



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 07:52 PM EDT

PDB ID : 3GB4 / pdb_00003gb4
Title : Crystal Structure of Dicamba Monooxygenase with Non-heme Cobalt and Dicamba
Authors : Rydel, T.J.; Sturman, E.J.; Moshiri, F.; Brown, G.R.; Qi, Y.
Deposited on : 2009-02-18
Resolution : 2.05 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

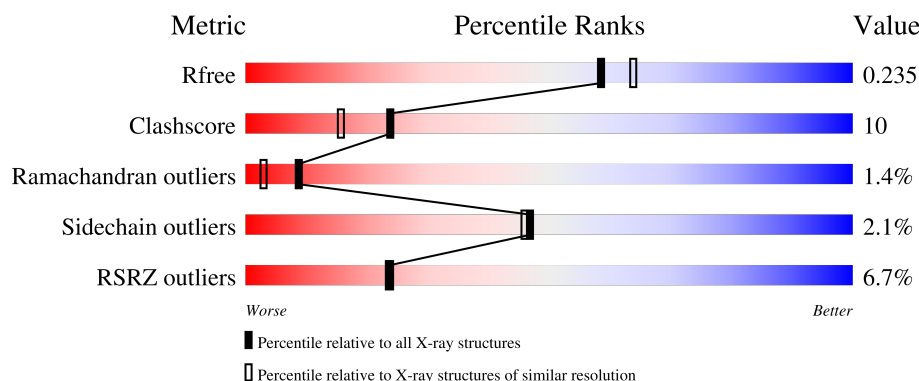
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	349	4% 76% 19% • •
1	B	349	6% 71% 25% • •
1	C	349	9% 72% 21% • •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8489 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 DdmC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			
1	B	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			
1	C	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			

There are 33 discrepancies between the modelled and reference sequences:

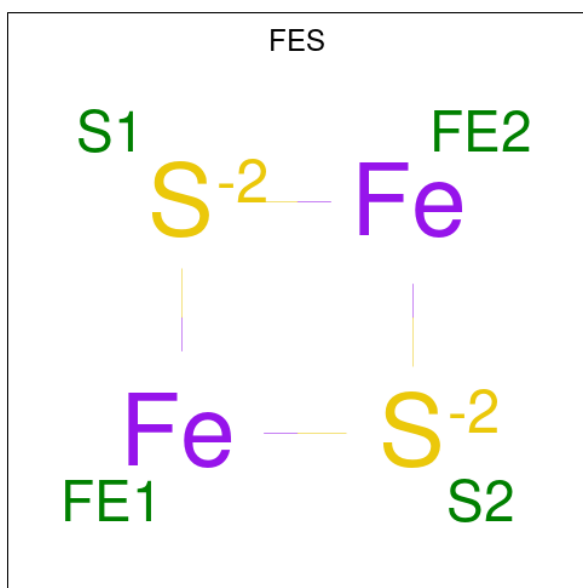
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q5S3I3
A	2	ALA	MET	engineered mutation	UNP Q5S3I3
A	341	ARG	-	expression tag	UNP Q5S3I3
A	342	LEU	-	expression tag	UNP Q5S3I3
A	343	GLU	-	expression tag	UNP Q5S3I3
A	344	HIS	-	expression tag	UNP Q5S3I3
A	345	HIS	-	expression tag	UNP Q5S3I3
A	346	HIS	-	expression tag	UNP Q5S3I3
A	347	HIS	-	expression tag	UNP Q5S3I3
A	348	HIS	-	expression tag	UNP Q5S3I3
A	349	HIS	-	expression tag	UNP Q5S3I3
B	1	MET	-	expression tag	UNP Q5S3I3
B	2	ALA	MET	engineered mutation	UNP Q5S3I3
B	341	ARG	-	expression tag	UNP Q5S3I3
B	342	LEU	-	expression tag	UNP Q5S3I3
B	343	GLU	-	expression tag	UNP Q5S3I3
B	344	HIS	-	expression tag	UNP Q5S3I3
B	345	HIS	-	expression tag	UNP Q5S3I3
B	346	HIS	-	expression tag	UNP Q5S3I3
B	347	HIS	-	expression tag	UNP Q5S3I3
B	348	HIS	-	expression tag	UNP Q5S3I3
B	349	HIS	-	expression tag	UNP Q5S3I3
C	1	MET	-	expression tag	UNP Q5S3I3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	2	ALA	MET	engineered mutation	UNP Q5S3I3
C	341	ARG	-	expression tag	UNP Q5S3I3
C	342	LEU	-	expression tag	UNP Q5S3I3
C	343	GLU	-	expression tag	UNP Q5S3I3
C	344	HIS	-	expression tag	UNP Q5S3I3
C	345	HIS	-	expression tag	UNP Q5S3I3
C	346	HIS	-	expression tag	UNP Q5S3I3
C	347	HIS	-	expression tag	UNP Q5S3I3
C	348	HIS	-	expression tag	UNP Q5S3I3
C	349	HIS	-	expression tag	UNP Q5S3I3

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe S 4 2 2	0	0
2	B	1	Total Fe S 4 2 2	0	0
2	C	1	Total Fe S 4 2 2	0	0

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

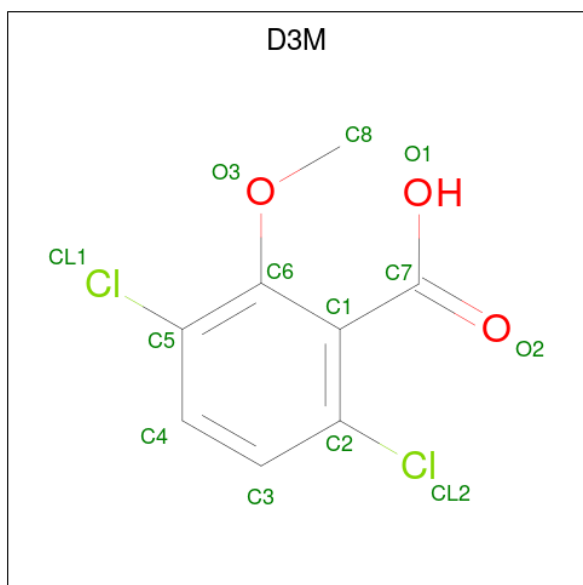
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Co 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Co	0	0
			3	3		
3	C	2	Total	Co	0	0
			2	2		

- Molecule 4 is 3,6-dichloro-2-methoxybenzoic acid (CCD ID: D3M) (formula: $C_8H_6Cl_2O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	Cl	O	0	0
			13	8	2	3		
4	B	1	Total	C	Cl	O	0	0
			13	8	2	3		

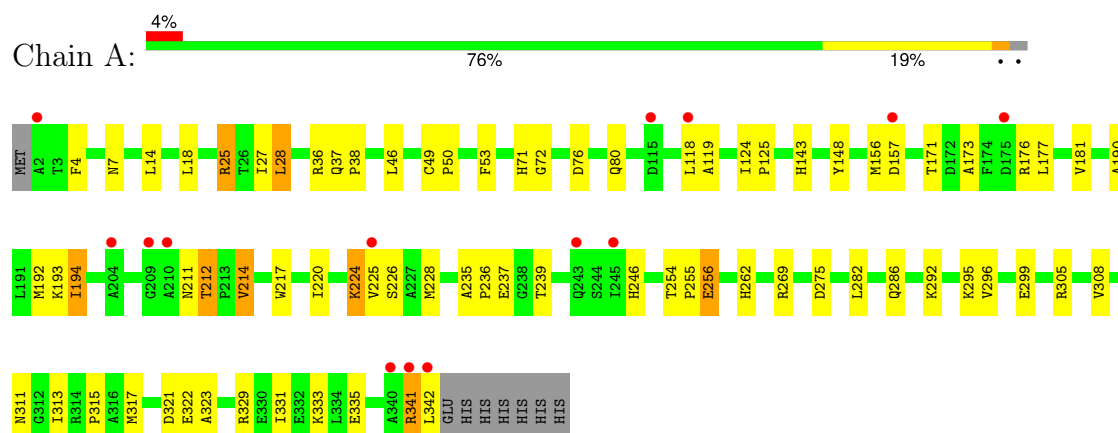
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	179	Total	O	0	0
			179	179		
5	B	174	Total	O	0	0
			174	174		
5	C	168	Total	O	0	0
			168	168		

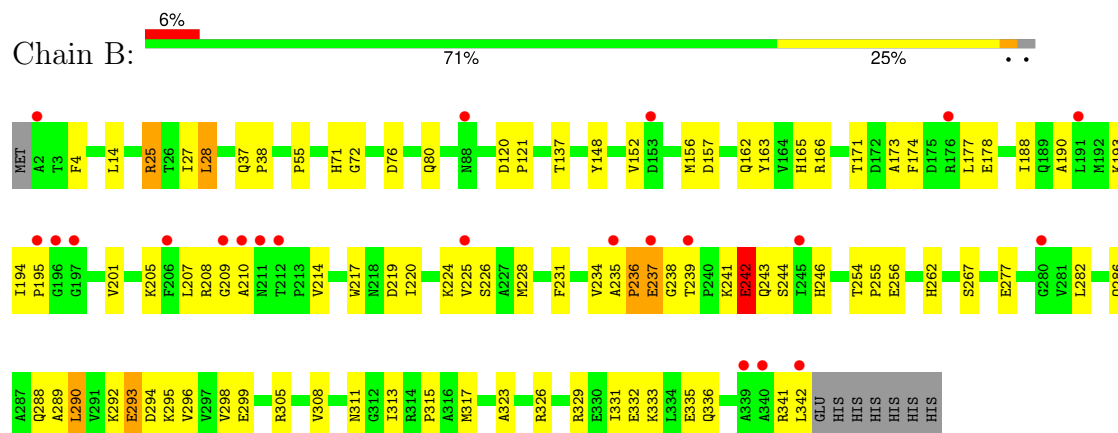
3 Residue-property plots [i](#)

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.

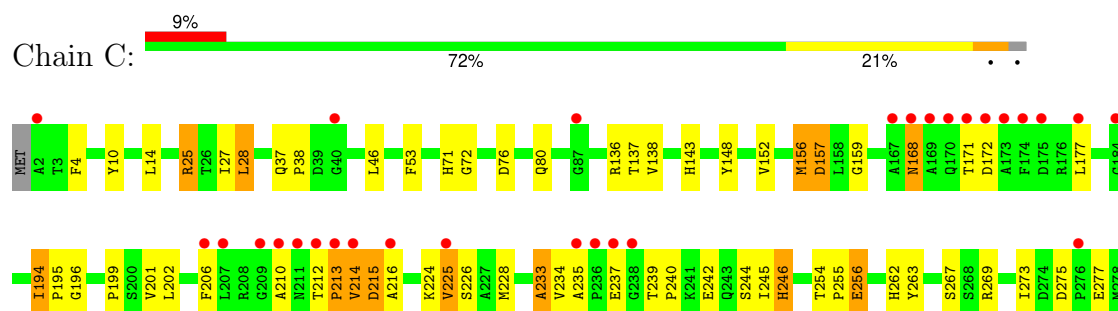
• Molecule 1: DdmC



• Molecule 1: DdmC



• Molecule 1: DdmC





4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	81.55Å 81.55Å 161.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.05 20.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	82.3 (20.00-2.05) 82.3 (20.00-2.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.05Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.221 , 0.262 0.200 , 0.235	Depositor DCC
R_{free} test set	6177 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.006 for -h,-k,l 0.039 for h,-h-k,-l 0.021 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8489	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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: FES, D3M, CO

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	0.39	0/2706	0.93	14/3685 (0.4%)
1	B	0.38	0/2706	0.92	15/3685 (0.4%)
1	C	0.40	1/2706 (0.0%)	0.93	16/3685 (0.4%)
All	All	0.39	1/8118 (0.0%)	0.93	45/11055 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	156	MET	SD-CE	-5.21	1.66	1.79

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	291	VAL	N-CA-C	7.49	117.61	110.42
1	B	226	SER	N-CA-C	7.30	119.02	110.41
1	A	226	SER	N-CA-C	7.24	118.95	110.41
1	A	72	GLY	N-CA-C	7.14	125.22	115.30
1	C	215	ASP	N-CA-C	7.01	120.70	109.83
1	B	72	GLY	N-CA-C	7.00	125.02	115.30
1	C	72	GLY	N-CA-C	6.77	124.72	115.30
1	A	256	GLU	N-CA-C	-6.52	103.20	112.13
1	B	14	LEU	N-CA-C	-6.42	102.05	110.39
1	B	246	HIS	N-CA-C	6.41	118.52	108.96
1	B	256	GLU	N-CA-C	-6.25	103.56	112.13
1	C	246	HIS	N-CA-C	6.24	118.26	108.96
1	A	14	LEU	N-CA-C	-6.24	102.28	110.39
1	A	311	ASN	N-CA-C	-6.18	100.84	110.17
1	A	246	HIS	N-CA-C	6.17	118.16	108.96
1	C	157	ASP	N-CA-C	6.16	118.95	108.90
1	C	233	ALA	N-CA-C	6.12	118.09	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	152	VAL	N-CA-C	-6.09	104.80	110.53
1	B	311	ASN	N-CA-C	-6.08	100.98	110.17
1	C	256	GLU	N-CA-C	-6.08	103.38	111.96
1	B	293	GLU	N-CA-C	6.08	117.59	110.97
1	B	244	SER	N-CA-C	6.07	116.25	108.24
1	C	311	ASN	N-CA-C	-5.97	101.63	110.46
1	B	194	ILE	CA-C-N	5.96	126.86	120.13
1	B	194	ILE	C-N-CA	5.96	126.86	120.13
1	B	341	ARG	N-CA-C	5.94	117.57	110.44
1	C	152	VAL	N-CA-C	-5.94	104.95	110.53
1	A	53	PHE	N-CA-C	5.88	120.40	113.23
1	C	14	LEU	N-CA-C	-5.77	102.35	110.31
1	C	226	SER	N-CA-C	5.77	118.77	109.83
1	C	46	LEU	N-CA-C	-5.66	101.47	109.96
1	C	159	GLY	N-CA-C	5.64	120.98	114.16
1	C	313	ILE	N-CA-C	5.63	116.33	108.89
1	B	290	LEU	N-CA-C	5.49	117.12	111.03
1	A	275	ASP	CA-C-N	5.44	125.26	119.28
1	A	275	ASP	C-N-CA	5.44	125.26	119.28
1	C	53	PHE	N-CA-C	5.41	119.83	113.23
1	A	194	ILE	CA-C-N	5.34	125.80	120.14
1	A	194	ILE	C-N-CA	5.34	125.80	120.14
1	B	313	ILE	N-CA-C	5.29	115.87	108.89
1	A	313	ILE	N-CA-C	5.24	115.81	108.89
1	A	46	LEU	N-CA-C	-5.23	102.31	110.10
1	B	195	PRO	N-CA-C	5.19	119.00	110.55
1	C	321	ASP	N-CA-C	5.13	119.52	113.16
1	A	321	ASP	N-CA-C	5.10	119.48	113.16

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	2641	0	2577	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2641	0	2577	51	0
1	C	2641	0	2577	61	0
2	A	4	0	0	1	0
2	B	4	0	0	1	0
2	C	4	0	0	1	0
3	A	2	0	0	0	0
3	B	3	0	0	0	0
3	C	2	0	0	0	0
4	A	13	0	5	0	0
4	B	13	0	5	0	0
5	A	179	0	0	5	0
5	B	174	0	0	1	0
5	C	168	0	0	4	0
All	All	8489	0	7741	158	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 10.

All (158) 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:156:MET:HE3	1:C:228:MET:SD	1.97	1.04
1:B:156:MET:HE3	1:B:228:MET:SD	2.15	0.86
1:A:156:MET:HE3	1:A:228:MET:SD	2.19	0.82
1:A:171:THR:HG22	1:A:173:ALA:H	1.46	0.80
1:B:288:GLN:HA	1:B:292:LYS:HE3	1.67	0.76
1:C:206:PHE:HB2	1:C:245:ILE:HD13	1.69	0.74
1:C:199:PRO:HB3	1:C:234:VAL:HG21	1.72	0.72
1:B:178:GLU:HG3	1:B:193:LYS:HB3	1.74	0.70
1:A:7:ASN:ND2	1:A:341:ARG:HH12	1.92	0.67
1:C:201:VAL:HG13	1:C:202:LEU:HD22	1.75	0.66
1:C:233:ALA:HB2	1:C:246:HIS:HB3	1.75	0.66
1:B:235:ALA:HB1	1:B:239:THR:HB	1.79	0.64
1:A:156:MET:HE1	1:A:220:ILE:HG12	1.80	0.62
1:B:156:MET:HE1	1:B:220:ILE:HG12	1.82	0.61
1:B:55:PRO:HB3	1:C:304:ARG:CZ	2.33	0.59
1:C:4:PHE:HB3	1:C:28:LEU:HD13	1.84	0.59
1:A:4:PHE:HB3	1:A:28:LEU:HD13	1.85	0.58
1:C:194:ILE:HG23	1:C:216:ALA:HB3	1.86	0.58
1:B:282:LEU:O	1:B:286:GLN:HG3	2.04	0.58
1:B:4:PHE:HB3	1:B:28:LEU:HD13	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:LYS:HB2	5:A:1104:HOH:O	2.06	0.56
1:C:275:ASP:OD1	1:C:277:GLU:HG2	2.06	0.56
1:B:162:GLN:HG3	1:B:174:PHE:CD1	2.40	0.56
1:C:213:PRO:O	1:C:214:VAL:HB	2.06	0.56
1:B:171:THR:HG22	1:B:173:ALA:H	1.71	0.55
1:B:237:GLU:HG3	1:B:238:GLY:H	1.71	0.55
1:C:240:PRO:HB2	1:C:242:GLU:OE1	2.07	0.55
1:C:76:ASP:OD2	1:C:80:GLN:HB3	2.06	0.55
1:A:171:THR:HG21	1:A:194:ILE:HG21	1.90	0.54
1:A:173:ALA:HB2	1:A:176:ARG:HH21	1.71	0.54
1:B:27:ILE:HD12	1:B:254:THR:HG21	1.90	0.54
1:A:7:ASN:HD21	1:A:341:ARG:HH12	1.56	0.53
1:A:171:THR:HG22	1:A:173:ALA:N	2.21	0.53
1:A:76:ASP:OD2	1:A:80:GLN:HB3	2.08	0.53
1:C:215:ASP:HB2	1:C:235:ALA:O	2.09	0.53
1:C:214:VAL:HG13	1:C:237:GLU:O	2.09	0.53
1:C:137:THR:HA	1:C:267:SER:O	2.08	0.53
1:C:157:ASP:HB2	5:C:823:HOH:O	2.09	0.53
1:C:235:ALA:HB1	1:C:239:THR:OG1	2.09	0.52
1:A:156:MET:HA	1:A:156:MET:HE2	1.91	0.52
1:B:162:GLN:HE22	1:B:171:THR:H	1.57	0.52
1:A:214:VAL:HG22	1:A:235:ALA:O	2.09	0.52
1:A:236:PRO:HG2	1:A:239:THR:HG21	1.90	0.52
1:B:207:LEU:HD11	1:B:234:VAL:HG22	1.90	0.52
1:C:37:GLN:HB3	1:C:38:PRO:HD2	1.91	0.52
1:B:157:ASP:O	1:B:157:ASP:OD1	2.27	0.51
1:A:27:ILE:HD12	1:A:254:THR:HG21	1.92	0.51
1:B:217:TRP:CD1	1:B:241:LYS:HD3	2.45	0.51
1:C:143:HIS:HE1	1:C:256:GLU:OE1	1.94	0.51
1:C:196:GLY:N	1:C:213:PRO:HA	2.26	0.51
1:C:71:HIS:HB2	2:C:501:FES:S1	2.51	0.51
1:C:242:GLU:H	1:C:242:GLU:CD	2.19	0.50
1:C:156:MET:HE2	1:C:156:MET:HA	1.92	0.50
1:A:211:ASN:O	1:A:212:THR:HB	2.11	0.50
1:C:214:VAL:HG22	1:C:214:VAL:O	2.11	0.50
1:B:76:ASP:OD2	1:B:80:GLN:HB3	2.12	0.50
1:C:201:VAL:CG1	1:C:202:LEU:HD22	2.43	0.49
5:B:704:HOH:O	1:C:320:CYS:HB3	2.12	0.49
1:C:212:THR:HB	1:C:237:GLU:HG3	1.95	0.49
1:C:10:TYR:CE2	1:C:225:VAL:HG11	2.47	0.49
1:C:136:ARG:NH2	1:C:279:ASP:OD1	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:331:ILE:O	1:C:335:GLU:HG3	2.12	0.49
1:C:206:PHE:HB2	1:C:245:ILE:CD1	2.40	0.49
1:B:295:LYS:HE3	1:B:299:GLU:OE2	2.13	0.49
1:B:201:VAL:O	1:B:205:LYS:HG2	2.12	0.48
1:A:71:HIS:HB2	2:A:501:FES:S1	2.54	0.48
1:B:156:MET:HA	1:B:156:MET:HE2	1.95	0.48
1:B:209:GLY:O	1:B:210:ALA:HB3	2.14	0.48
1:B:37:GLN:HB3	1:B:38:PRO:HD2	1.95	0.47
1:A:157:ASP:O	1:A:157:ASP:OD1	2.31	0.47
1:C:315:PRO:HB2	1:C:317:MET:HE3	1.95	0.47
1:B:242:GLU:CD	1:B:243:GLN:H	2.21	0.47
1:A:181:VAL:HG21	1:A:322:GLU:HG2	1.96	0.47
1:B:188:ILE:HD13	1:B:326:ARG:HB3	1.96	0.47
1:B:331:ILE:O	1:B:335:GLU:HG3	2.15	0.47
1:C:214:VAL:HG13	1:C:214:VAL:O	2.15	0.47
1:B:295:LYS:O	1:B:299:GLU:HG3	2.15	0.47
1:C:206:PHE:O	1:C:245:ILE:HD11	2.15	0.46
1:A:37:GLN:HB3	1:A:38:PRO:HD2	1.98	0.46
1:A:315:PRO:HB2	1:A:317:MET:HE3	1.98	0.46
1:B:162:GLN:HG3	1:B:174:PHE:CE1	2.51	0.46
1:A:329:ARG:O	1:A:333:LYS:HG3	2.15	0.46
1:A:157:ASP:HB3	5:A:869:HOH:O	2.15	0.46
1:B:165:HIS:ND1	1:B:293:GLU:OE2	2.44	0.46
1:C:234:VAL:O	1:C:244:SER:HB2	2.16	0.46
1:B:71:HIS:HB2	2:B:501:FES:S1	2.56	0.45
1:B:207:LEU:O	1:B:208:ARG:HB2	2.14	0.45
1:C:292:LYS:C	1:C:293:GLU:HG3	2.41	0.45
1:B:55:PRO:HB3	1:C:304:ARG:NH2	2.31	0.45
1:A:118:LEU:O	1:A:119:ALA:C	2.59	0.45
1:A:224:LYS:HB3	5:A:1124:HOH:O	2.17	0.45
1:B:190:ALA:HB2	1:B:323:ALA:CB	2.45	0.45
1:A:25:ARG:HD2	1:A:262:HIS:CE1	2.51	0.45
1:A:282:LEU:O	1:A:286:GLN:HG3	2.17	0.45
1:A:173:ALA:CB	1:A:176:ARG:HH21	2.30	0.44
1:B:292:LYS:O	1:B:296:VAL:HG23	2.18	0.44
1:C:25:ARG:HD2	1:C:262:HIS:CE1	2.52	0.44
1:C:305:ARG:O	1:C:308:VAL:HG22	2.17	0.44
1:A:143:HIS:HE1	1:A:256:GLU:OE1	2.00	0.44
1:B:207:LEU:HD13	1:B:210:ALA:HB2	1.99	0.44
1:B:294:ASP:O	1:B:298:VAL:HG22	2.18	0.44
1:C:288:GLN:NE2	1:C:292:LYS:HB2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:LYS:HE3	1:A:299:GLU:OE2	2.18	0.44
1:B:305:ARG:O	1:B:308:VAL:HG22	2.17	0.44
1:A:193:LYS:HD2	1:A:217:TRP:HE1	1.83	0.44
1:C:225:VAL:HB	5:C:845:HOH:O	2.18	0.44
1:C:295:LYS:O	1:C:299:GLU:HG3	2.18	0.44
1:A:269:ARG:HD3	1:A:282:LEU:HD11	2.00	0.44
1:C:235:ALA:HB1	1:C:239:THR:HG1	1.82	0.44
1:B:28:LEU:N	1:B:28:LEU:HD23	2.33	0.43
1:B:315:PRO:HB2	1:B:317:MET:HE3	1.99	0.43
1:C:295:LYS:HE3	1:C:299:GLU:OE2	2.17	0.43
1:B:137:THR:HA	1:B:267:SER:O	2.19	0.43
1:C:233:ALA:CB	1:C:246:HIS:HB3	2.43	0.43
1:B:305:ARG:HA	1:B:308:VAL:HG22	2.01	0.43
1:C:27:ILE:HD12	1:C:254:THR:HG21	2.00	0.43
1:B:120:ASP:HA	1:B:121:PRO:HD2	1.91	0.43
1:B:163:TYR:O	1:B:166:ARG:HB2	2.19	0.43
1:B:148:TYR:CD2	1:B:255:PRO:HG3	2.54	0.43
1:C:171:THR:HB	1:C:172:ASP:H	1.73	0.42
1:C:194:ILE:O	1:C:216:ALA:HB3	2.19	0.42
1:C:332:GLU:O	1:C:336:GLN:HG3	2.20	0.42
1:C:148:TYR:CD2	1:C:255:PRO:HG3	2.55	0.42
1:A:148:TYR:CD2	1:A:255:PRO:HG3	2.55	0.42
1:B:162:GLN:NE2	1:B:171:THR:H	2.16	0.42
1:A:177:LEU:HD21	1:A:192:MET:HE3	2.02	0.42
1:A:190:ALA:HB2	1:A:323:ALA:CB	2.50	0.42
1:A:295:LYS:O	1:A:299:GLU:HG3	2.19	0.42
1:A:342:LEU:H	1:A:342:LEU:HD23	1.84	0.42
1:B:236:PRO:O	1:B:237:GLU:C	2.61	0.42
1:A:157:ASP:CG	5:A:1080:HOH:O	2.63	0.42
1:B:289:ALA:HA	1:B:293:GLU:OE2	2.20	0.42
1:A:235:ALA:HB1	1:A:236:PRO:HD2	2.02	0.42
1:A:305:ARG:O	1:A:308:VAL:HG22	2.19	0.42
1:B:332:GLU:O	1:B:336:GLN:HG3	2.20	0.42
1:C:233:ALA:HB2	1:C:246:HIS:CB	2.48	0.42
1:A:341:ARG:HG2	1:A:341:ARG:HH11	1.84	0.41
1:B:25:ARG:HD2	1:B:262:HIS:CE1	2.54	0.41
1:A:331:ILE:O	1:A:335:GLU:HG3	2.20	0.41
1:C:263:TYR:CD1	1:C:290:LEU:HD22	2.55	0.41
1:B:235:ALA:CB	1:B:239:THR:HB	2.50	0.41
1:B:329:ARG:O	1:B:333:LYS:HG3	2.19	0.41
1:C:339:ALA:C	1:C:341:ARG:H	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:269:ARG:HD3	1:C:282:LEU:CD1	2.51	0.41
1:C:168:ASN:HD22	1:C:168:ASN:HA	1.54	0.41
1:C:194:ILE:HA	1:C:195:PRO:HD2	1.86	0.41
1:A:49:CYS:HA	1:A:50:PRO:HD3	1.90	0.41
1:B:219:ASP:HB2	1:B:231:PHE:HB3	2.02	0.41
1:C:28:LEU:N	1:C:28:LEU:HD23	2.36	0.41
1:C:136:ARG:HE	1:C:279:ASP:CG	2.29	0.41
1:C:293:GLU:HG2	5:C:857:HOH:O	2.19	0.41
1:A:292:LYS:O	1:A:296:VAL:HG23	2.21	0.41
1:C:138:VAL:HG11	1:C:279:ASP:HB3	2.03	0.41
1:A:124:ILE:HA	1:A:125:PRO:HD3	1.94	0.40
1:C:273:ILE:HD11	5:C:918:HOH:O	2.22	0.40
1:A:18:LEU:HB2	1:A:36:ARG:HH12	1.87	0.40
1:A:305:ARG:HA	1:A:308:VAL:HG22	2.03	0.40
1:A:341:ARG:NH2	5:A:1215:HOH:O	2.48	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	339/349 (97%)	317 (94%)	18 (5%)	4 (1%)	10	4
1	B	339/349 (97%)	316 (93%)	18 (5%)	5 (2%)	8	2
1	C	339/349 (97%)	311 (92%)	23 (7%)	5 (2%)	8	2
All	All	1017/1047 (97%)	944 (93%)	59 (6%)	14 (1%)	9	3

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	237	GLU
1	C	210	ALA

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Mol	Chain	Res	Type
1	C	214	VAL
1	A	225	VAL
1	A	237	GLU
1	B	225	VAL
1	B	242	GLU
1	C	213	PRO
1	C	225	VAL
1	A	224	LYS
1	B	224	LYS
1	C	224	LYS
1	B	236	PRO
1	A	212	THR

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	276/284 (97%)	272 (99%)	4 (1%)	59	60
1	B	276/284 (97%)	268 (97%)	8 (3%)	37	33
1	C	276/284 (97%)	271 (98%)	5 (2%)	51	51
All	All	828/852 (97%)	811 (98%)	17 (2%)	47	46

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	28	LEU
1	A	214	VAL
1	A	341	ARG
1	B	25	ARG
1	B	28	LEU
1	B	177	LEU
1	B	214	VAL
1	B	242	GLU
1	B	277	GLU

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Mol	Chain	Res	Type
1	B	290	LEU
1	B	342	LEU
1	C	25	ARG
1	C	28	LEU
1	C	168	ASN
1	C	177	LEU
1	C	194	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	143	HIS
1	A	154	ASN
1	A	211	ASN
1	A	218	ASN
1	A	288	GLN
1	B	80	GLN
1	B	143	HIS
1	B	170	GLN
1	B	336	GLN
1	C	143	HIS
1	C	154	ASN
1	C	168	ASN
1	C	211	ASN
1	C	223	ASN
1	C	336	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

Of 12 ligands modelled in this entry, 7 are monoatomic - leaving 5 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	D3M	B	601	-	13,13,13	1.58	2 (15%)	18,18,18	1.14	2 (11%)
2	FES	B	501	1	0,4,4	-	-	-	-	-
4	D3M	A	601	-	13,13,13	1.50	2 (15%)	18,18,18	1.14	1 (5%)
2	FES	A	501	1	0,4,4	-	-	-	-	-
2	FES	C	501	1	0,4,4	-	-	-	-	-

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	D3M	B	601	-	-	0/6/6/6	0/1/1/1
2	FES	B	501	1	-	-	0/1/1/1
4	D3M	A	601	-	-	0/6/6/6	0/1/1/1
2	FES	A	501	1	-	-	0/1/1/1
2	FES	C	501	1	-	-	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	D3M	C1-C2	3.12	1.44	1.39
4	A	601	D3M	C1-C2	2.78	1.43	1.39
4	B	601	D3M	C1-C6	2.07	1.45	1.40
4	A	601	D3M	C4-C3	2.01	1.42	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	D3M	C6-C5-CL1	2.83	122.14	118.40

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	601	D3M	C6-C5-CL1	2.70	121.96	118.40
4	B	601	D3M	C1-C2-CL2	2.07	122.29	119.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	FES	1	0
2	A	501	FES	1	0
2	C	501	FES	1	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	341/349 (97%)	0.33	14 (4%) 41 41	27, 42, 65, 84	0
1	B	341/349 (97%)	0.53	22 (6%) 25 24	26, 43, 70, 90	0
1	C	341/349 (97%)	0.55	33 (9%) 13 13	27, 40, 83, 90	0
All	All	1023/1047 (97%)	0.47	69 (6%) 24 23	26, 42, 73, 90	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	213	PRO	8.4
1	C	212	THR	8.0
1	C	173	ALA	7.7
1	C	167	ALA	6.1
1	C	174	PHE	6.0
1	C	177	LEU	5.7
1	C	172	ASP	5.3
1	B	191	LEU	4.5
1	C	171	THR	4.3
1	C	169	ALA	4.3
1	C	342	LEU	4.2
1	A	342	LEU	3.9
1	B	212	THR	3.9
1	A	245	ILE	3.8
1	C	2	ALA	3.7
1	C	209	GLY	3.6
1	B	239	THR	3.4
1	B	342	LEU	3.4
1	A	225	VAL	3.4
1	A	2	ALA	3.4
1	C	214	VAL	3.4
1	B	211	ASN	3.3
1	C	175	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	237	GLU	3.2
1	C	312	GLY	3.2
1	C	170	GLN	3.2
1	B	210	ALA	3.1
1	B	225	VAL	3.1
1	C	207	LEU	2.9
1	A	157	ASP	2.9
1	A	175	ASP	2.8
1	C	276	PRO	2.8
1	A	118	LEU	2.8
1	C	235	ALA	2.7
1	C	236	PRO	2.7
1	B	2	ALA	2.7
1	B	340	ALA	2.6
1	C	168	ASN	2.6
1	C	206	PHE	2.6
1	B	209	GLY	2.6
1	B	206	PHE	2.5
1	A	340	ALA	2.5
1	C	225	VAL	2.5
1	B	153	ASP	2.4
1	C	210	ALA	2.4
1	C	216	ALA	2.4
1	C	87	GLY	2.4
1	B	197	GLY	2.3
1	A	115	ASP	2.3
1	C	341	ARG	2.3
1	B	235	ALA	2.3
1	B	196	GLY	2.3
1	C	339	ALA	2.3
1	B	176	ARG	2.3
1	B	245	ILE	2.3
1	B	339	ALA	2.3
1	B	237	GLU	2.2
1	B	88	ASN	2.2
1	C	211	ASN	2.2
1	A	243	GLN	2.1
1	C	238	GLY	2.1
1	B	280	GLY	2.1
1	C	184	GLY	2.1
1	A	204	ALA	2.1
1	A	210	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	341	ARG	2.0
1	B	195	PRO	2.0
1	A	209	GLY	2.0
1	C	40	GLY	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
4	D3M	B	601	13/13	0.77	0.13	51,52,55,55	0
4	D3M	A	601	13/13	0.85	0.10	49,51,52,54	0
3	CO	B	603	1/1	0.91	0.09	82,82,82,82	0
3	CO	B	602	1/1	0.95	0.09	62,62,62,62	0
3	CO	A	602	1/1	0.95	0.06	66,66,66,66	0
3	CO	C	602	1/1	0.96	0.09	56,56,56,56	0
3	CO	B	502	1/1	0.97	0.04	40,40,40,40	0
2	FES	A	501	4/4	0.98	0.04	32,32,34,34	0
2	FES	B	501	4/4	0.99	0.04	28,30,30,31	0
2	FES	C	501	4/4	0.99	0.03	28,29,29,29	0
3	CO	C	502	1/1	0.99	0.05	40,40,40,40	0
3	CO	A	502	1/1	1.00	0.07	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.