



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 01:33 PM EDT

PDB ID : 6YMZ
Title : Structure of the CheB methylsterase from *P. atrosepticum* SCRI1043
Authors : Gavira, J.A.; Krell, T.; Velando-Soriano, F.; Matilla, M.A.
Deposited on : 2020-04-10
Resolution : 2.30 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 2.37.1
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) : 2.37.1

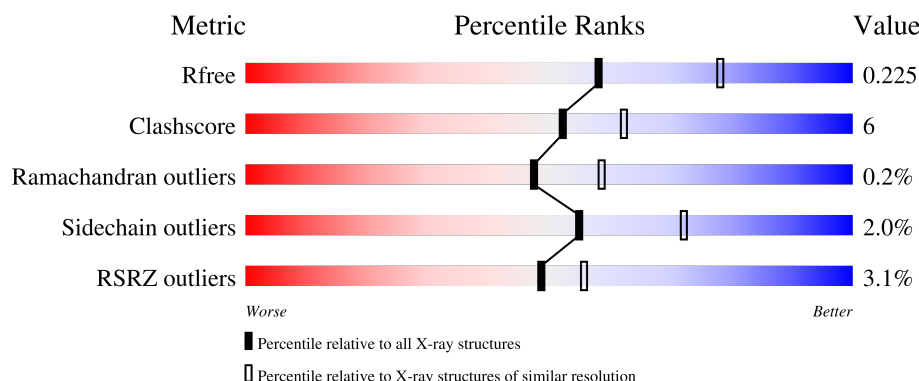
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	356	 2% 85% 11% . .
1	B	356	 4% 85% 10% . .
1	C	356	 3% 84% 11% . .
1	D	356	 3% 84% 9% . 6%
1	E	356	 3% 71% 21% 8%

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
4	GOL	E	403[A]	-	-	-	X
4	GOL	E	403[B]	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13986 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 Protein-glutamate methyltransferase/protein-glutamine glutaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	18	0
			2753	1729	507	496	21			
1	B	341	Total	C	N	O	S	0	13	0
			2707	1696	495	495	21			
1	C	341	Total	C	N	O	S	0	14	0
			2713	1707	496	489	21			
1	D	336	Total	C	N	O	S	0	10	0
			2636	1654	479	481	22			
1	E	328	Total	C	N	O	S	0	15	0
			2624	1645	481	478	20			

There are 30 discrepancies between the modelled and reference sequences:

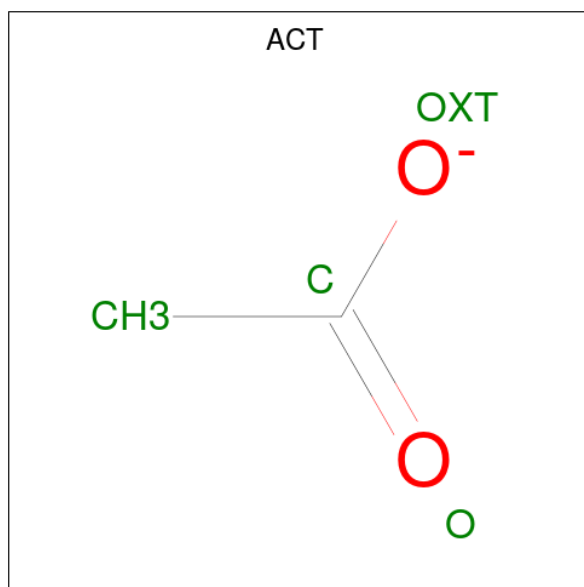
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP Q6D6I7
A	-4	SER	-	expression tag	UNP Q6D6I7
A	-3	HIS	-	expression tag	UNP Q6D6I7
A	-2	MET	-	expression tag	UNP Q6D6I7
A	-1	ALA	-	expression tag	UNP Q6D6I7
A	0	SER	-	expression tag	UNP Q6D6I7
B	-5	GLY	-	expression tag	UNP Q6D6I7
B	-4	SER	-	expression tag	UNP Q6D6I7
B	-3	HIS	-	expression tag	UNP Q6D6I7
B	-2	MET	-	expression tag	UNP Q6D6I7
B	-1	ALA	-	expression tag	UNP Q6D6I7
B	0	SER	-	expression tag	UNP Q6D6I7
C	-5	GLY	-	expression tag	UNP Q6D6I7
C	-4	SER	-	expression tag	UNP Q6D6I7
C	-3	HIS	-	expression tag	UNP Q6D6I7
C	-2	MET	-	expression tag	UNP Q6D6I7
C	-1	ALA	-	expression tag	UNP Q6D6I7
C	0	SER	-	expression tag	UNP Q6D6I7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-5	GLY	-	expression tag	UNP Q6D6I7
D	-4	SER	-	expression tag	UNP Q6D6I7
D	-3	HIS	-	expression tag	UNP Q6D6I7
D	-2	MET	-	expression tag	UNP Q6D6I7
D	-1	ALA	-	expression tag	UNP Q6D6I7
D	0	SER	-	expression tag	UNP Q6D6I7
E	-5	GLY	-	expression tag	UNP Q6D6I7
E	-4	SER	-	expression tag	UNP Q6D6I7
E	-3	HIS	-	expression tag	UNP Q6D6I7
E	-2	MET	-	expression tag	UNP Q6D6I7
E	-1	ALA	-	expression tag	UNP Q6D6I7
E	0	SER	-	expression tag	UNP Q6D6I7

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Na 2 2	0	0
3	B	2	Total Na 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total	Na	0	0
			2	2		
3	D	1	Total	Na	0	0
			1	1		
3	E	1	Total	Na	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	1
			12	6	6		
4	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	O	S	0	0
			5	4	1		

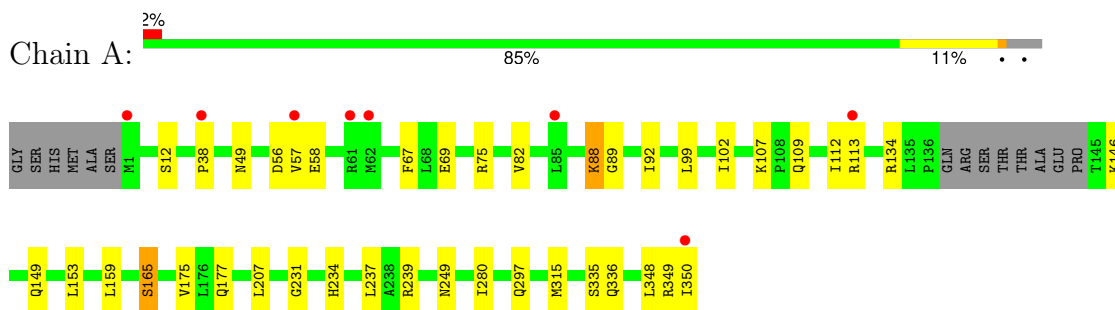
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	105	Total	O	0	0
			105	105		
6	B	99	Total	O	0	0
			99	99		
6	C	121	Total	O	0	0
			121	121		
6	D	99	Total	O	0	0
			99	99		
6	E	64	Total	O	0	0
			64	64		

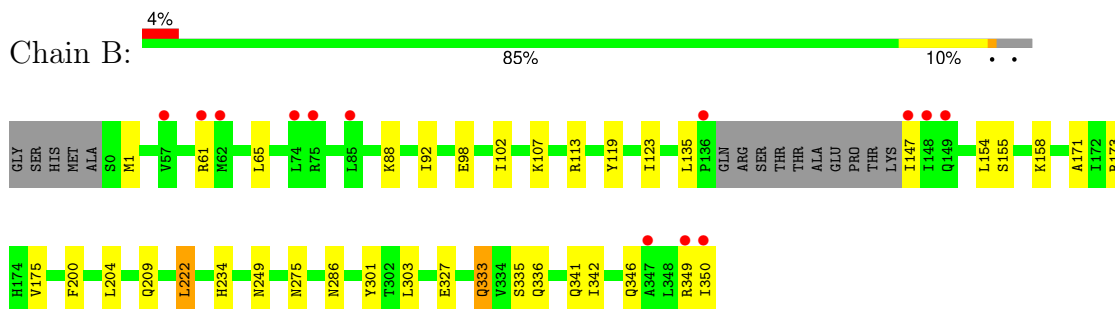
3 Residue-property plots [i](#)

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.

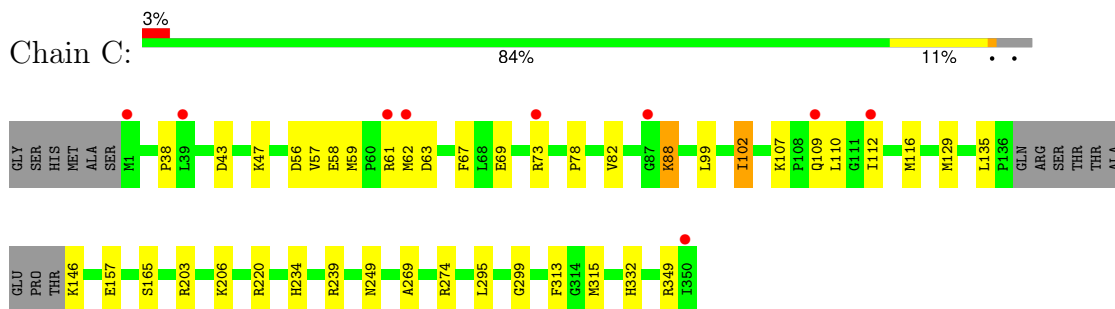
- Molecule 1: Protein-glutamate methyltransferase/protein-glutamine glutaminase



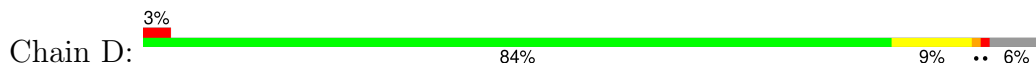
- Molecule 1: Protein-glutamate methyltransferase/protein-glutamine glutaminase

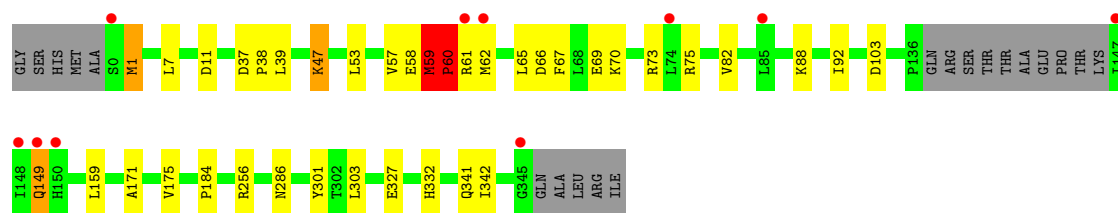


- Molecule 1: Protein-glutamate methyltransferase/protein-glutamine glutaminase

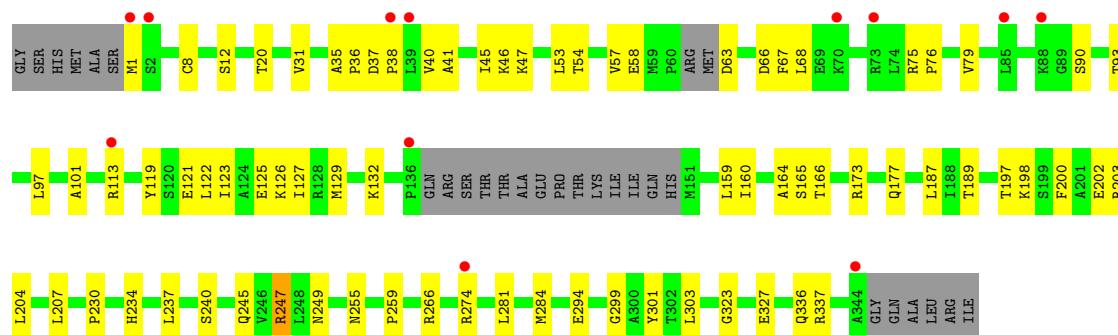


- Molecule 1: Protein-glutamate methyltransferase/protein-glutamine glutaminase





● Molecule 1: Protein-glutamate methylesterase/protein-glutamine glutaminase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	148.74Å 148.74Å 206.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.31 – 2.30 63.31 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.5 (63.31-2.30) 98.5 (63.31-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.192 , 0.226 0.193 , 0.225	Depositor DCC
R_{free} test set	4881 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 37.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.028 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k 0.028 for -1/2*h+1/2*k+1/2*l,1/2*h-1/2*k+1/2*l,h+k 0.037 for -1/2*h-1/2*k+1/2*l,-1/2*h-1/2*k-1/2*l,h-k 0.038 for -1/2*h-1/2*k-1/2*l,-1/2*h-1/2*k+1/2*l,-h+k 0.417 for -h,k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13986	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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: SO4, ACT, GOL, 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	0.29	0/2803	0.52	0/3782
1	B	0.33	1/2748 (0.0%)	0.50	0/3713
1	C	0.29	0/2757	0.48	0/3721
1	D	0.42	3/2678 (0.1%)	0.55	3/3619 (0.1%)
1	E	0.28	0/2663	0.48	0/3595
All	All	0.32	4/13649 (0.0%)	0.50	3/18430 (0.0%)

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	D	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	60	PRO	N-CA	12.52	1.68	1.47
1	D	60	PRO	C-N	-7.67	1.16	1.34
1	B	222	LEU	C-N	7.53	1.48	1.34
1	D	59	MET	C-N	5.35	1.44	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	60	PRO	O-C-N	-10.23	106.34	122.70
1	D	60	PRO	CA-N-CD	-7.03	101.66	111.50
1	D	159	LEU	CA-CB-CG	5.83	128.70	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	58[A]	GLU	Mainchain
1	D	58[B]	GLU	Mainchain
1	D	60	PRO	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	2753	0	2874	25	0
1	B	2707	0	2795	31	0
1	C	2713	0	2831	29	0
1	D	2636	0	2720	27	0
1	E	2624	0	2713	45	0
2	A	4	0	3	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
4	A	12	0	16	0	0
4	C	12	0	16	3	0
4	D	6	0	8	2	0
4	E	18	0	24	0	0
5	E	5	0	0	0	0
6	A	105	0	0	5	0
6	B	99	0	0	6	0
6	C	121	0	0	4	2
6	D	99	0	0	3	2
6	E	64	0	0	8	0
All	All	13986	0	14000	152	3

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 (152) 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:D:60:PRO:CA	1:D:60:PRO:N	1.68	1.50
1:C:313:PHE:H	4:C:404:GOL:H32	1.23	1.02
1:A:348:LEU:HG	1:B:222:LEU:HD11	1.48	0.95
1:B:154:LEU:HA	1:B:346:GLN:HG2	1.53	0.90
1:B:158:LYS:HE2	6:B:503:HOH:O	1.72	0.87
1:C:313:PHE:N	4:C:404:GOL:H32	2.00	0.77
1:E:37:ASP:HB3	6:E:513:HOH:O	1.87	0.75
1:E:247[A]:ARG:NH2	6:E:501:HOH:O	2.08	0.70
1:B:158:LYS:HE3	1:B:346:GLN:OE1	1.92	0.69
1:C:38:PRO:HD2	1:C:62:MET:HG2	1.73	0.68
1:E:159:LEU:HD21	1:E:187:LEU:HG	1.75	0.68
1:E:234:HIS:HB2	1:E:249:ASN:HB3	1.74	0.68
1:E:202[B]:GLU:OE1	6:E:502:HOH:O	2.12	0.67
1:C:313:PHE:H	4:C:404:GOL:C3	2.04	0.67
1:B:275:ASN:OD1	6:B:503:HOH:O	2.13	0.67
1:B:155:SER:O	6:B:504:HOH:O	2.13	0.66
1:E:323:GLY:O	6:E:503:HOH:O	2.13	0.66
1:A:336:GLN:NE2	6:A:501:HOH:O	2.12	0.65
1:A:149:GLN:NE2	6:A:504:HOH:O	2.30	0.65
1:E:35:ALA:HB1	1:E:40:VAL:HG23	1.78	0.65
1:E:76:PRO:O	6:E:504:HOH:O	2.15	0.65
1:B:98:GLU:HG2	1:B:173[B]:ARG:HH12	1.60	0.64
1:B:102:ILE:HD11	1:B:135:LEU:HD23	1.79	0.64
1:B:341:GLN:NE2	6:B:502:HOH:O	2.04	0.64
1:B:61:ARG:NH1	6:B:507:HOH:O	2.33	0.61
1:B:119:TYR:CZ	1:B:123:ILE:HD11	2.36	0.61
1:C:56:ASP:O	6:C:502:HOH:O	2.17	0.60
1:E:119:TYR:CE2	1:E:123:ILE:HD11	2.36	0.60
1:E:173:ARG:HD3	1:E:203:ARG:HH21	1.67	0.60
1:E:187:LEU:HD11	1:E:237:LEU:HD13	1.84	0.60
1:C:109[A]:GLN:HB2	1:C:112:ILE:HD12	1.85	0.59
1:E:38:PRO:HB3	1:E:67:PHE:CD1	2.38	0.59
1:D:7:LEU:HB3	1:D:53:LEU:HD12	1.84	0.59
1:D:75[B]:ARG:NH1	6:D:501:HOH:O	2.07	0.58
1:C:38:PRO:HD3	1:C:59:MET:SD	2.44	0.58
1:C:274:ARG:HG2	1:C:299:GLY:HA3	1.86	0.58
1:E:36:PRO:HD2	1:E:40:VAL:HG21	1.85	0.57
1:E:173:ARG:HD3	1:E:203:ARG:NH2	2.19	0.57
1:C:61[B]:ARG:NH2	6:C:501:HOH:O	2.07	0.56
1:E:274[A]:ARG:HG2	1:E:299:GLY:HA3	1.87	0.56
1:A:234:HIS:HB2	1:A:249:ASN:HB3	1.88	0.55
1:E:123:ILE:O	1:E:127[B]:ILE:HG22	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:PRO:HB3	1:C:67:PHE:CG	2.43	0.54
1:E:240:SER:HB3	1:E:245:GLN:OE1	2.07	0.54
1:E:20:THR:HA	1:E:31:VAL:HG21	1.89	0.54
1:A:109:GLN:HB3	1:A:112:ILE:HD12	1.89	0.54
1:E:177:GLN:HE21	1:E:207:LEU:HB3	1.74	0.53
1:A:165:SER:HB2	1:A:315:MET:SD	2.48	0.53
1:A:134:ARG:NH1	6:A:507:HOH:O	2.35	0.53
1:B:303:LEU:HD23	1:B:327:GLU:HB3	1.91	0.53
1:E:189:THR:HG23	1:E:259:PRO:HB2	1.89	0.53
1:E:121:GLU:O	1:E:125:GLU:HG2	2.08	0.53
1:B:65:LEU:HD11	1:B:92:ILE:HG23	1.91	0.52
1:C:269:ALA:HB2	1:C:295:LEU:HA	1.91	0.52
1:E:197:THR:HG21	1:E:230:PRO:HA	1.91	0.52
1:A:56:ASP:OD1	1:A:58:GLU:HG2	2.09	0.52
1:D:301:TYR:CG	1:D:341:GLN:HG2	2.44	0.52
1:D:332:HIS:ND1	6:D:505:HOH:O	2.34	0.52
1:A:348:LEU:HD22	1:A:350:ILE:HD11	1.92	0.51
1:D:37:ASP:HB2	1:D:62[B]:MET:SD	2.50	0.51
1:A:297:GLN:HA	1:B:350:ILE:O	2.11	0.51
1:B:88[A]:LYS:O	1:B:92:ILE:HG13	2.11	0.51
1:B:346:GLN:HG3	6:B:554:HOH:O	2.10	0.51
1:C:69:GLU:HG2	1:C:99:LEU:HD22	1.92	0.51
1:C:157:GLU:OE1	1:C:239:ARG:NH2	2.44	0.51
1:D:70:LYS:HA	1:D:73:ARG:HG2	1.93	0.50
1:A:348:LEU:HG	1:B:222:LEU:CD1	2.31	0.50
1:B:234:HIS:HB2	1:B:249:ASN:HB3	1.92	0.50
1:C:102[B]:ILE:HD11	1:C:129:MET:SD	2.51	0.50
1:E:58:GLU:O	6:E:505:HOH:O	2.18	0.50
1:C:234:HIS:HB2	1:C:249:ASN:HB3	1.92	0.50
1:C:57:VAL:HG23	1:C:58:GLU:HG2	1.93	0.50
1:A:159[A]:LEU:HD21	1:A:237:LEU:HD22	1.94	0.49
1:B:301:TYR:CD2	1:B:341:GLN:HG3	2.47	0.49
1:E:301:TYR:HE2	1:E:303:LEU:HD21	1.78	0.49
1:E:266:ARG:NH1	1:E:294:GLU:OE1	2.37	0.49
1:A:82:VAL:HG12	1:A:107:LYS:HE3	1.95	0.49
1:C:38:PRO:HB3	1:C:67:PHE:CD1	2.48	0.49
1:A:69:GLU:HG2	1:A:99:LEU:HD22	1.94	0.49
1:C:110:LEU:HD23	1:C:116:MET:HA	1.94	0.48
1:D:184:PRO:HG3	1:D:342:ILE:HG21	1.94	0.48
1:D:60:PRO:N	1:D:60:PRO:C	2.60	0.48
1:B:88[B]:LYS:O	1:B:92:ILE:HG13	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:ASP:HB2	1:D:59:MET:HE2	1.94	0.48
1:A:177:GLN:HE21	1:A:207:LEU:HB3	1.77	0.48
1:C:332:HIS:ND1	6:C:506:HOH:O	2.35	0.48
1:E:57:VAL:O	1:E:57:VAL:HG23	2.14	0.48
1:A:146:LYS:O	1:A:146:LYS:HG3	2.13	0.48
1:D:103:ASP:OD1	4:D:402:GOL:H12	2.14	0.47
1:B:175:VAL:O	1:B:335:SER:OG	2.32	0.47
1:E:129:MET:HG3	1:E:284:MET:SD	2.54	0.47
1:E:46:LYS:HG2	1:E:75:ARG:HH12	1.79	0.47
1:D:38:PRO:HB3	1:D:67:PHE:CG	2.49	0.47
1:E:164:ALA:HB2	1:E:281:LEU:HB2	1.97	0.47
1:A:49:ASN:OD1	1:A:75:ARG:NH2	2.48	0.47
1:E:165:SER:OG	1:E:166:THR:N	2.47	0.47
1:E:200:PHE:CZ	1:E:204:LEU:HD11	2.50	0.47
1:E:113[A]:ARG:H	1:E:113[A]:ARG:HD2	1.80	0.46
1:B:158:LYS:NZ	1:B:342:ILE:O	2.47	0.46
1:B:286:ASN:HB3	1:D:1:MET:SD	2.55	0.46
1:D:11:ASP:HB2	1:D:59:MET:CE	2.45	0.46
1:E:90:SER:OG	6:E:506:HOH:O	2.21	0.46
1:C:203:ARG:HA	1:C:206[A]:LYS:HE3	1.97	0.46
1:E:177:GLN:NE2	1:E:207:LEU:HD22	2.31	0.45
1:A:88[A]:LYS:HG2	1:A:89:GLY:N	2.31	0.45
1:D:256:ARG:HB2	1:D:256:ARG:NH1	2.31	0.45
1:A:57:VAL:HB	1:A:92:ILE:HG21	1.98	0.45
1:E:93:THR:O	1:E:97:LEU:HG	2.17	0.45
1:B:171:ALA:O	1:B:175:VAL:HG23	2.17	0.44
1:D:65:LEU:HD11	1:D:92:ILE:HG23	1.99	0.44
1:D:303:LEU:HD23	1:D:327:GLU:HB3	2.00	0.44
1:C:88[B]:LYS:HA	1:C:88[B]:LYS:HD3	1.78	0.44
1:B:107:LYS:H	1:B:107:LYS:HG2	1.62	0.44
1:A:231:GLY:O	6:A:502:HOH:O	2.21	0.44
1:B:119:TYR:O	1:B:123:ILE:HG12	2.17	0.43
1:D:70:LYS:HA	1:D:73:ARG:CG	2.48	0.43
1:D:38:PRO:HD2	1:D:62[B]:MET:SD	2.58	0.43
1:A:102:ILE:O	6:A:503:HOH:O	2.22	0.43
1:E:122:LEU:HD11	1:E:126:LYS:HE2	2.00	0.43
1:E:53:LEU:HB3	1:E:79:VAL:HG22	2.01	0.42
1:D:61[B]:ARG:H	1:D:61[B]:ARG:HD2	1.84	0.42
1:B:147:ILE:HD13	1:B:209:GLN:OE1	2.18	0.42
1:E:68:LEU:HD21	1:E:101:ALA:HB2	2.01	0.42
1:A:153:LEU:O	1:C:220:ARG:NH2	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:SER:HB2	1:C:315:MET:SD	2.59	0.42
1:C:82:VAL:HG12	1:C:107:LYS:HE2	2.02	0.42
1:E:41:ALA:O	1:E:45:ILE:HG12	2.20	0.42
1:C:43:ASP:HB2	6:C:504:HOH:O	2.19	0.42
1:C:110:LEU:CD2	1:C:116:MET:HA	2.50	0.42
1:D:69:GLU:O	1:D:73:ARG:HG2	2.20	0.42
1:D:39:LEU:HD13	1:D:39:LEU:HA	1.95	0.41
1:A:38:PRO:HB3	1:A:67:PHE:CG	2.55	0.41
1:B:1:MET:SD	1:D:286:ASN:HB3	2.60	0.41
4:D:402:GOL:H2	6:D:543:HOH:O	2.21	0.41
1:E:47:LYS:HD3	1:E:47:LYS:HA	1.80	0.41
1:D:57:VAL:HG13	1:D:82:VAL:O	2.20	0.41
1:D:171:ALA:O	1:D:175:VAL:HG23	2.20	0.41
1:D:47[B]:LYS:HE3	1:D:47[B]:LYS:HB2	1.77	0.41
1:E:337:ARG:HD2	1:E:337:ARG:HA	1.95	0.41
1:A:280:ILE:HD11	1:A:315:MET:HE3	2.02	0.41
1:A:175:VAL:O	1:A:335:SER:OG	2.39	0.41
1:C:78:PRO:HG3	1:C:135:LEU:HD21	2.01	0.41
1:B:200:PHE:CZ	1:B:204:LEU:HD11	2.55	0.41
1:B:349[B]:ARG:NH2	1:B:349[B]:ARG:HB2	2.35	0.41
1:D:38:PRO:HB3	1:D:67:PHE:CD1	2.56	0.41
1:E:303:LEU:CD2	1:E:327:GLU:HB3	2.51	0.41
1:E:159:LEU:HD23	1:E:160:ILE:N	2.36	0.40
1:C:56:ASP:OD1	1:C:58:GLU:N	2.52	0.40
1:E:58:GLU:HB3	6:E:505:HOH:O	2.20	0.40
1:B:333:GLN:O	1:B:336[B]:GLN:HG3	2.21	0.40
1:C:62:MET:SD	1:C:63:ASP:N	2.94	0.40
1:E:8:CYS:HA	1:E:54[A]:THR:OG1	2.22	0.40

All (3) 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)
6:C:583:HOH:O	6:C:618:HOH:O[4_555]	2.08	0.12
6:D:575:HOH:O	6:D:597:HOH:O[4_555]	2.13	0.07
6:C:557:HOH:O	6:D:570:HOH:O[8_555]	2.16	0.04

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	356/356 (100%)	347 (98%)	9 (2%)	0	100	100
1	B	350/356 (98%)	344 (98%)	6 (2%)	0	100	100
1	C	351/356 (99%)	348 (99%)	3 (1%)	0	100	100
1	D	342/356 (96%)	336 (98%)	3 (1%)	3 (1%)	17	20
1	E	337/356 (95%)	329 (98%)	6 (2%)	2 (1%)	25	31
All	All	1736/1780 (98%)	1704 (98%)	27 (2%)	5 (0%)	47	50

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	60	PRO
1	D	149[A]	GLN
1	D	149[B]	GLN
1	E	255[A]	ASN
1	E	255[B]	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	298/291 (102%)	290 (97%)	8 (3%)	44	61
1	B	292/291 (100%)	290 (99%)	2 (1%)	84	92
1	C	293/291 (101%)	284 (97%)	9 (3%)	40	55
1	D	285/291 (98%)	276 (97%)	9 (3%)	39	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	283/291 (97%)	273 (96%)	10 (4%)	36	50
All	All	1451/1455 (100%)	1413 (97%)	38 (3%)	55	63

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	SER
1	A	88[A]	LYS
1	A	88[B]	LYS
1	A	113[A]	ARG
1	A	113[B]	ARG
1	A	165	SER
1	A	239	ARG
1	A	349	ARG
1	B	113	ARG
1	B	333	GLN
1	C	47[A]	LYS
1	C	47[B]	LYS
1	C	73	ARG
1	C	88[A]	LYS
1	C	88[B]	LYS
1	C	102[A]	ILE
1	C	102[B]	ILE
1	C	146	LYS
1	C	349	ARG
1	D	1	MET
1	D	47[A]	LYS
1	D	47[B]	LYS
1	D	59	MET
1	D	66	ASP
1	D	88[A]	LYS
1	D	88[B]	LYS
1	D	149[A]	GLN
1	D	149[B]	GLN
1	E	1	MET
1	E	12	SER
1	E	63	ASP
1	E	66	ASP
1	E	132[A]	LYS
1	E	132[B]	LYS
1	E	198	LYS
1	E	247[A]	ARG

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Mol	Chain	Res	Type
1	E	247[B]	ARG
1	E	336	GLN

Sometimes 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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 18 ligands modelled in this entry, 8 are monoatomic - leaving 10 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	GOL	E	403[B]	-	5,5,5	0.94	0	5,5,5	1.07	0
2	ACT	A	401	-	3,3,3	1.37	0	3,3,3	1.53	0
4	GOL	A	403	-	5,5,5	0.98	0	5,5,5	1.08	0
4	GOL	A	405	-	5,5,5	1.00	0	5,5,5	1.00	0
4	GOL	D	402	-	5,5,5	0.91	0	5,5,5	1.00	0
4	GOL	E	403[A]	-	5,5,5	0.97	0	5,5,5	1.07	0
4	GOL	C	404	-	5,5,5	0.21	0	5,5,5	0.79	0
5	SO4	E	402	-	4,4,4	0.23	0	6,6,6	0.05	0
4	GOL	C	401	-	5,5,5	0.94	0	5,5,5	1.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	E	404	-	5,5,5	0.91	0	5,5,5	1.09	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	GOL	E	403[B]	-	-	2/4/4/4	-
4	GOL	D	402	-	-	0/4/4/4	-
4	GOL	A	403	-	-	1/4/4/4	-
4	GOL	A	405	-	-	0/4/4/4	-
4	GOL	E	403[A]	-	-	2/4/4/4	-
4	GOL	C	404	-	-	4/4/4/4	-
4	GOL	C	401	-	-	2/4/4/4	-
4	GOL	E	404	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	401	GOL	O1-C1-C2-C3
4	C	404	GOL	O1-C1-C2-O2
4	C	404	GOL	O1-C1-C2-C3
4	E	403[A]	GOL	C1-C2-C3-O3
4	E	403[B]	GOL	C1-C2-C3-O3
4	E	404	GOL	O1-C1-C2-C3
4	C	404	GOL	C1-C2-C3-O3
4	C	401	GOL	O1-C1-C2-O2
4	C	404	GOL	O2-C2-C3-O3
4	E	403[A]	GOL	O2-C2-C3-O3
4	E	403[B]	GOL	O2-C2-C3-O3
4	E	404	GOL	O2-C2-C3-O3
4	A	403	GOL	O2-C2-C3-O3
4	E	404	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	402	GOL	2	0
4	C	404	GOL	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	60:PRO	C	61[A]:ARG	N	1.16

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	342/356 (96%)	-0.06	8 (2%) 60 67	14, 31, 68, 113	0
1	B	341/356 (95%)	0.01	13 (3%) 40 47	17, 34, 81, 109	0
1	C	341/356 (95%)	-0.08	9 (2%) 56 63	14, 30, 71, 107	0
1	D	336/356 (94%)	-0.05	10 (2%) 50 57	17, 32, 76, 112	0
1	E	328/356 (92%)	0.48	12 (3%) 41 48	37, 57, 95, 131	0
All	All	1688/1780 (94%)	0.06	52 (3%) 49 56	14, 36, 82, 131	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61[A]	ARG	7.1
1	D	147	ILE	6.0
1	E	73[A]	ARG	5.7
1	E	1	MET	5.7
1	D	345	GLY	5.4
1	B	347	ALA	4.7
1	B	349[A]	ARG	4.7
1	B	147	ILE	4.1
1	E	344	ALA	4.1
1	C	39[A]	LEU	4.0
1	A	62	MET	3.9
1	E	136	PRO	3.7
1	B	136	PRO	3.6
1	C	62	MET	3.5
1	B	75[A]	ARG	3.4
1	B	148	ILE	3.2
1	E	2	SER	3.2
1	E	39	LEU	3.1
1	B	149[A]	GLN	3.0
1	C	350	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	61[A]	ARG	3.0
1	A	57	VAL	3.0
1	A	85[A]	LEU	2.9
1	D	85	LEU	2.9
1	B	62	MET	2.9
1	D	0	SER	2.8
1	E	38	PRO	2.7
1	E	274[A]	ARG	2.8
1	D	148	ILE	2.6
1	E	85	LEU	2.6
1	C	61[A]	ARG	2.6
1	A	38	PRO	2.6
1	A	350	ILE	2.6
1	E	88[A]	LYS	2.5
1	D	149[A]	GLN	2.5
1	A	1	MET	2.5
1	B	57	VAL	2.5
1	C	87	GLY	2.4
1	B	350	ILE	2.3
1	C	112	ILE	2.3
1	B	74	LEU	2.3
1	C	109[A]	GLN	2.2
1	D	62[A]	MET	2.2
1	D	150	HIS	2.2
1	C	73	ARG	2.2
1	E	70	LYS	2.1
1	E	113[A]	ARG	2.1
1	B	85	LEU	2.1
1	B	61	ARG	2.1
1	D	74	LEU	2.1
1	C	1	MET	2.1
1	A	113[A]	ARG	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides 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
4	GOL	E	403[A]	6/6	0.59	0.44	104,110,111,111	6
4	GOL	E	403[B]	6/6	0.59	0.44	110,111,112,112	6
3	NA	B	401	1/1	0.81	0.08	55,55,55,55	0
4	GOL	D	402	6/6	0.85	0.16	82,83,84,86	0
4	GOL	A	403	6/6	0.87	0.31	76,78,81,82	0
4	GOL	E	404	6/6	0.91	0.23	61,63,68,70	0
3	NA	C	402	1/1	0.92	0.14	42,42,42,42	0
2	ACT	A	401	4/4	0.92	0.11	51,60,61,62	0
4	GOL	A	405	6/6	0.92	0.12	65,69,72,73	0
4	GOL	C	404	6/6	0.92	0.17	60,62,63,67	0
3	NA	E	401	1/1	0.93	0.09	46,46,46,46	0
3	NA	A	404	1/1	0.94	0.12	57,57,57,57	0
5	SO4	E	402	5/5	0.94	0.23	102,106,107,108	0
3	NA	D	401	1/1	0.95	0.05	44,44,44,44	0
3	NA	B	402	1/1	0.95	0.05	54,54,54,54	0
4	GOL	C	401	6/6	0.95	0.15	50,52,56,57	0
3	NA	C	403	1/1	0.96	0.06	34,34,34,34	0
3	NA	A	402	1/1	0.96	0.11	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.