



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:17 PM UTC

PDB ID : 3PGY / pdb_00003pgy
Title : Serine hydroxymethyltransferase from Staphylococcus aureus, S95P mutant.
Authors : Osipiuk, J.; Makowska-Grzyska, M.; Kwon, K.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2010-11-02
Resolution : 1.92 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

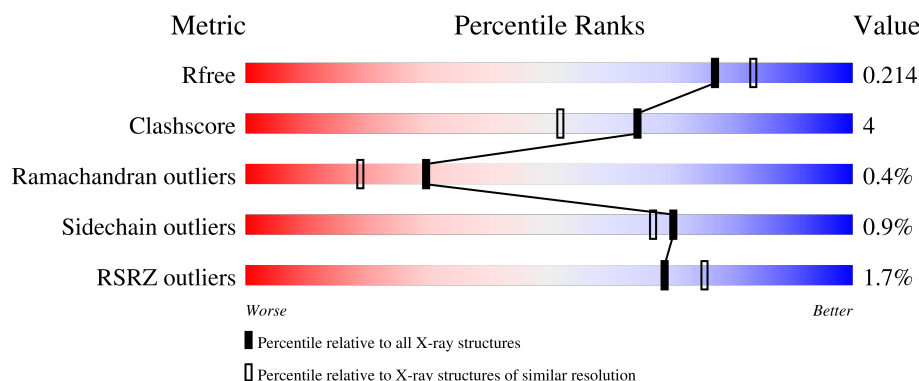
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	415	0% 88% 8% .
1	B	415	2% 90% 7% .
1	C	415	0% 89% 7% .
1	D	415	2% 87% 9% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CIT	B	501	-	X	-	-
2	CIT	B	506	-	X	-	-
2	CIT	D	503	-	X	-	-
3	GLY	B	509	-	X	-	-
3	GLY	B	511	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13941 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.

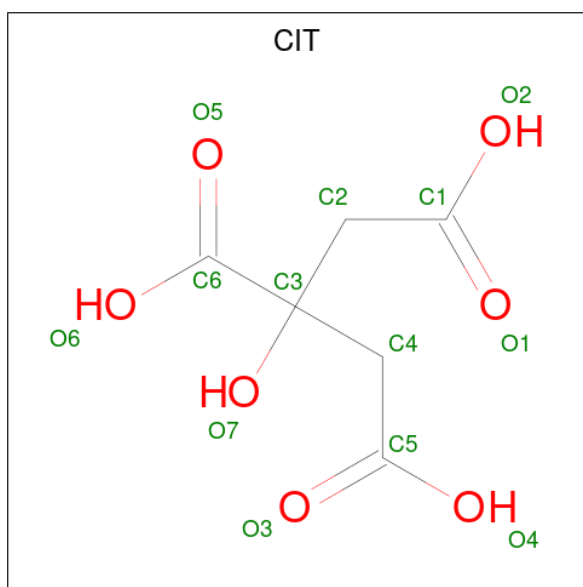
- Molecule 1 is a protein called Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	11	0
			3172	2017	532	612	11			
1	B	404	Total	C	N	O	S	0	8	0
			3172	2016	532	613	11			
1	C	398	Total	C	N	O	S	0	22	0
			3201	2044	535	611	11			
1	D	402	Total	C	N	O	S	0	17	0
			3212	2046	537	618	11			

There are 16 discrepancies between the modelled and reference sequences:

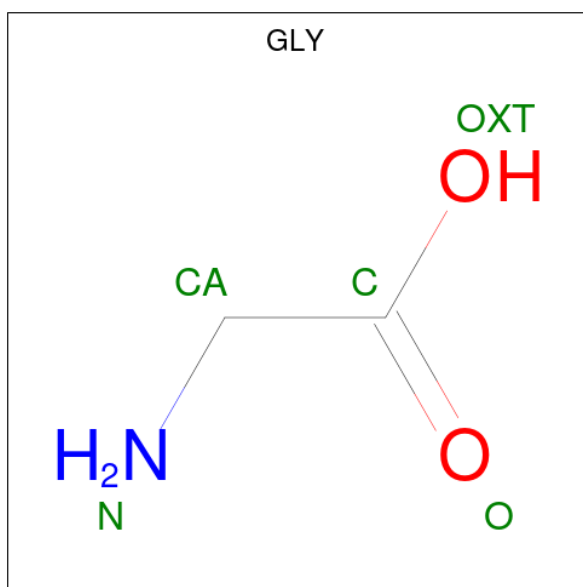
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q5HE87
A	-1	ASN	-	expression tag	UNP Q5HE87
A	0	ALA	-	expression tag	UNP Q5HE87
A	95	PRO	SER	engineered mutation	UNP Q5HE87
B	-2	SER	-	expression tag	UNP Q5HE87
B	-1	ASN	-	expression tag	UNP Q5HE87
B	0	ALA	-	expression tag	UNP Q5HE87
B	95	PRO	SER	engineered mutation	UNP Q5HE87
C	-2	SER	-	expression tag	UNP Q5HE87
C	-1	ASN	-	expression tag	UNP Q5HE87
C	0	ALA	-	expression tag	UNP Q5HE87
C	95	PRO	SER	engineered mutation	UNP Q5HE87
D	-2	SER	-	expression tag	UNP Q5HE87
D	-1	ASN	-	expression tag	UNP Q5HE87
D	0	ALA	-	expression tag	UNP Q5HE87
D	95	PRO	SER	engineered mutation	UNP Q5HE87

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



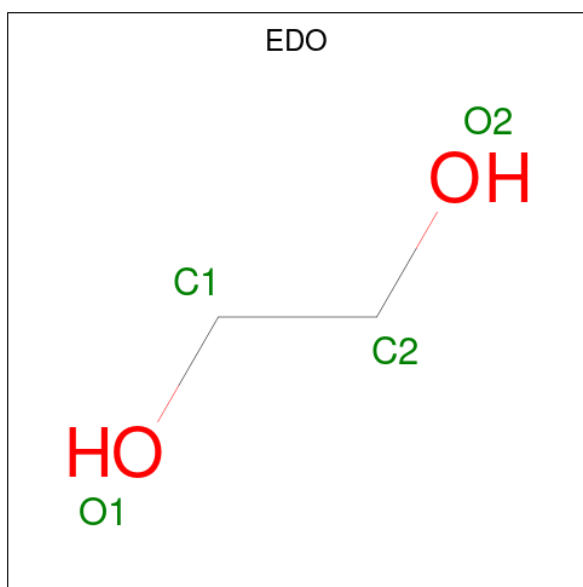
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	C	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is GLYCINE (CCD ID: GLY) (formula: $\text{C}_2\text{H}_5\text{NO}_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	B	1	Total	C	N	O	0	0
			5	2	1	2		
3	B	1	Total	C	N	O	0	0
			5	2	1	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

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

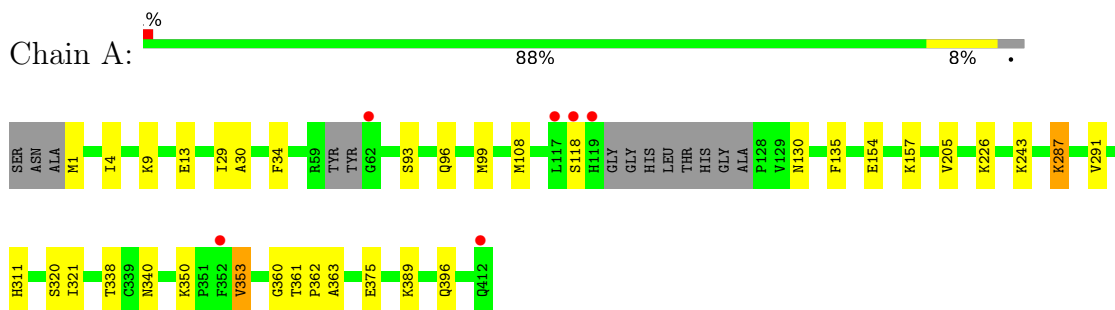
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	247	Total	O	0	0
			247	247		
5	B	289	Total	O	0	0
			289	289		
5	C	279	Total	O	0	0
			279	279		
5	D	242	Total	O	0	0
			242	242		

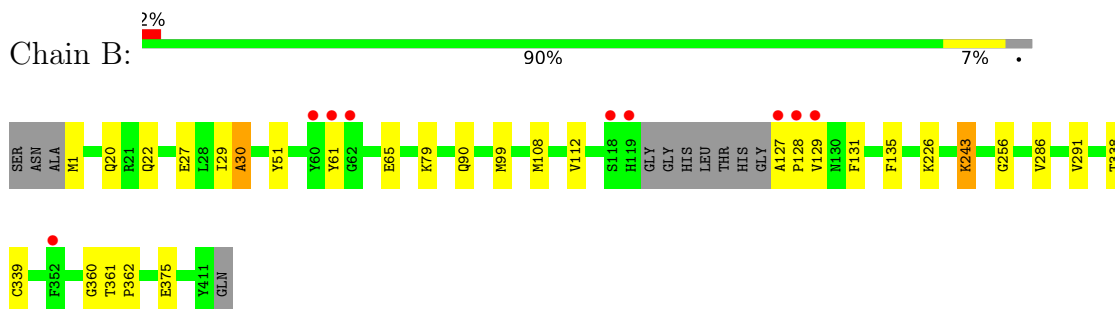
3 Residue-property plots [i](#)

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.

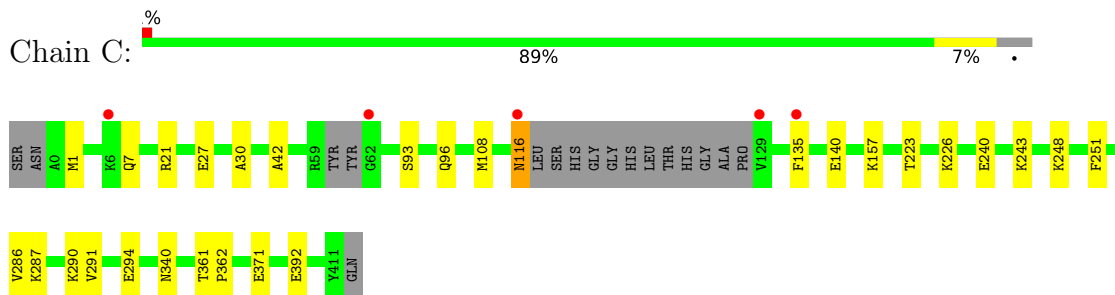
- Molecule 1: Serine hydroxymethyltransferase



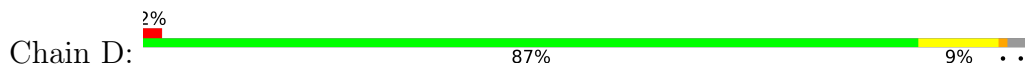
- Molecule 1: Serine hydroxymethyltransferase

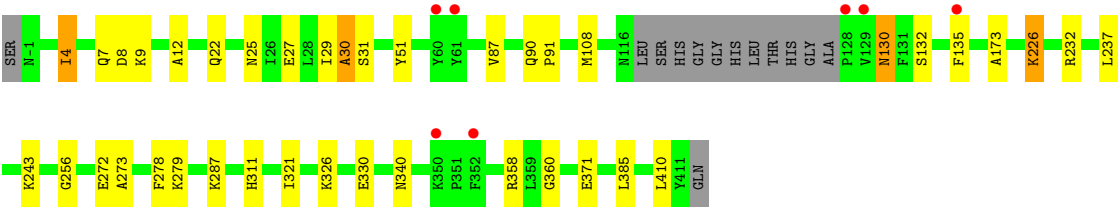


- Molecule 1: Serine hydroxymethyltransferase



- Molecule 1: Serine hydroxymethyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.08Å 86.53Å 301.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.70 – 1.92 47.70 – 1.92	Depositor EDS
% Data completeness (in resolution range)	97.8 (47.70-1.92) 97.7 (47.70-1.92)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.166 , 0.208 0.173 , 0.214	Depositor DCC
R_{free} test set	6819 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.592	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13941	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.59% 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: CIT, EDO

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	0.99	1/3257 (0.0%)	0.92	0/4388
1	B	0.99	0/3251	0.94	2/4386 (0.0%)
1	C	1.00	1/3318 (0.0%)	0.93	0/4466
1	D	0.95	0/3318	0.91	1/4474 (0.0%)
All	All	0.98	2/13144 (0.0%)	0.92	3/17714 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	205	VAL	C-O	5.97	1.30	1.24
1	C	223	THR	N-CA	-5.39	1.39	1.45

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	4	ILE	CB-CA-C	-7.30	102.28	112.14
1	B	90	GLN	CA-C-N	-5.92	113.44	119.83
1	B	90	GLN	C-N-CA	-5.92	113.44	119.83

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	3172	0	3207	21	0
1	B	3172	0	3186	24	0
1	C	3201	0	3281	24	1
1	D	3212	0	3253	33	0
2	A	13	0	5	1	0
2	B	39	0	15	8	0
2	C	13	0	5	1	0
2	D	39	0	15	5	0
3	A	5	0	2	0	0
3	B	10	0	4	0	0
4	C	8	0	12	0	0
5	A	247	0	0	1	1
5	B	289	0	0	8	0
5	C	279	0	0	5	0
5	D	242	0	0	5	0
All	All	13941	0	12985	100	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 4.

All (100) 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:506:CIT:O2	2:B:506:CIT:H41	1.51	0.96
1:C:290:LYS:HE3	5:C:819:HOH:O	1.82	0.78
2:D:507:CIT:O5	5:D:800:HOH:O	2.06	0.74
1:D:4:ILE:HG22	1:D:12:ALA:HB2	1.69	0.72
1:B:286:VAL:HG13	5:B:481:HOH:O	1.88	0.72
1:D:25[A]:ASN:ND2	5:D:774:HOH:O	2.23	0.71
1:D:173:ALA:HB3	1:D:358[B]:ARG:CZ	2.22	0.69
1:C:21:ARG:NH1	1:C:27:GLU:OE2	2.26	0.69
1:B:22:GLN:O	1:B:27[B]:GLU:OE2	2.11	0.68
1:C:240:GLU:OE1	1:C:243[A]:LYS:HD2	1.95	0.67
1:B:29:ILE:HD11	1:B:338:THR:HG22	1.77	0.66
1:D:226:LYS:NZ	2:D:507:CIT:O3	2.31	0.63
1:C:286:VAL:HG13	5:C:460:HOH:O	1.99	0.63
1:D:29:ILE:HG23	1:D:340:ASN:HD21	1.64	0.62
2:B:501:CIT:O1	5:B:424:HOH:O	2.16	0.61
1:A:9:LYS:HE2	1:A:13[A]:GLU:OE2	2.00	0.61
1:C:108:MET:HE3	1:D:108:MET:HE1	1.81	0.61
1:C:243[A]:LYS:NZ	2:C:508:CIT:O7	2.33	0.61
1:B:108:MET:HE1	1:B:135:PHE:CD1	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154[B]:GLU:OE2	1:A:157[B]:LYS:HD2	2.02	0.60
1:B:129:VAL:HG23	5:B:728:HOH:O	2.01	0.60
1:A:99:MET:HE3	5:B:510:HOH:O	2.00	0.60
1:A:1:MET:HE3	1:A:4:ILE:HG21	1.86	0.58
1:A:118:SER:HB3	1:A:130:ASN:ND2	2.18	0.58
1:B:256:GLY:HA3	2:B:506:CIT:H42	1.84	0.58
1:D:108:MET:HE1	1:D:135:PHE:CD2	2.41	0.56
1:A:243:LYS:HE2	2:A:505:CIT:O7	2.07	0.55
1:B:20:GLN:NE2	5:B:644:HOH:O	2.40	0.54
1:A:108:MET:SD	1:A:135:PHE:HB3	2.48	0.54
1:C:108:MET:HE3	1:D:108:MET:CE	2.39	0.53
1:D:31:SER:HB3	1:D:226:LYS:HB3	1.90	0.53
1:D:287[B]:LYS:HD2	1:D:371:GLU:OE2	2.09	0.52
1:D:51:TYR:HE1	2:D:503:CIT:H41	1.74	0.52
1:D:130:ASN:C	1:D:130:ASN:OD1	2.53	0.52
1:C:287[B]:LYS:HD2	5:C:818:HOH:O	2.09	0.51
1:D:22:GLN:O	1:D:27[A]:GLU:OE2	2.28	0.51
1:B:1:MET:CE	5:B:547:HOH:O	2.58	0.50
1:C:157[A]:LYS:HD2	5:C:1054:HOH:O	2.11	0.50
1:D:232:ARG:NH1	2:D:507:CIT:O4	2.45	0.50
1:A:108:MET:HE3	1:B:108:MET:HE1	1.94	0.49
1:A:108:MET:SD	1:B:108:MET:HE2	2.53	0.49
1:C:42:ALA:HB2	1:D:4:ILE:HD11	1.95	0.48
1:C:93:SER:OG	1:C:96:GLN:NE2	2.46	0.48
1:D:287[A]:LYS:NZ	5:D:972:HOH:O	2.46	0.47
1:D:243:LYS:HD2	5:D:1055:HOH:O	2.12	0.47
1:B:1:MET:HE3	5:B:547:HOH:O	2.15	0.47
1:D:30:ALA:O	1:D:360:GLY:HA3	2.15	0.47
1:B:127:ALA:HB1	1:B:128:PRO:CD	2.44	0.47
1:D:256:GLY:HA3	2:D:503:CIT:H22	1.97	0.46
1:D:279:LYS:HE3	5:D:455:HOH:O	2.14	0.46
1:D:87:VAL:HG22	1:D:237:LEU:HD22	1.98	0.46
1:A:287[A]:LYS:HG3	5:A:539:HOH:O	2.16	0.46
1:B:30:ALA:O	1:B:360:GLY:HA3	2.16	0.46
1:B:79:LYS:NZ	2:B:502:CIT:O5	2.49	0.46
1:B:361:THR:N	1:B:362:PRO:CD	2.79	0.46
1:C:7[B]:GLN:HE21	1:C:7[B]:GLN:HB2	1.46	0.46
1:B:361:THR:N	1:B:362:PRO:HD3	2.31	0.45
1:D:7[B]:GLN:HG3	1:D:8:ASP:HB2	1.99	0.45
1:B:243:LYS:HE2	2:B:502:CIT:O7	2.17	0.45
1:D:273:ALA:HA	1:D:278:PHE:CG	2.53	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ALA:O	1:A:360:GLY:HA3	2.17	0.44
1:C:1:MET:HB3	1:C:1:MET:HE3	1.77	0.44
1:A:29[B]:ILE:HD12	1:A:340:ASN:ND2	2.33	0.44
1:A:311:HIS:HB2	1:A:362:PRO:HD3	1.99	0.44
1:C:116:ASN:HB3	1:C:140:GLU:OE2	2.18	0.43
1:B:51:TYR:HE1	2:B:506:CIT:H22	1.84	0.43
1:D:410:LEU:HD23	1:D:410:LEU:HA	1.91	0.43
1:C:240:GLU:OE1	1:C:243[A]:LYS:CD	2.64	0.43
1:C:361:THR:N	1:C:362:PRO:CD	2.82	0.43
1:D:29:ILE:HG23	1:D:340:ASN:ND2	2.33	0.43
1:D:326:LYS:O	1:D:330:GLU:HG3	2.19	0.43
1:A:29[B]:ILE:HD11	1:A:338:THR:HG22	2.01	0.42
1:C:108:MET:SD	1:C:135:PHE:HB3	2.59	0.42
1:D:4:ILE:HG21	1:D:4:ILE:HD13	1.64	0.42
1:A:320:SER:O	1:A:321[A]:ILE:HD13	2.19	0.42
1:C:287[B]:LYS:CD	5:C:818:HOH:O	2.66	0.42
1:C:291:VAL:HG21	1:C:371[B]:GLU:HG2	2.00	0.42
1:B:1:MET:HE2	5:B:547:HOH:O	2.19	0.42
1:D:130:ASN:OD1	1:D:132:SER:N	2.52	0.42
1:A:34:PHE:CE2	1:A:363:ALA:HB1	2.55	0.42
1:A:361:THR:N	1:A:362:PRO:CD	2.83	0.41
1:B:51:TYR:HE1	2:B:506:CIT:C2	2.33	0.41
2:B:506:CIT:O3	2:B:506:CIT:O7	2.32	0.41
1:D:4:ILE:HG22	1:D:4:ILE:O	2.18	0.41
1:B:51:TYR:CD2	1:B:61:TYR:CD2	3.08	0.41
1:C:7[B]:GLN:OE1	1:D:272:GLU:OE2	2.38	0.41
1:A:291:VAL:CG1	1:A:375[B]:GLU:HG2	2.50	0.41
1:D:90:GLN:N	1:D:91:PRO:CD	2.83	0.41
1:B:99:MET:HE2	1:B:131[A]:PHE:CG	2.56	0.41
1:C:248:LYS:HA	1:C:251:PHE:O	2.20	0.41
1:A:93:SER:OG	1:A:96:GLN:NE2	2.54	0.41
1:D:311:HIS:HD1	1:D:311:HIS:H	1.68	0.41
1:A:350:LYS:O	1:A:353:VAL:HG13	2.20	0.40
1:D:321:ILE:HD11	1:D:385:LEU:HB3	2.03	0.40
1:B:29:ILE:HD13	1:B:339:CYS:HA	2.03	0.40
1:B:291:VAL:HG13	1:B:375[A]:GLU:HG3	2.04	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)
1:C:294[B]:GLU:OE2	5:A:859:HOH:O[3_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	407/415 (98%)	397 (98%)	9 (2%)	1 (0%)	43	34
1	B	408/415 (98%)	397 (97%)	9 (2%)	2 (0%)	24	14
1	C	414/415 (100%)	402 (97%)	10 (2%)	2 (0%)	24	14
1	D	415/415 (100%)	402 (97%)	11 (3%)	2 (0%)	24	14
All	All	1644/1660 (99%)	1598 (97%)	39 (2%)	7 (0%)	30	19

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	226	LYS
1	B	30	ALA
1	C	226	LYS
1	D	30	ALA
1	D	226	LYS
1	B	226	LYS
1	C	30	ALA

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	340/337 (101%)	335 (98%)	5 (2%)	57	47
1	B	338/337 (100%)	334 (99%)	4 (1%)	63	57
1	C	346/337 (103%)	343 (99%)	3 (1%)	70	67
1	D	345/337 (102%)	342 (99%)	3 (1%)	70	67
All	All	1369/1348 (102%)	1354 (99%)	15 (1%)	70	60

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	287[A]	LYS
1	A	287[B]	LYS
1	A	353	VAL
1	A	389	LYS
1	A	396	GLN
1	B	65[A]	GLU
1	B	65[B]	GLU
1	B	112	VAL
1	B	243	LYS
1	C	116	ASN
1	C	340	ASN
1	C	392	GLU
1	D	9[A]	LYS
1	D	9[B]	LYS
1	D	130	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	340	ASN
1	B	7	GLN
1	B	20	GLN
1	B	25	ASN
1	B	137	ASN
1	B	342	ASN
1	B	396	GLN
1	C	96	GLN
1	C	137	ASN
1	D	162	HIS
1	D	340	ASN
1	D	396	GLN

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 ⓘ

13 ligands are 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$
2	CIT	A	505	-	12,12,12	0.95	0	17,17,17	2.01	6 (35%)
2	CIT	D	504	-	12,12,12	1.13	0	17,17,17	1.14	1 (5%)
2	CIT	C	508	-	12,12,12	0.99	0	17,17,17	1.99	5 (29%)
3	GLY	B	511	-	4,4,4	1.10	1 (25%)	3,4,4	1.61	1 (33%)
3	GLY	B	509	-	4,4,4	1.11	0	3,4,4	1.79	2 (66%)
2	CIT	D	507	-	12,12,12	1.60	3 (25%)	17,17,17	2.24	7 (41%)
4	EDO	C	512	-	3,3,3	0.92	0	2,2,2	0.64	0
2	CIT	B	501	-	12,12,12	1.74	1 (8%)	17,17,17	3.27	9 (52%)
2	CIT	D	503	-	12,12,12	2.24	4 (33%)	17,17,17	3.19	9 (52%)
4	EDO	C	513	-	3,3,3	0.89	0	2,2,2	0.33	0
3	GLY	A	510	-	4,4,4	1.51	1 (25%)	3,4,4	0.72	0
2	CIT	B	506	-	12,12,12	2.48	4 (33%)	17,17,17	3.63	12 (70%)
2	CIT	B	502	-	12,12,12	1.00	0	17,17,17	1.91	4 (23%)

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
2	CIT	A	505	-	-	4/16/16/16	-
2	CIT	D	504	-	-	0/16/16/16	-
2	CIT	C	508	-	-	2/16/16/16	-
3	GLY	B	511	-	-	2/2/2/2	-
3	GLY	B	509	-	-	2/2/2/2	-
2	CIT	D	507	-	-	5/16/16/16	-
4	EDO	C	512	-	-	1/1/1/1	-
2	CIT	B	501	-	-	9/16/16/16	-
2	CIT	D	503	-	-	7/16/16/16	-
4	EDO	C	513	-	-	0/1/1/1	-
3	GLY	A	510	-	-	2/2/2/2	-
2	CIT	B	506	-	-	5/16/16/16	-
2	CIT	B	502	-	-	2/16/16/16	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	506	CIT	C4-C3	-5.83	1.46	1.54
2	D	503	CIT	C2-C3	-5.66	1.47	1.54
2	B	501	CIT	C2-C3	-4.29	1.48	1.54
2	B	506	CIT	C2-C3	-4.07	1.48	1.54
2	D	503	CIT	O2-C1	-3.19	1.20	1.30
3	A	510	GLY	OXT-C	-2.68	1.21	1.30
2	B	506	CIT	O4-C5	-2.62	1.22	1.30
2	B	506	CIT	C4-C5	-2.49	1.43	1.50
2	D	507	CIT	O4-C5	-2.33	1.23	1.30
2	D	507	CIT	C3-C6	-2.32	1.51	1.53
2	D	503	CIT	O5-C6	2.30	1.29	1.22
2	D	507	CIT	O7-C3	2.12	1.47	1.43
3	B	511	GLY	OXT-C	-2.06	1.24	1.30
2	D	503	CIT	C2-C1	-2.02	1.44	1.50

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	503	CIT	C2-C3-C6	-9.36	89.33	110.03
2	B	506	CIT	C4-C3-C2	-7.66	89.65	109.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	CIT	O1-C1-C2	-7.39	102.02	122.95
2	B	506	CIT	C2-C3-C6	7.17	125.90	110.03
2	B	501	CIT	O2-C1-C2	6.00	133.35	114.35
2	B	501	CIT	C2-C3-C6	-5.50	97.86	110.03
2	B	501	CIT	O7-C3-C6	4.89	115.90	108.96
2	B	502	CIT	O6-C6-C3	4.82	122.38	113.14
2	A	505	CIT	O6-C6-C3	4.61	121.99	113.14
2	B	506	CIT	C4-C3-C6	-4.36	100.38	110.03
2	D	503	CIT	C4-C3-C6	4.35	119.65	110.03
2	D	507	CIT	C3-C4-C5	-4.34	102.06	113.92
2	C	508	CIT	O6-C6-C3	4.24	121.27	113.14
2	B	506	CIT	O2-C1-C2	4.09	127.29	114.35
2	B	506	CIT	O2-C1-O1	-4.04	112.93	123.33
2	D	503	CIT	O6-C6-C3	4.03	120.86	113.14
2	D	507	CIT	O7-C3-C6	-3.73	103.67	108.96
2	B	506	CIT	O6-C6-C3	3.50	119.86	113.14
2	B	506	CIT	O7-C3-C4	3.48	117.30	109.38
2	C	508	CIT	C3-C4-C5	-3.27	104.98	113.92
2	B	506	CIT	C3-C2-C1	-3.24	105.06	113.92
2	C	508	CIT	O7-C3-C6	-3.17	104.47	108.96
2	D	503	CIT	O4-C5-C4	3.13	124.27	114.35
2	D	507	CIT	C4-C3-C6	-3.08	103.22	110.03
2	D	507	CIT	O7-C3-C4	3.04	116.31	109.38
2	D	503	CIT	O7-C3-C6	3.04	113.27	108.96
2	D	507	CIT	C2-C3-C6	3.01	116.69	110.03
2	A	505	CIT	O1-C1-C2	-2.97	114.55	122.95
2	D	503	CIT	O4-C5-O3	-2.95	115.74	123.33
2	A	505	CIT	O2-C1-C2	2.94	123.67	114.35
2	B	506	CIT	O7-C3-C6	2.76	112.88	108.96
2	A	505	CIT	O4-C5-C4	2.74	123.04	114.35
2	B	502	CIT	O2-C1-C2	2.49	122.23	114.35
2	B	501	CIT	O7-C3-C4	2.48	115.04	109.38
2	B	501	CIT	C4-C3-C2	-2.46	102.99	109.31
2	C	508	CIT	O2-C1-C2	2.45	122.11	114.35
2	A	505	CIT	O6-C6-O5	-2.43	116.07	123.86
2	B	506	CIT	O4-C5-O3	2.32	129.30	123.33
2	D	507	CIT	O6-C6-C3	2.30	117.55	113.14
2	D	503	CIT	O2-C1-C2	-2.27	107.14	114.35
2	A	505	CIT	C4-C3-C6	2.25	115.02	110.03
3	B	509	GLY	OXT-C-CA	2.24	122.28	113.38
2	B	501	CIT	O5-C6-C3	2.20	126.35	122.09
2	D	503	CIT	O5-C6-C3	-2.18	117.86	122.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	508	CIT	O3-C5-C4	-2.18	116.77	122.95
2	D	503	CIT	C3-C2-C1	-2.18	107.96	113.92
2	D	507	CIT	O2-C1-C2	2.15	121.17	114.35
3	B	511	GLY	OXT-C-CA	2.15	121.93	113.38
2	B	501	CIT	C3-C4-C5	-2.15	108.06	113.92
2	B	506	CIT	O4-C5-C4	-2.09	107.71	114.35
2	B	502	CIT	O5-C6-C3	-2.08	118.07	122.09
2	B	506	CIT	C3-C4-C5	-2.07	108.27	113.92
2	B	502	CIT	O7-C3-C6	-2.04	106.06	108.96
3	B	509	GLY	OXT-C-O	-2.03	118.11	123.33
2	D	504	CIT	O2-C1-C2	2.02	120.75	114.35
2	B	501	CIT	C3-C2-C1	2.00	119.40	113.92

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	CIT	C1-C2-C3-O7
2	B	501	CIT	C1-C2-C3-C4
2	B	501	CIT	C1-C2-C3-C6
2	B	501	CIT	C2-C3-C6-O5
2	B	501	CIT	C2-C3-C6-O6
2	B	501	CIT	O7-C3-C6-O5
2	B	501	CIT	O7-C3-C6-O6
2	B	506	CIT	C2-C3-C6-O5
2	B	506	CIT	C2-C3-C6-O6
2	D	503	CIT	O7-C3-C6-O5
2	D	503	CIT	O7-C3-C6-O6
2	D	503	CIT	C4-C3-C6-O5
2	D	503	CIT	C4-C3-C6-O6
2	D	507	CIT	C2-C3-C4-C5
2	D	507	CIT	O7-C3-C4-C5
2	D	507	CIT	C6-C3-C4-C5
3	B	509	GLY	O-C-CA-N
3	B	509	GLY	OXT-C-CA-N
3	B	511	GLY	OXT-C-CA-N
3	A	510	GLY	O-C-CA-N
2	A	505	CIT	C3-C4-C5-O4
2	A	505	CIT	C3-C4-C5-O3
2	B	506	CIT	C2-C3-C4-C5
2	B	506	CIT	C6-C3-C4-C5
3	B	511	GLY	O-C-CA-N

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Mol	Chain	Res	Type	Atoms
2	B	501	CIT	O1-C1-C2-C3
2	B	501	CIT	O2-C1-C2-C3
2	D	503	CIT	C2-C3-C6-O5
2	D	503	CIT	C2-C3-C6-O6
2	D	507	CIT	O2-C1-C2-C3
2	B	506	CIT	O7-C3-C4-C5
2	A	505	CIT	C2-C3-C6-O5
2	A	505	CIT	C2-C3-C6-O6
2	B	502	CIT	O2-C1-C2-C3
2	D	507	CIT	O1-C1-C2-C3
2	B	502	CIT	O1-C1-C2-C3
3	A	510	GLY	OXT-C-CA-N
2	C	508	CIT	O1-C1-C2-C3
2	C	508	CIT	O2-C1-C2-C3
4	C	512	EDO	O1-C1-C2-O2
2	D	503	CIT	C2-C3-C4-C5

There are no ring outliers.

7 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	505	CIT	1	0
2	C	508	CIT	1	0
2	D	507	CIT	3	0
2	B	501	CIT	1	0
2	D	503	CIT	2	0
2	B	506	CIT	5	0
2	B	502	CIT	2	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 ⓘ

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	402/415 (96%)	-0.33	6 (1%) 72 79	16, 26, 45, 63	11 (2%)
1	B	404/415 (97%)	-0.26	9 (2%) 62 69	15, 28, 40, 56	8 (1%)
1	C	398/415 (95%)	-0.33	5 (1%) 75 82	14, 25, 41, 47	23 (5%)
1	D	402/415 (96%)	-0.18	7 (1%) 69 75	14, 30, 45, 56	17 (4%)
All	All	1606/1660 (96%)	-0.27	27 (1%) 69 75	14, 27, 44, 63	59 (3%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	129	VAL	9.1
1	D	352[A]	PHE	5.2
1	B	127	ALA	4.3
1	D	129	VAL	4.1
1	B	60	TYR	4.0
1	B	61	TYR	3.7
1	D	60	TYR	3.6
1	D	128	PRO	3.2
1	A	352	PHE	3.2
1	B	352	PHE	3.1
1	B	119	HIS	3.1
1	A	412	GLN	3.0
1	A	62	GLY	2.9
1	B	118	SER	2.9
1	A	117	LEU	2.9
1	D	61	TYR	2.9
1	D	350[A]	LYS	2.8
1	C	135	PHE	2.7
1	B	62	GLY	2.6
1	A	119	HIS	2.5
1	C	62	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	129	VAL	2.5
1	A	118	SER	2.3
1	D	135	PHE	2.2
1	B	128	PRO	2.2
1	C	6[A]	LYS	2.1
1	C	116	ASN	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(Å ²)	Q<0.9
3	GLY	B	509	5/5	0.80	0.18	55,55,56,56	0
2	CIT	D	507	13/13	0.83	0.12	48,51,59,61	0
2	CIT	C	508	13/13	0.84	0.10	39,48,61,63	0
2	CIT	D	504	13/13	0.86	0.10	37,46,54,57	0
2	CIT	B	501	13/13	0.87	0.10	41,45,51,52	0
2	CIT	D	503	13/13	0.88	0.11	38,40,57,59	0
2	CIT	B	506	13/13	0.88	0.12	37,43,49,56	0
3	GLY	B	511	5/5	0.88	0.15	67,67,68,68	0
4	EDO	C	512	4/4	0.88	0.13	30,36,38,38	0
2	CIT	A	505	13/13	0.89	0.10	28,37,55,58	0
2	CIT	B	502	13/13	0.90	0.09	42,48,60,62	0
3	GLY	A	510	5/5	0.92	0.13	51,52,54,54	0
4	EDO	C	513	4/4	0.95	0.08	31,35,35,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.