



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:49 AM GMT

PDB ID : 3JUS
Title : Crystal structure of human lanosterol 14alpha-demethylase (CYP51) in complex with econazole
Authors : Strushkevich, N.; MacKenzie, F.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Weigelt, J.; Usanov, S.A.; Park, H.; Structural Genomics Consortium (SGC)
Deposited on : 2009-09-15
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

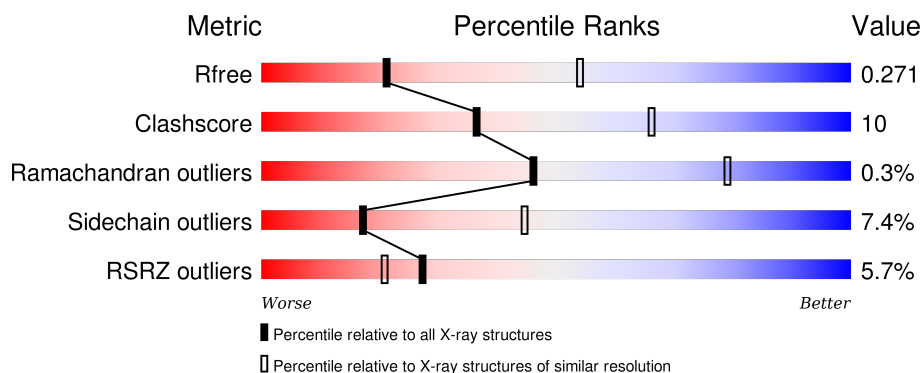
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	461	
1	B	461	

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
3	ECN	A	602[A]	-	-	X	-
3	ECN	B	602[A]	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7553 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 Lanosterol 14-alpha demethylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	0	0
			3589	2310	611	652	16			
1	B	445	Total	C	N	O	S	0	0	0
			3589	2310	611	652	16			

There are 24 discrepancies between the modelled and reference sequences:

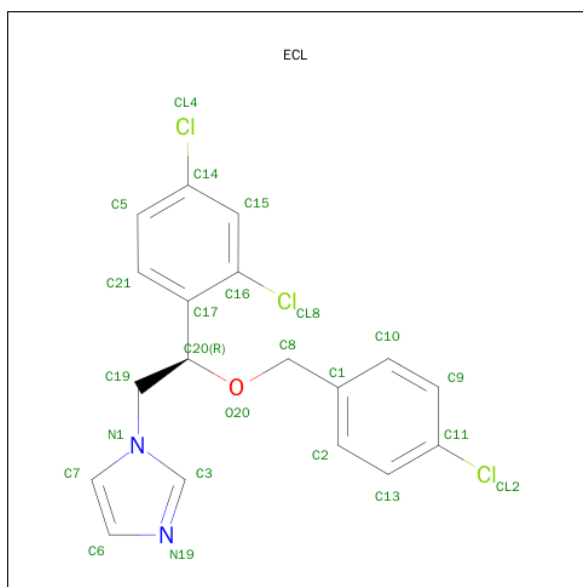
Chain	Residue	Modelled	Actual	Comment	Reference
A	49	MET	-	EXPRESSION TAG	UNP Q16850
A	50	ALA	-	EXPRESSION TAG	UNP Q16850
A	51	LYS	-	EXPRESSION TAG	UNP Q16850
A	52	LYS	-	EXPRESSION TAG	UNP Q16850
A	53	THR	-	EXPRESSION TAG	UNP Q16850
A	503	THR	-	EXPRESSION TAG	UNP Q16850
A	504	HIS	-	EXPRESSION TAG	UNP Q16850
A	505	HIS	-	EXPRESSION TAG	UNP Q16850
A	506	HIS	-	EXPRESSION TAG	UNP Q16850
A	507	HIS	-	EXPRESSION TAG	UNP Q16850
A	508	HIS	-	EXPRESSION TAG	UNP Q16850
A	509	HIS	-	EXPRESSION TAG	UNP Q16850
B	49	MET	-	EXPRESSION TAG	UNP Q16850
B	50	ALA	-	EXPRESSION TAG	UNP Q16850
B	51	LYS	-	EXPRESSION TAG	UNP Q16850
B	52	LYS	-	EXPRESSION TAG	UNP Q16850
B	53	THR	-	EXPRESSION TAG	UNP Q16850
B	503	THR	-	EXPRESSION TAG	UNP Q16850
B	504	HIS	-	EXPRESSION TAG	UNP Q16850
B	505	HIS	-	EXPRESSION TAG	UNP Q16850
B	506	HIS	-	EXPRESSION TAG	UNP Q16850
B	507	HIS	-	EXPRESSION TAG	UNP Q16850
B	508	HIS	-	EXPRESSION TAG	UNP Q16850
B	509	HIS	-	EXPRESSION TAG	UNP Q16850

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- Chemical structure of HEM (Heme) showing a central iron atom coordinated by four nitrogen atoms in a porphyrin-like ring, with various side chains and a central heme group.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

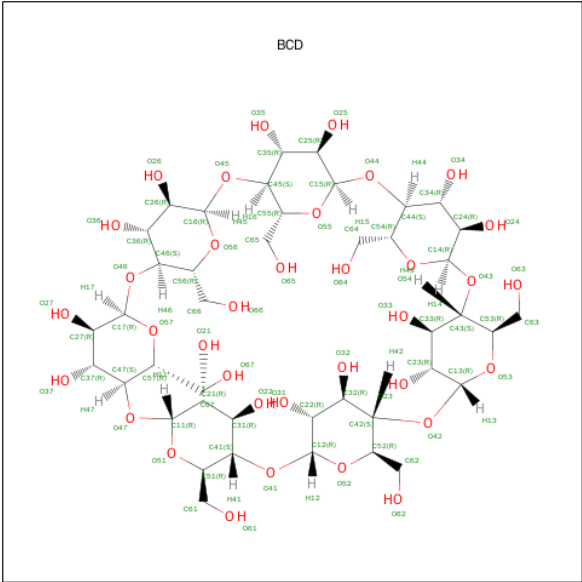
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	1
			24	18	3	2	1		
3	B	1	Total	C	Cl	N	O	0	1
			24	18	3	2	1		

- Molecule 4 is 1-[(2R)-2-[(4-CHLOROBENZYL)OXY]-2-(2,4-DICHLOROPHENYL)ETHYL]-1H-IMIDAZOLE (three-letter code: ECL) (formula: C₁₈H₁₅Cl₃N₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	0	1
			24	18	3	2	1		
4	B	1	Total	C	Cl	N	O	0	1
			24	18	3	2	1		

- Molecule 5 is SUGAR (BETA-CYCLODEXTRIN) (three-letter code: BCD) (formula: C₄₂H₇₀O₃₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			77	42	35		
5	B	1	Total	C	O	0	0
			77	42	35		

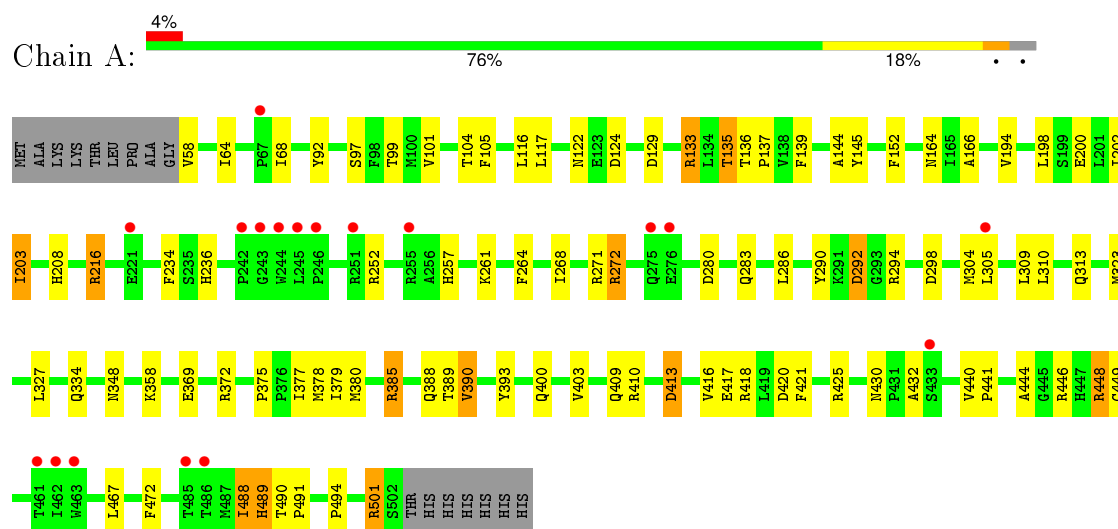
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	29	Total	O	0	0
			29	29		
6	B	10	Total	O	0	0
			10	10		

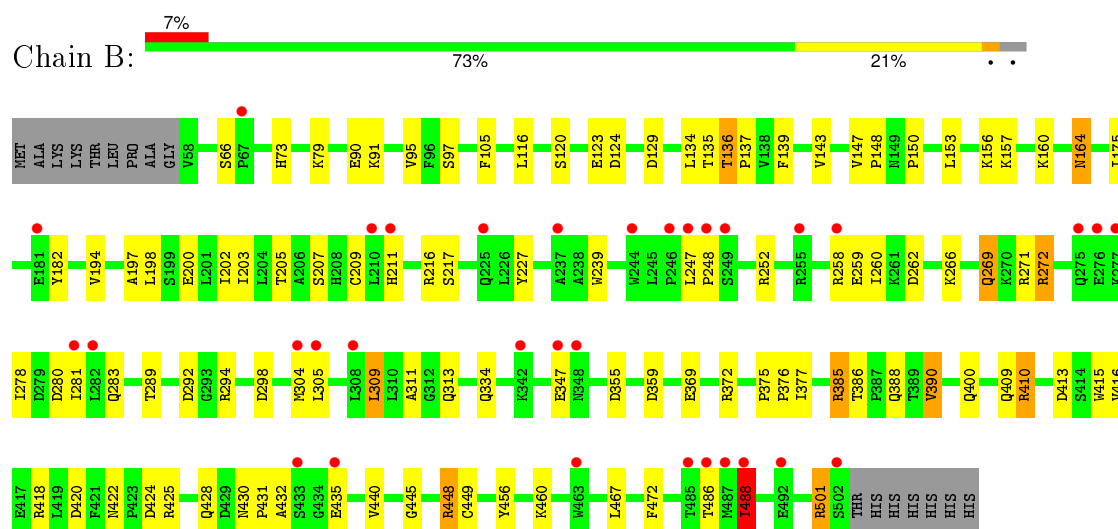
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Lanosterol 14-alpha demethylase



- Molecule 1: Lanosterol 14-alpha demethylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	146.63 Å 146.63 Å 110.56 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.52 – 2.90 24.50 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.2 (24.52-2.90) 99.2 (24.50-2.89)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 2.89 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.203 , 0.272 0.203 , 0.271	Depositor DCC
R_{free} test set	1356 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	66.4	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 27428 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7553	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: BCD, HEM, ECL, ECN

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	0.48	0/3682	0.59	0/4988
1	B	0.45	0/3682	0.58	0/4988
All	All	0.47	0/7364	0.59	0/9976

There are no bond length outliers.

There are no bond angle outliers.

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	3589	0	3581	66	0
1	B	3589	0	3581	58	0
2	A	43	0	30	7	0
2	B	43	0	30	6	0
3	A	24	0	15	8	0
3	B	24	0	15	7	0
4	A	24	0	15	0	0
4	B	24	0	15	1	0
5	A	77	0	70	1	0
5	B	77	0	70	0	0
6	A	29	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	10	0	0	2	0
All	All	7553	0	7422	142	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 (142) 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)
3:A:602[A]:ECN:H8C1	3:A:602[A]:ECN:C3	1.49	1.40
3:B:602[A]:ECN:C3	3:B:602[A]:ECN:H8C1	1.65	1.25
3:A:602[A]:ECN:C8	3:A:602[A]:ECN:C3	2.39	1.01
3:A:602[A]:ECN:H8C1	3:A:602[A]:ECN:N1	1.71	0.93
1:B:334:GLN:HE22	1:B:472:PHE:H	1.23	0.86
1:A:409:GLN:HE21	1:A:440:VAL:H	1.24	0.85
3:B:602[A]:ECN:H3	3:B:602[A]:ECN:H8C1	1.57	0.83
3:A:602[A]:ECN:H3	3:A:602[A]:ECN:H8C1	1.63	0.81
3:B:602[A]:ECN:C3	3:B:602[A]:ECN:C8	2.56	0.79
1:B:409:GLN:HE21	1:B:440:VAL:H	1.32	0.77
1:A:198:LEU:O	1:A:202:ILE:HG12	1.86	0.76
3:B:602[A]:ECN:H191	3:B:602[A]:ECN:CL4	2.24	0.74
1:B:124:ASP:HB3	1:B:388:GLN:NE2	2.03	0.73
1:B:200:GLU:O	1:B:203:ILE:HG22	1.89	0.73
1:A:124:ASP:HB3	1:A:388:GLN:HE21	1.53	0.71
1:A:416:VAL:O	1:A:425:ARG:NH2	2.24	0.70
1:A:372:ARG:NH2	1:A:418:ARG:O	2.20	0.69
1:B:309:LEU:O	1:B:313:GLN:HB2	1.93	0.69
3:B:602[A]:ECN:N1	3:B:602[A]:ECN:H8C1	2.05	0.66
1:B:271:ARG:HD3	1:B:283:GLN:OE1	1.96	0.66
2:B:601:HEM:HMB2	2:B:601:HEM:HBB2	1.79	0.64
1:A:430:ASN:ND2	1:A:432:ALA:HB3	2.12	0.64
1:B:372:ARG:NH2	1:B:418:ARG:O	2.27	0.64
1:A:334:GLN:HE22	1:A:472:PHE:H	1.46	0.64
1:A:467:LEU:O	1:A:501:ARG:HD2	1.98	0.63
3:B:602[A]:ECN:H3	3:B:602[A]:ECN:C8	2.22	0.63
1:B:116:LEU:HB2	1:B:390:VAL:HG21	1.81	0.63
1:A:271:ARG:HD3	1:A:283:GLN:OE1	1.99	0.62
1:A:257:HIS:CE1	1:A:261:LYS:HE3	2.34	0.62
1:B:269:GLN:HA	1:B:272:ARG:HB2	1.80	0.62
1:A:124:ASP:HB3	1:A:388:GLN:NE2	2.16	0.60
1:A:292:ASP:HB3	1:A:294:ARG:H	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:ARG:O	1:B:410:ARG:HG2	2.00	0.60
1:B:292:ASP:HB3	1:B:294:ARG:H	1.66	0.60
1:B:430:ASN:ND2	1:B:432:ALA:HB3	2.18	0.59
1:B:456:TYR:O	1:B:460:LYS:HB2	2.03	0.59
1:A:117:LEU:HD13	1:A:403:VAL:HG23	1.85	0.58
3:A:602[A]:ECN:C8	3:A:602[A]:ECN:H3	2.23	0.58
1:B:422:ASN:HD21	1:B:424:ASP:HB2	1.69	0.58
1:A:309:LEU:O	1:A:313:GLN:HB2	2.04	0.58
1:A:272:ARG:NH1	1:A:298:ASP:OD1	2.37	0.58
1:B:449:CYS:HA	2:B:601:HEM:C4D	2.40	0.57
1:B:416:VAL:O	1:B:425:ARG:NH2	2.38	0.57
1:A:116:LEU:HB2	1:A:390:VAL:HG21	1.87	0.56
1:A:234:PHE:CE2	3:A:602[A]:ECN:H17	2.40	0.56
1:B:135:THR:HG23	1:B:139:PHE:CD1	2.41	0.56
1:B:334:GLN:NE2	1:B:472:PHE:H	1.99	0.56
1:A:99:THR:HG22	1:A:104:THR:OG1	2.06	0.56
1:B:409:GLN:NE2	1:B:440:VAL:H	2.02	0.56
1:A:135:THR:HG21	1:A:145:TYR:HD1	1.70	0.55
1:A:271:ARG:HH12	1:A:280:ASP:CG	2.09	0.55
1:B:377:ILE:HD12	2:B:601:HEM:C4A	2.42	0.55
1:A:271:ARG:NH1	1:A:280:ASP:OD1	2.38	0.55
1:A:264:PHE:HE2	1:A:305:LEU:HB3	1.71	0.55
1:A:116:LEU:HB2	1:A:390:VAL:CG2	2.36	0.55
3:B:602[A]:ECN:CL4	3:B:602[A]:ECN:O20	2.62	0.55
1:B:227:TYR:CE2	1:B:309:LEU:HD22	2.43	0.54
1:B:430:ASN:HD22	1:B:432:ALA:HB3	1.72	0.54
1:B:150:PRO:HA	1:B:153:LEU:HD12	1.90	0.54
1:B:135:THR:HG22	6:B:2:HOH:O	2.08	0.54
1:B:129:ASP:OD2	1:B:385:ARG:HD3	2.08	0.54
1:B:278:ILE:O	1:B:283:GLN:HG2	2.08	0.53
1:A:290:TYR:HB2	1:A:292:ASP:HB2	1.90	0.53
1:A:416:VAL:HG12	1:A:417:GLU:HG3	1.90	0.53
1:A:444:ALA:HA	1:A:448:ARG:HG2	1.91	0.53
1:A:369:GLU:OE1	1:A:372:ARG:HD3	2.09	0.53
1:A:449:CYS:HB2	2:A:601:HEM:NA	2.24	0.52
1:B:415:TRP:CE2	1:B:431:PRO:HG2	2.44	0.52
1:B:281:ILE:HD12	1:B:281:ILE:N	2.25	0.52
1:A:369:GLU:OE1	1:A:369:GLU:HA	2.09	0.52
5:A:603:BCD:H52	5:A:603:BCD:O53	2.09	0.52
1:A:283:GLN:HA	1:A:286:LEU:HD12	1.91	0.51
1:A:410:ARG:HG2	1:A:418:ARG:NH1	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ALA:HA	1:A:152:PHE:CD2	2.46	0.51
1:A:208:HIS:CA	1:A:216:ARG:HG3	2.41	0.51
1:A:491:PRO:HB2	1:A:494:PRO:HB3	1.93	0.50
1:B:116:LEU:HB2	1:B:390:VAL:CG2	2.42	0.49
1:B:175:ILE:HG12	1:B:205:THR:HB	1.95	0.49
1:A:348:ASN:N	1:A:348:ASN:OD1	2.45	0.49
2:B:601:HEM:HBB2	2:B:601:HEM:CMB	2.42	0.48
1:A:380:MET:HG2	2:A:601:HEM:CGA	2.43	0.48
1:A:144:ALA:HA	1:A:152:PHE:CE2	2.48	0.48
1:B:369:GLU:HA	1:B:369:GLU:OE1	2.13	0.48
1:B:211:HIS:CD2	1:B:305:LEU:HD22	2.48	0.48
1:B:271:ARG:NH1	1:B:280:ASP:OD1	2.43	0.48
2:A:601:HEM:HMB2	2:A:601:HEM:HBB2	1.94	0.48
1:A:122:ASN:H	1:A:446:ARG:NH2	2.12	0.47
1:A:97:SER:HA	1:A:105:PHE:O	2.14	0.47
1:B:271:ARG:HH12	1:B:280:ASP:CG	2.18	0.47
1:A:389:THR:HA	1:A:393:TYR:O	2.14	0.47
1:A:409:GLN:NE2	1:A:440:VAL:H	2.04	0.46
1:B:269:GLN:HA	1:B:272:ARG:CB	2.45	0.46
1:B:136:THR:HB	1:B:137:PRO:HD3	1.96	0.46
3:A:602[A]:ECN:C8	3:A:602[A]:ECN:N1	2.59	0.46
1:B:97:SER:HA	1:B:105:PHE:O	2.16	0.46
1:B:79:LYS:HA	1:B:79:LYS:HD3	1.75	0.46
1:B:182:TYR:OH	1:B:197:ALA:HA	2.15	0.45
1:A:200:GLU:O	1:A:203:ILE:HG22	2.17	0.45
1:A:488:ILE:HG13	1:A:488:ILE:H	1.46	0.45
1:A:449:CYS:HB2	2:A:601:HEM:C1A	2.51	0.45
1:B:445:GLY:O	1:B:448:ARG:HB2	2.17	0.45
1:A:101:VAL:HG12	1:A:101:VAL:O	2.17	0.45
1:A:448:ARG:HD2	1:A:449:CYS:O	2.17	0.44
1:B:486:THR:HB	1:B:488:ILE:HG13	1.99	0.44
1:A:377:ILE:HD12	2:A:601:HEM:C4A	2.52	0.44
1:B:147:VAL:HG23	1:B:148:PRO:O	2.18	0.44
1:A:64:ILE:HG12	1:A:92:TYR:CE1	2.52	0.44
1:B:467:LEU:O	1:B:501:ARG:CD	2.65	0.44
1:A:375:PRO:O	1:A:489:HIS:HD2	2.00	0.44
1:B:209:CYS:O	1:B:281:ILE:HD13	2.18	0.44
1:B:66:SER:HB2	1:B:73:HIS:NE2	2.32	0.44
1:A:377:ILE:HD12	2:A:601:HEM:C3A	2.53	0.43
1:B:431:PRO:HB3	1:B:435:GLU:OE2	2.18	0.43
1:A:129:ASP:OD2	1:A:385:ARG:HD3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:LEU:HD11	1:B:239:TRP:CD1	2.53	0.43
1:B:211:HIS:CE1	1:B:309:LEU:HB2	2.53	0.43
1:A:208:HIS:HA	1:A:216:ARG:HG3	2.01	0.43
1:A:133:ARG:HG3	6:A:3:HOH:O	2.19	0.43
3:A:602[A]:ECN:CL4	3:A:602[A]:ECN:O20	2.74	0.42
1:A:409:GLN:NE2	1:A:441:PRO:HD3	2.34	0.42
1:A:136:THR:N	1:A:137:PRO:CD	2.83	0.42
1:B:156:LYS:HE3	2:B:601:HEM:HAD2	2.02	0.42
1:A:139:PHE:HE2	1:A:310:LEU:HD12	1.84	0.42
1:B:120:SER:HB2	6:B:22:HOH:O	2.20	0.42
1:A:164:ASN:HD22	1:A:166:ALA:H	1.67	0.42
1:A:413:ASP:N	1:A:413:ASP:OD1	2.53	0.42
1:B:248:PRO:HB2	1:B:252:ARG:NH2	2.35	0.42
1:A:327:LEU:HB3	1:A:334:GLN:HG3	2.00	0.41
1:A:268:ILE:HG23	1:A:286:LEU:HD21	2.02	0.41
1:A:449:CYS:HA	2:A:601:HEM:C4D	2.56	0.41
1:A:133:ARG:HG3	1:A:133:ARG:H	1.58	0.41
1:B:449:CYS:HA	2:B:601:HEM:CHA	2.50	0.41
1:A:309:LEU:O	1:A:309:LEU:HG	2.20	0.41
1:B:375:PRO:HA	1:B:376:PRO:HD3	1.98	0.41
1:B:198:LEU:O	1:B:202:ILE:HG12	2.21	0.41
1:A:369:GLU:HG3	1:A:421:PHE:CD1	2.56	0.41
1:A:430:ASN:HD22	1:A:432:ALA:HB3	1.82	0.40
1:B:95:VAL:HG21	1:B:390:VAL:HG12	2.02	0.40
1:B:164:ASN:HD22	1:B:164:ASN:C	2.23	0.40
1:B:311:ALA:HB2	4:B:600[B]:ECL:H9	2.03	0.40
1:B:258:ARG:O	1:B:262:ASP:HB2	2.20	0.40
1:A:378:MET:O	1:A:379:ILE:HG13	2.22	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	443/461 (96%)	424 (96%)	17 (4%)	2 (0%)	34	71
1	B	443/461 (96%)	416 (94%)	26 (6%)	1 (0%)	52	84
All	All	886/922 (96%)	840 (95%)	43 (5%)	3 (0%)	46	79

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	292	ASP
1	A	489	HIS
1	B	488	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	390/403 (97%)	368 (94%)	22 (6%)	26	60
1	B	390/403 (97%)	354 (91%)	36 (9%)	11	33
All	All	780/806 (97%)	722 (93%)	58 (7%)	17	44

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	VAL
1	A	68	ILE
1	A	133	ARG
1	A	135	THR
1	A	194	VAL
1	A	203	ILE
1	A	216	ARG
1	A	236	HIS
1	A	252	ARG
1	A	272	ARG
1	A	304	MET
1	A	323	MET
1	A	358	LYS

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Mol	Chain	Res	Type
1	A	385	ARG
1	A	390	VAL
1	A	400	GLN
1	A	413	ASP
1	A	420	ASP
1	A	448	ARG
1	A	488	ILE
1	A	490	THR
1	A	501	ARG
1	B	90	GLU
1	B	91	LYS
1	B	123	GLU
1	B	136	THR
1	B	143	VAL
1	B	157	LYS
1	B	160	LYS
1	B	164	ASN
1	B	194	VAL
1	B	207	SER
1	B	216	ARG
1	B	217	SER
1	B	247	LEU
1	B	259	GLU
1	B	260	ILE
1	B	266	LYS
1	B	269	GLN
1	B	272	ARG
1	B	289	THR
1	B	298	ASP
1	B	304	MET
1	B	309	LEU
1	B	347	GLU
1	B	355	ASP
1	B	359	ASP
1	B	385	ARG
1	B	386	THR
1	B	390	VAL
1	B	400	GLN
1	B	410	ARG
1	B	413	ASP
1	B	420	ASP
1	B	428	GLN

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Mol	Chain	Res	Type
1	B	448	ARG
1	B	488	ILE
1	B	501	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	149	ASN
1	A	164	ASN
1	A	170	GLN
1	A	257	HIS
1	A	275	GLN
1	A	334	GLN
1	A	388	GLN
1	A	409	GLN
1	A	422	ASN
1	A	430	ASN
1	A	489	HIS
1	B	164	ASN
1	B	334	GLN
1	B	388	GLN
1	B	409	GLN
1	B	422	ASN
1	B	430	ASN

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

8 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$
4	ECL	A	600[B]	2	24,26,26	1.08	3 (12%)	31,35,35	1.26	4 (12%)
2	HEM	A	601	1,3,4	30,50,50	2.22	7 (23%)	24,82,82	2.35	7 (29%)
3	ECN	A	602[A]	2	24,26,26	0.99	3 (12%)	31,35,35	1.46	3 (9%)
5	BCD	A	603	-	84,84,84	0.47	0	126,126,126	1.14	10 (7%)
4	ECL	B	600[B]	2	24,26,26	1.35	6 (25%)	31,35,35	1.33	3 (9%)
2	HEM	B	601	1,3,4	30,50,50	2.07	6 (20%)	24,82,82	2.23	5 (20%)
3	ECN	B	602[A]	2	24,26,26	1.17	3 (12%)	31,35,35	1.50	3 (9%)
5	BCD	B	603	-	84,84,84	0.46	0	126,126,126	1.27	14 (11%)

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	ECL	A	600[B]	2	-	0/13/13/13	0/3/3/3
2	HEM	A	601	1,3,4	-	0/10/54/54	0/0/8/8
3	ECN	A	602[A]	2	-	2/13/13/13	0/3/3/3
5	BCD	A	603	-	-	0/42/182/182	0/0/8/8
4	ECL	B	600[B]	2	-	0/13/13/13	0/3/3/3
2	HEM	B	601	1,3,4	-	0/10/54/54	0/0/8/8
3	ECN	B	602[A]	2	-	2/13/13/13	0/3/3/3
5	BCD	B	603	-	-	0/42/182/182	0/0/8/8

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HEM	C3B-C4B	-7.44	1.45	1.51
2	B	601	HEM	C3B-C4B	-6.55	1.46	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HEM	C3D-C4D	-5.48	1.44	1.51
2	B	601	HEM	C3D-C4D	-4.93	1.45	1.51
2	A	601	HEM	C2C-C1C	-3.95	1.45	1.52
2	B	601	HEM	C2C-C1C	-3.87	1.45	1.52
3	B	602[A]	ECN	C19-N1	-3.00	1.45	1.48
4	B	600[B]	ECL	C19-N1	-2.50	1.45	1.48
2	A	601	HEM	C2D-C1D	-2.18	1.44	1.51
4	B	600[B]	ECL	C7-N1	-2.07	1.33	1.37
2	B	601	HEM	C2B-C1B	-2.00	1.45	1.51
3	A	602[A]	ECN	C7-N1	2.10	1.41	1.37
3	A	602[A]	ECN	C11-CL2	2.17	1.79	1.74
4	B	600[B]	ECL	C14-CL4	2.17	1.79	1.74
2	A	601	HEM	CAA-C2A	2.20	1.55	1.52
2	B	601	HEM	FE-ND	2.21	2.09	1.97
3	B	602[A]	ECN	C11-CL2	2.21	1.79	1.74
4	A	600[B]	ECL	C14-CL4	2.30	1.79	1.74
3	A	602[A]	ECN	C16-CL8	2.37	1.79	1.74
4	A	600[B]	ECL	C16-CL8	2.45	1.79	1.73
3	B	602[A]	ECN	C14-CL4	2.46	1.79	1.73
2	B	601	HEM	C3B-CAB	2.54	1.56	1.51
4	B	600[B]	ECL	C11-CL2	2.56	1.80	1.74
4	A	600[B]	ECL	C11-CL2	2.63	1.80	1.74
4	B	600[B]	ECL	C16-CL8	2.69	1.80	1.73
2	A	601	HEM	C3B-CAB	2.75	1.56	1.51
4	B	600[B]	ECL	O20-C8	2.76	1.50	1.42
2	A	601	HEM	FE-ND	2.78	2.12	1.97

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602[A]	ECN	O20-C20-C5	-5.19	104.55	112.41
2	A	601	HEM	C3C-CAC-CBC	-4.52	117.53	124.46
3	B	602[A]	ECN	O20-C20-C5	-4.36	105.80	112.41
3	B	602[A]	ECN	C14-C5-C20	-3.85	117.30	122.20
4	B	600[B]	ECL	O20-C8-C1	-3.60	101.40	109.99
4	A	600[B]	ECL	O20-C8-C1	-3.24	102.25	109.99
5	B	603	BCD	C14-O43-C43	-3.13	109.83	118.01
5	B	603	BCD	C61-C51-C41	-2.51	105.94	113.25
3	A	602[A]	ECN	C7-N1-C3	-2.44	105.84	108.20
4	A	600[B]	ECL	O20-C20-C17	-2.27	108.97	112.41
5	A	603	BCD	C13-O42-C42	-2.26	112.09	118.01
5	B	603	BCD	C12-O41-C41	-2.25	112.13	118.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	600[B]	ECL	C15-C16-C17	-2.19	119.62	122.46
5	B	603	BCD	O42-C13-O53	-2.17	105.18	110.68
5	B	603	BCD	O36-C36-C46	-2.09	104.93	109.87
5	B	603	BCD	C32-C42-C52	2.02	115.40	110.84
5	B	603	BCD	C22-C32-C42	2.11	114.24	109.60
5	A	603	BCD	O51-C51-C61	2.12	111.70	106.36
5	B	603	BCD	O51-C51-C41	2.13	114.25	109.75
5	A	603	BCD	O52-C12-C22	2.22	114.83	110.28
5	B	603	BCD	C26-C36-C46	2.29	114.63	109.60
3	A	602[A]	ECN	C21-C5-C14	2.34	119.32	116.79
5	A	603	BCD	C21-C31-C41	2.40	114.86	109.60
5	B	603	BCD	C15-C25-C35	2.40	114.70	109.97
5	A	603	BCD	O47-C11-C21	2.43	114.01	108.10
4	A	600[B]	ECL	C17-C16-CL8	2.49	123.17	120.42
4	A	600[B]	ECL	C21-C17-C16	2.53	119.53	116.79
2	A	601	HEM	C2D-C3D-C4D	2.54	105.81	101.50
2	A	601	HEM	CMD-C2D-C3D	2.57	125.72	114.35
2	B	601	HEM	CMD-C2D-C3D	2.95	127.39	114.35
5	A	603	BCD	O56-C56-C46	3.01	116.11	109.75
5	A	603	BCD	C27-C37-C47	3.08	116.37	109.60
5	A	603	BCD	C16-O56-C56	3.10	119.76	113.75
5	A	603	BCD	O52-C52-C42	3.13	116.36	109.75
5	B	603	BCD	C17-C27-C37	3.24	116.35	109.97
5	B	603	BCD	C12-O52-C52	3.55	120.64	113.75
4	B	600[B]	ECL	C21-C17-C16	3.86	120.96	116.79
5	A	603	BCD	C12-O52-C52	3.88	121.28	113.75
3	B	602[A]	ECN	C21-C5-C14	4.06	121.18	116.79
5	B	603	BCD	C11-O51-C51	4.16	121.81	113.75
5	B	603	BCD	C27-C37-C47	4.21	118.85	109.60
2	B	601	HEM	CAD-C3D-C4D	4.22	127.34	112.47
2	A	601	HEM	CAD-C3D-C2D	4.28	125.53	113.22
2	B	601	HEM	CMB-C2B-C3B	4.42	127.57	116.53
2	A	601	HEM	CAD-C3D-C4D	4.58	128.64	112.47
2	A	601	HEM	CMB-C2B-C3B	4.60	128.02	116.53
2	A	601	HEM	CMC-C2C-C3C	4.66	128.15	116.53
2	B	601	HEM	CMC-C2C-C3C	4.68	128.22	116.53
2	B	601	HEM	CAD-C3D-C2D	5.17	128.08	113.22

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	602[A]	ECN	C8-O20-C20-C19
3	A	602[A]	ECN	C8-O20-C20-C5
3	B	602[A]	ECN	C8-O20-C20-C5
3	A	602[A]	ECN	C8-O20-C20-C19

There are no ring outliers.

6 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	7	0
3	A	602[A]	ECN	8	0
5	A	603	BCD	1	0
4	B	600[B]	ECL	1	0
2	B	601	HEM	6	0
3	B	602[A]	ECN	7	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	445/461 (96%)	0.09	18 (4%) 42 35	39, 56, 82, 98	0
1	B	445/461 (96%)	0.29	33 (7%) 17 11	47, 71, 95, 106	0
All	All	890/922 (96%)	0.19	51 (5%) 27 21	39, 62, 92, 106	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	276	GLU	4.8
1	B	485	THR	4.4
1	A	244	TRP	4.3
1	B	277	LYS	4.0
1	B	275	GLN	3.6
1	A	463	TRP	3.5
1	B	308	LEU	3.5
1	B	305	LEU	3.3
1	B	246	PRO	3.2
1	A	275	GLN	3.2
1	B	249	SER	3.0
1	B	488	ILE	3.0
1	B	304	MET	2.9
1	B	67	PRO	2.9
1	B	248	PRO	2.9
1	B	348	ASN	2.9
1	A	486	THR	2.8
1	A	276	GLU	2.8
1	A	221	GLU	2.8
1	A	433	SER	2.7
1	B	282	LEU	2.7
1	B	347	GLU	2.6
1	B	211	HIS	2.6
1	B	502	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	255	ARG	2.5
1	B	244	TRP	2.4
1	B	210	LEU	2.4
1	A	243	GLY	2.4
1	B	492	GLU	2.3
1	B	433	SER	2.3
1	A	245	LEU	2.3
1	A	67	PRO	2.3
1	A	242	PRO	2.3
1	A	461	THR	2.3
1	B	486	THR	2.3
1	A	462	ILE	2.2
1	B	342	LYS	2.2
1	A	246	PRO	2.2
1	B	237	ALA	2.2
1	A	255	ARG	2.1
1	B	247	LEU	2.1
1	B	487	MET	2.1
1	A	485	THR	2.1
1	B	225	GLN	2.1
1	B	463	TRP	2.1
1	B	258	ARG	2.1
1	A	251	ARG	2.1
1	A	305	LEU	2.0
1	B	181	GLU	2.0
1	B	435	GLU	2.0
1	B	281	ILE	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

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	LLDF	B-factors(Å ²)	Q<0.9
3	ECN	A	602[A]	24/24	0.94	0.20	0.29	54,59,62,65	24
4	ECL	A	600[B]	24/24	0.94	0.19	0.12	60,63,69,72	24
3	ECN	B	602[A]	24/24	0.89	0.24	0.06	74,78,81,82	24
4	ECL	B	600[B]	24/24	0.89	0.22	-0.08	66,69,72,73	24
5	BCD	B	603	77/77	0.92	0.17	-0.22	58,68,71,72	0
2	HEM	B	601	43/43	0.98	0.15	-0.67	47,51,52,54	0
5	BCD	A	603	77/77	0.93	0.15	-0.84	69,75,82,83	0
2	HEM	A	601	43/43	0.98	0.13	-0.98	42,47,52,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.