



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 01:53 pm BST

PDB ID : 5FJU  
Title : N-acyl amino acid racemase from *Amycolatopsis* sp. Ts-1-60: Q26A M50I  
G291D F323Y mutant in complex with N-acetyl phenylalanine  
Authors : Sanchez Carron, G.; Campopiano, D.; Grogan, G.  
Deposited on : 2015-10-13  
Resolution : 2.52 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

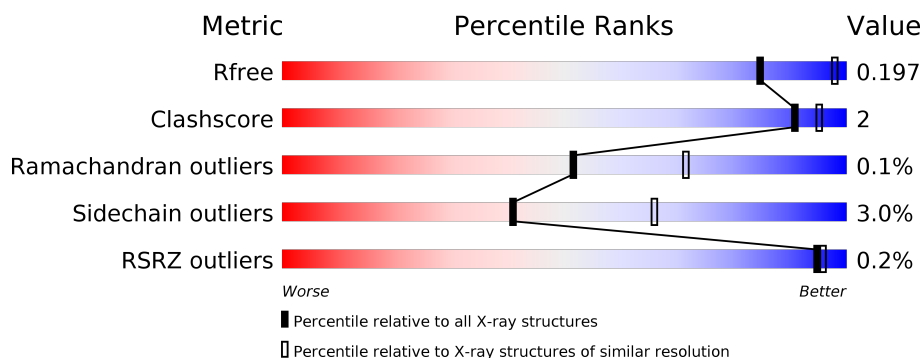
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.52 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))
$R_{free}$	130704	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (2.54-2.50)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	368	<div> <div style="width: 93%;"></div> <div style="width: 6%;"></div> <div style="width: 1%;"></div> </div> <div>93% 6% .</div>
1	B	368	<div> <div style="width: 94%;"></div> <div style="width: 6%;"></div> <div style="width: 0%;"></div> </div> <div>94% 6%</div>
1	C	368	<div> <div style="width: 94%;"></div> <div style="width: 5%;"></div> <div style="width: 1%;"></div> </div> <div>94% 5% .</div>
1	D	368	<div> <div style="width: 93%;"></div> <div style="width: 6%;"></div> <div style="width: 1%;"></div> </div> <div>93% 6% .</div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11561 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 O-SUCCINYLBENZOATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2755	1753	477	514	11			
1	B	367	Total	C	N	O	S	0	0	0
			2736	1744	478	503	11			
1	C	367	Total	C	N	O	S	0	0	0
			2734	1743	477	503	11			
1	D	367	Total	C	N	O	S	0	0	0
			2743	1746	477	509	11			

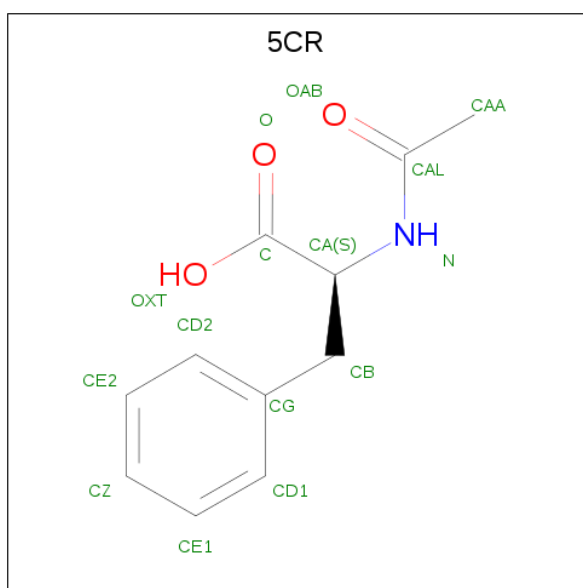
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	ALA	GLN	engineered mutation	UNP Q44244
A	50	ILE	MET	engineered mutation	UNP Q44244
A	291	ASP	GLY	engineered mutation	UNP Q44244
A	323	TYR	PHE	engineered mutation	UNP Q44244
B	26	ALA	GLN	engineered mutation	UNP Q44244
B	50	ILE	MET	engineered mutation	UNP Q44244
B	291	ASP	GLY	engineered mutation	UNP Q44244
B	323	TYR	PHE	engineered mutation	UNP Q44244
C	26	ALA	GLN	engineered mutation	UNP Q44244
C	50	ILE	MET	engineered mutation	UNP Q44244
C	291	ASP	GLY	engineered mutation	UNP Q44244
C	323	TYR	PHE	engineered mutation	UNP Q44244
D	26	ALA	GLN	engineered mutation	UNP Q44244
D	50	ILE	MET	engineered mutation	UNP Q44244
D	291	ASP	GLY	engineered mutation	UNP Q44244
D	323	TYR	PHE	engineered mutation	UNP Q44244

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

- Molecule 3 is N-acetyl-L-phenylalanine (three-letter code: 5CR) (formula:  $C_{11}H_{13}NO_3$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 15 11 1 3	0	0
3	A	1	Total C N O 15 11 1 3	0	0
3	B	1	Total C N O 15 11 1 3	0	0
3	B	1	Total C N O 15 11 1 3	0	0
3	C	1	Total C N O 15 11 1 3	0	0
3	C	1	Total C N O 15 11 1 3	0	0
3	D	1	Total C N O 15 11 1 3	0	0
3	D	1	Total C N O 15 11 1 3	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	166	Total 166	O 166	0	0
4	B	111	Total 111	O 111	0	0
4	C	88	Total 88	O 88	0	0
4	D	104	Total 104	O 104	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: O-SUCCINYLBENZOATE SYNTHASE

Chain A:  93% 6%

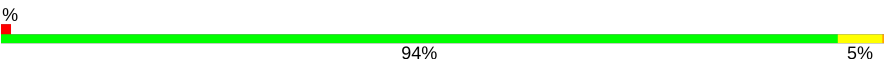


#### • Molecule 1: O-SUCCINYLBENZOATE SYNTHASE

Chain B:  94% 6%



#### • Molecule 1: O-SUCCINYLBENZOATE SYNTHASE

Chain C:  94% 5%



#### • Molecule 1: O-SUCCINYLBENZOATE SYNTHASE

Chain D:  93% 6%



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	216.90Å 216.90Å 261.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	107.35 – 2.52 107.35 – 2.52	Depositor EDS
% Data completeness (in resolution range)	100.0 (107.35-2.52) 100.0 (107.35-2.52)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.07 (at 2.52Å)	Xtriage
Refinement program	REFMAC 5.8.0123	Depositor
R, $R_{free}$	0.156 , 0.190 0.164 , 0.197	Depositor DCC
$R_{free}$ test set	3783 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	47.4	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 26.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -2/3*h-1/3*k+2/3*l,-1/3*h-2/3*k-2/3*l,2/3*h-2/3*k+1/3*l 0.018 for -h,1/3*h-1/3*k+2/3*l,2/3*h+4/3*k+1/3*l 0.013 for -1/3*h+1/3*k-2/3*l,-k,-4/3*h-2/3*k+1/3*l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	11561	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.13% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: MG, 5CR

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	0.80	1/2811 (0.0%)	0.94	8/3835 (0.2%)
1	B	0.79	1/2792 (0.0%)	0.92	8/3809 (0.2%)
1	C	0.78	1/2790 (0.0%)	0.93	9/3807 (0.2%)
1	D	0.81	0/2799	0.92	8/3819 (0.2%)
All	All	0.79	3/11192 (0.0%)	0.93	33/15270 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	218	GLU	CD-OE2	6.48	1.32	1.25
1	A	7	GLU	CD-OE1	6.05	1.32	1.25
1	B	250	ASP	CB-CG	5.25	1.62	1.51

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	291	ASP	CB-CG-OD1	8.41	125.87	118.30
1	C	113	ARG	NE-CZ-NH1	8.28	124.44	120.30
1	D	113	ARG	NE-CZ-NH1	8.23	124.42	120.30
1	C	204	ARG	NE-CZ-NH2	-8.22	116.19	120.30
1	D	113	ARG	NE-CZ-NH2	-8.07	116.27	120.30

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	2755	0	2774	10	0
1	B	2736	0	2770	6	0
1	C	2734	0	2763	9	0
1	D	2743	0	2769	12	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	30	0	0	0	0
3	B	30	0	0	0	0
3	C	30	0	0	0	0
3	D	30	0	0	0	0
4	A	166	0	0	0	0
4	B	111	0	0	1	0
4	C	88	0	0	1	0
4	D	104	0	0	1	0
All	All	11561	0	11076	37	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 2.

The worst 5 of 37 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:C:7:GLU:OE1	1:C:35:ARG:NH1	2.12	0.81
1:D:214:GLU:OE2	4:D:2063:HOH:O	2.00	0.78
1:D:10:ARG:NH2	1:D:61:ASP:OD1	2.16	0.78
1:D:1:MET:O	1:D:80:ASP:O	2.12	0.67
1:A:11:VAL:HG13	1:A:359:VAL:CG1	2.26	0.66

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	366/368 (100%)	358 (98%)	8 (2%)	0	100	100
1	B	365/368 (99%)	357 (98%)	8 (2%)	0	100	100
1	C	365/368 (99%)	358 (98%)	7 (2%)	0	100	100
1	D	365/368 (99%)	355 (97%)	9 (2%)	1 (0%)	41	59
All	All	1461/1472 (99%)	1428 (98%)	32 (2%)	1 (0%)	51	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	81	ILE

### 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	285/291 (98%)	276 (97%)	9 (3%)	39	63
1	B	282/291 (97%)	275 (98%)	7 (2%)	47	72
1	C	281/291 (97%)	272 (97%)	9 (3%)	39	63
1	D	284/291 (98%)	275 (97%)	9 (3%)	39	63
All	All	1132/1164 (97%)	1098 (97%)	34 (3%)	41	66

5 of 34 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	339	LEU
1	C	116	GLU
1	D	179	ARG
1	C	21	THR
1	A	322	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no

such sidechains identified.

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
3	5CR	C	1370	2	12,15,15	1.81	4 (33%)	14,19,19	1.71	3 (21%)
3	5CR	B	1369	2	12,15,15	1.69	4 (33%)	14,19,19	1.71	3 (21%)
3	5CR	A	1370	2	12,15,15	1.63	3 (25%)	14,19,19	1.31	3 (21%)
3	5CR	A	1371	-	12,15,15	2.06	4 (33%)	14,19,19	3.45	6 (42%)
3	5CR	B	1370	-	12,15,15	2.49	5 (41%)	14,19,19	3.28	7 (50%)
3	5CR	C	1369	-	12,15,15	1.77	3 (25%)	14,19,19	1.85	3 (21%)
3	5CR	D	1370	-	12,15,15	1.58	2 (16%)	14,19,19	1.55	1 (7%)
3	5CR	D	1369	-	12,15,15	2.75	3 (25%)	14,19,19	3.44	8 (57%)

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	5CR	C	1370	2	-	2/8/12/12	0/1/1/1
3	5CR	B	1369	2	-	2/8/12/12	0/1/1/1
3	5CR	A	1370	2	-	3/8/12/12	0/1/1/1
3	5CR	A	1371	-	-	1/8/12/12	0/1/1/1
3	5CR	B	1370	-	-	2/8/12/12	0/1/1/1
3	5CR	C	1369	-	-	1/8/12/12	0/1/1/1
3	5CR	D	1370	-	-	3/8/12/12	0/1/1/1
3	5CR	D	1369	-	-	2/8/12/12	0/1/1/1

The worst 5 of 28 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1369	5CR	CA-N	6.17	1.54	1.46
3	D	1369	5CR	CB-CA	5.94	1.62	1.53
3	B	1370	5CR	CB-CA	-5.21	1.46	1.53
3	B	1370	5CR	CB-CG	-4.24	1.41	1.51
3	A	1371	5CR	CB-CA	-4.19	1.48	1.53

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1371	5CR	CA-N-CAL	9.65	133.04	122.44
3	D	1369	5CR	CA-N-CAL	9.64	133.03	122.44
3	B	1370	5CR	CA-N-CAL	7.94	131.16	122.44
3	B	1370	5CR	CG-CB-CA	5.61	122.52	112.97
3	D	1370	5CR	CG-CB-CA	-5.19	104.14	112.97

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1370	5CR	C-CA-CB-CG
3	A	1371	5CR	CB-CA-N-CAL
3	C	1369	5CR	CB-CA-N-CAL
3	D	1370	5CR	CB-CA-N-CAL
3	D	1370	5CR	C-CA-CB-CG

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	368/368 (100%)	-0.32	1 (0%) 94 94	30, 41, 60, 88	0
1	B	367/368 (99%)	-0.29	0 100 100	32, 45, 67, 92	0
1	C	367/368 (99%)	-0.23	2 (0%) 91 92	31, 50, 70, 101	0
1	D	367/368 (99%)	-0.32	0 100 100	30, 45, 63, 92	0
All	All	1469/1472 (99%)	-0.29	3 (0%) 95 96	30, 45, 67, 101	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	40	ALA	2.4
1	C	1	MET	2.3
1	A	1	MET	2.3

### 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 carbohydrates 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, 95<sup>th</sup> 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	MG	B	1368	1/1	0.83	0.16	56,56,56,56	0
2	MG	C	1368	1/1	0.89	0.10	51,51,51,51	0
3	5CR	D	1369	15/15	0.91	0.20	44,60,82,83	0
2	MG	A	1369	1/1	0.92	0.16	50,50,50,50	0
3	5CR	C	1369	15/15	0.93	0.28	49,57,71,76	0
3	5CR	A	1371	15/15	0.95	0.30	48,53,69,75	0
3	5CR	B	1370	15/15	0.95	0.18	40,54,70,95	0
3	5CR	A	1370	15/15	0.96	0.16	40,44,51,56	0
3	5CR	B	1369	15/15	0.96	0.15	44,49,57,59	0
3	5CR	C	1370	15/15	0.97	0.16	50,53,63,69	0
2	MG	D	1368	1/1	0.98	0.10	51,51,51,51	0
3	5CR	D	1370	15/15	0.98	0.18	44,49,57,57	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.