
1	B	75	ILE
1	C	74	ARG

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	166/220 (76%)	164 (99%)	2 (1%)	63	83
1	B	174/220 (79%)	171 (98%)	3 (2%)	53	78
1	C	175/220 (80%)	172 (98%)	3 (2%)	53	78
All	All	515/660 (78%)	507 (98%)	8 (2%)	55	79

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	132	VAL
1	A	149	THR
1	B	107	ARG
1	B	132	VAL
1	B	133	LEU
1	C	29	ASN
1	C	71	GLN
1	C	110	HIS

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	215	HIS
1	A	241	HIS
1	B	12	GLN
1	B	29	ASN
1	C	12	GLN
1	C	13	HIS
1	C	30	HIS
1	C	31	HIS
1	C	110	HIS
1	C	215	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 [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	240/288 (83%)	0.71	17 (7%) 22 19	51, 68, 81, 89	0
1	B	242/288 (84%)	0.75	28 (11%) 9 8	44, 69, 84, 101	0
1	C	242/288 (84%)	0.70	19 (7%) 18 16	49, 69, 83, 101	0
All	All	724/864 (83%)	0.72	64 (8%) 15 14	44, 69, 83, 101	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	143	ASP	7.8
1	C	143	ASP	6.8
1	B	254	ASP	6.2
1	C	254	ASP	5.8
1	A	254	ASP	5.5
1	C	141	THR	4.9
1	B	75	ILE	4.9
1	A	262	ILE	4.6
1	A	110	HIS	4.5
1	A	108	GLN	4.0
1	C	7	ILE	3.8
1	A	69	PHE	3.6
1	B	262	ILE	3.6
1	A	75	ILE	3.6
1	C	258	ARG	3.5
1	B	146	ASP	3.4
1	A	127	ASP	3.3
1	B	143	ASP	3.3
1	C	75	ILE	3.2
1	A	258	ARG	3.1
1	A	29	ASN	2.8
1	A	222	SER	2.7
1	B	92	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	153	ILE	2.6
1	B	140	GLU	2.6
1	C	206	ILE	2.6
1	A	109	GLN	2.6
1	C	73	ALA	2.6
1	B	258	ARG	2.5
1	C	69	PHE	2.5
1	B	97	VAL	2.5
1	B	74	ARG	2.5
1	B	207	GLY	2.5
1	A	206	ILE	2.5
1	B	251	VAL	2.5
1	C	147	ALA	2.5
1	C	142	PRO	2.4
1	B	69	PHE	2.4
1	B	71	GLN	2.4
1	B	145	SER	2.4
1	B	98	GLU	2.4
1	A	20	ILE	2.4
1	C	214	ALA	2.4
1	B	141	THR	2.3
1	B	116	ILE	2.3
1	A	138	THR	2.3
1	C	148	VAL	2.3
1	B	126	GLU	2.2
1	A	13	HIS	2.2
1	B	20	ILE	2.2
1	B	113	THR	2.2
1	B	10	PRO	2.1
1	B	191	VAL	2.1
1	B	82	THR	2.1
1	C	28	THR	2.1
1	C	140	GLU	2.1
1	C	156	PRO	2.1
1	B	124	GLU	2.1
1	B	127	ASP	2.1
1	A	112	CYS	2.1
1	C	126	GLU	2.0
1	C	161	SER	2.0
1	B	147	ALA	2.0
1	C	139	GLU	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
2	NA	B	277	1/1	0.91	0.09	81,81,81,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.