



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:48 AM GMT

PDB ID : 2BO4
Title : DISSECTION OF MANNOSYLGLYCERATE SYNTHASE: AN ARCHETYPAL MANNOSYLTRANSFERASE
Authors : Flint, J.; Taylor, E.; Yang, M.; Bolam, D.N.; Tailford, L.E.; Martinez-Fleites, C.; Dodson, E.J.; Davis, B.G.; Gilbert, H.J.; Davies, G.J.
Deposited on : 2005-04-07
Resolution : 1.95 Å(reported)

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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

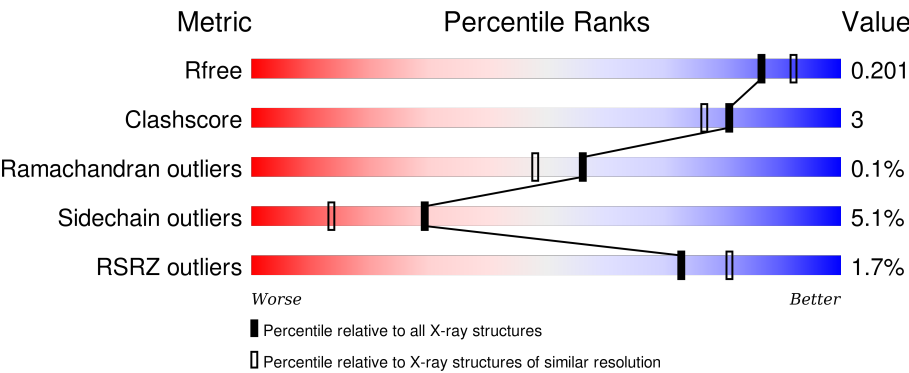
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	397	<div><div>2%</div><div>84%</div><div>10%</div><div>..</div></div>
1	B	397	<div><div>%</div><div>84%</div><div>10%</div><div>..</div></div>
1	C	397	<div><div>2%</div><div>83%</div><div>12%</div><div>..</div></div>
1	D	397	<div><div>2%</div><div>84%</div><div>10%</div><div>..</div></div>
1	E	397	<div><div>2%</div><div>84%</div><div>11%</div><div>..</div></div>

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Mol	Chain	Length	Quality of chain
1	F	397	<div> <div></div> <div>%</div> <div>83%</div> <div>11%</div> <div>• •</div> </div>

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	FLC	A	1382	-	-	-	X
2	FLC	B	1382	-	-	-	X
2	FLC	D	1382	-	-	-	X
2	FLC	F	1382	-	-	-	X

2 Entry composition [i](#)

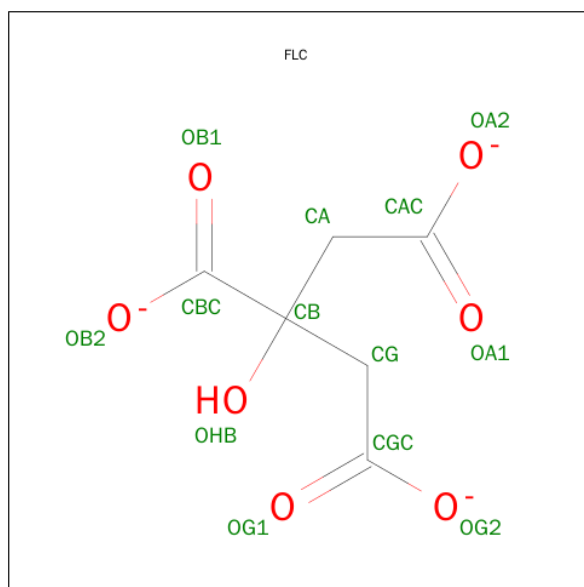
There are 3 unique types of molecules in this entry. The entry contains 20669 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MANNOSYLGLYCERATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	0	1
			3138	2006	558	559	15			
1	B	380	Total	C	N	O	S	0	0	0
			3137	2006	557	559	15			
1	C	381	Total	C	N	O	S	0	0	1
			3138	2006	558	559	15			
1	D	380	Total	C	N	O	S	0	0	0
			3137	2006	557	559	15			
1	E	380	Total	C	N	O	S	0	0	0
			3137	2006	557	559	15			
1	F	380	Total	C	N	O	S	0	0	0
			3137	2006	557	559	15			

- Molecule 2 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7^-$).



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

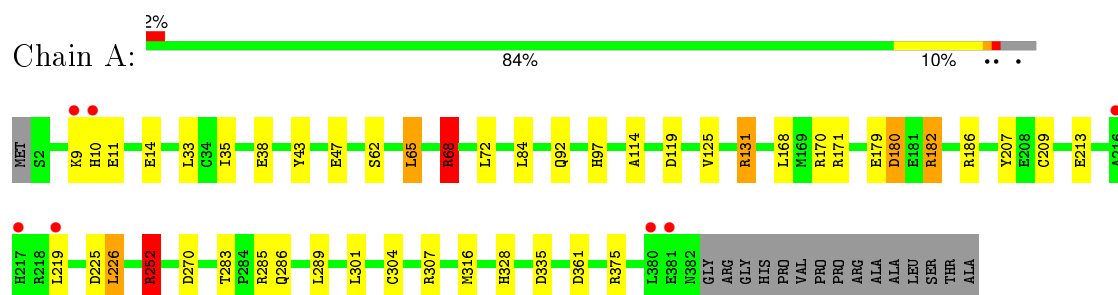
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	302	Total O 302 302	0	0
3	B	288	Total O 288 288	0	0
3	C	301	Total O 301 301	0	0
3	D	289	Total O 289 289	0	0
3	E	301	Total O 301 301	0	0
3	F	286	Total O 286 286	0	0

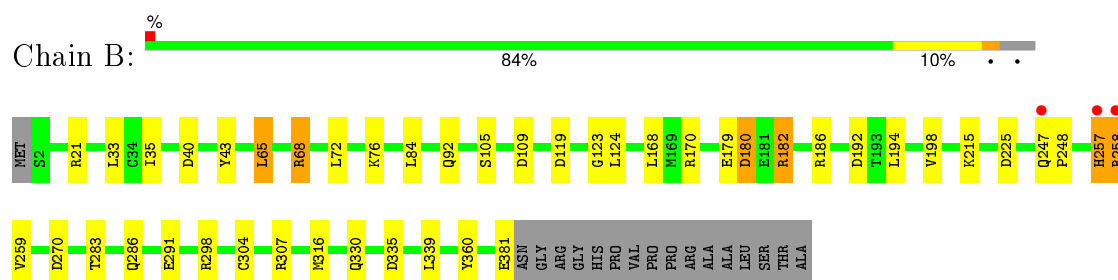
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

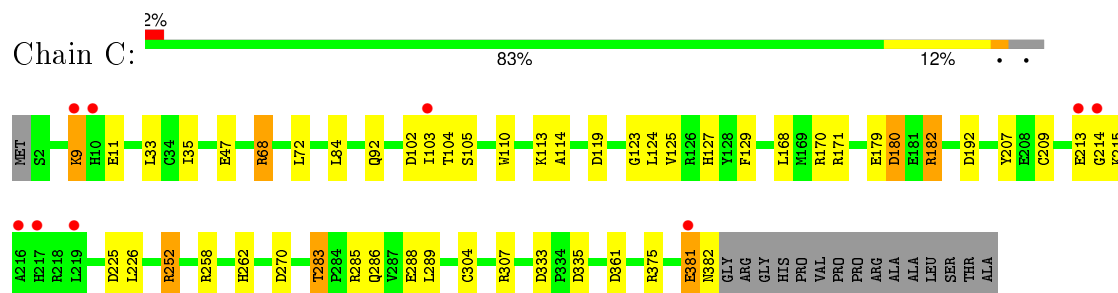
• Molecule 1: MANNOSYLGLYCERATE SYNTHASE



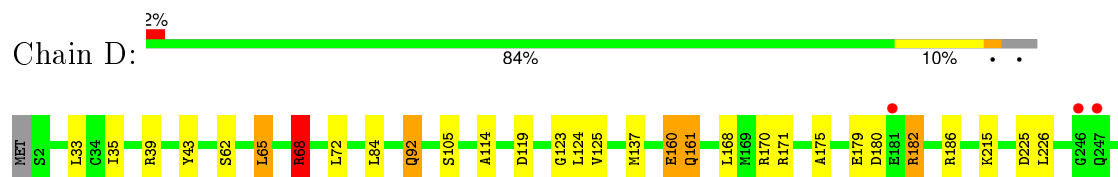
• Molecule 1: MANNOSYLGLYCERATE SYNTHASE



• Molecule 1: MANNOSYLGLYCERATE SYNTHASE

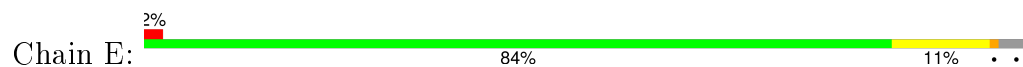


• Molecule 1: MANNOSYLGLYCERATE SYNTHASE

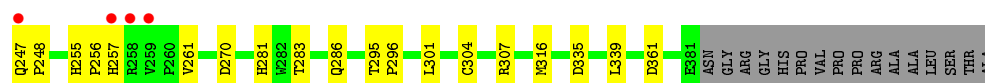
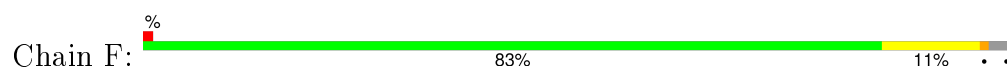




• Molecule 1: MANNOSYLGLYCERATE SYNTHASE



• Molecule 1: MANNOSYLGLYCERATE SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	261.15Å 151.29Å 153.24Å 90.00° 90.41° 90.00°	Depositor
Resolution (Å)	20.00 – 1.95 19.99 – 1.95	Depositor EDS
% Data completeness (in resolution range)	95.7 (20.00-1.95) 94.8 (19.99-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.176 , 0.192 0.186 , 0.201	Depositor DCC
R_{free} test set	20756 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
Estimated twinning fraction	0.008 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.009 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.377 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.357 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.010 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	4 of 413730 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	20669	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: FLC

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	0.46	0/3229	0.80	15/4394 (0.3%)
1	B	0.45	0/3228	0.77	10/4392 (0.2%)
1	C	0.48	0/3229	0.82	18/4394 (0.4%)
1	D	0.47	1/3228 (0.0%)	0.80	12/4392 (0.3%)
1	E	0.47	0/3228	0.81	14/4392 (0.3%)
1	F	0.44	0/3228	0.80	12/4392 (0.3%)
All	All	0.46	1/19370 (0.0%)	0.80	81/26356 (0.3%)

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	A	1	0
1	B	1	0
1	E	1	0
1	F	1	0
All	All	4	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	137	MET	SD-CE	-7.30	1.36	1.77

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	161	GLN	N-CA-CB	-11.26	90.33	110.60
1	A	375	ARG	NE-CZ-NH2	-10.97	114.81	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	375	ARG	NE-CZ-NH2	-10.45	115.07	120.30
1	D	161	GLN	N-CA-CB	-10.45	91.80	110.60
1	C	375	ARG	NE-CZ-NH2	-9.38	115.61	120.30
1	E	68	ARG	NE-CZ-NH2	9.36	124.98	120.30
1	A	375	ARG	NE-CZ-NH1	8.32	124.46	120.30
1	E	375	ARG	NE-CZ-NH1	8.21	124.40	120.30
1	F	68	ARG	NE-CZ-NH2	8.16	124.38	120.30
1	C	68	ARG	NE-CZ-NH2	8.04	124.32	120.30
1	C	270	ASP	CB-CG-OD2	8.01	125.51	118.30
1	D	68	ARG	NE-CZ-NH2	7.77	124.19	120.30
1	E	270	ASP	CB-CG-OD2	7.71	125.24	118.30
1	B	68	ARG	NE-CZ-NH2	7.63	124.11	120.30
1	A	68	ARG	NE-CZ-NH2	7.59	124.09	120.30
1	B	270	ASP	CB-CG-OD2	7.41	124.97	118.30
1	C	182	ARG	NE-CZ-NH1	7.14	123.87	120.30
1	D	182	ARG	NE-CZ-NH1	7.12	123.86	120.30
1	B	182	ARG	NE-CZ-NH1	7.11	123.85	120.30
1	C	225	ASP	CB-CG-OD2	7.06	124.66	118.30
1	F	119	ASP	CB-CG-OD2	7.01	124.61	118.30
1	D	119	ASP	CB-CG-OD2	6.99	124.59	118.30
1	C	361	ASP	CB-CG-OD2	6.96	124.56	118.30
1	D	160	GLU	N-CA-C	6.95	129.75	111.00
1	A	270	ASP	CB-CG-OD2	6.82	124.44	118.30
1	E	182	ARG	NE-CZ-NH2	-6.67	116.97	120.30
1	A	361	ASP	CB-CG-OD2	6.66	124.29	118.30
1	C	375	ARG	NE-CZ-NH1	6.65	123.62	120.30
1	F	182	ARG	NE-CZ-NH1	6.61	123.61	120.30
1	D	270	ASP	CB-CG-OD2	6.58	124.22	118.30
1	C	182	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	C	68	ARG	NE-CZ-NH1	-6.54	117.03	120.30
1	E	361	ASP	CB-CG-OD2	6.53	124.18	118.30
1	B	182	ARG	NE-CZ-NH2	-6.49	117.06	120.30
1	E	225	ASP	CB-CG-OD2	6.46	124.11	118.30
1	F	270	ASP	CB-CG-OD2	6.44	124.10	118.30
1	A	182	ARG	NE-CZ-NH2	-6.43	117.08	120.30
1	A	225	ASP	CB-CG-OD2	6.40	124.06	118.30
1	D	225	ASP	CB-CG-OD2	6.34	124.01	118.30
1	E	252	ARG	NE-CZ-NH1	6.30	123.45	120.30
1	B	119	ASP	CB-CG-OD2	6.27	123.95	118.30
1	C	252	ARG	NE-CZ-NH2	-6.26	117.17	120.30
1	E	68	ARG	NE-CZ-NH1	-6.20	117.20	120.30
1	E	252	ARG	NE-CZ-NH2	-6.09	117.26	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	258	ARG	NE-CZ-NH2	6.08	123.34	120.30
1	A	182	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	A	252	ARG	NE-CZ-NH2	-6.01	117.30	120.30
1	D	182	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	E	171	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	C	252	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	F	161	GLN	N-CA-C	-5.98	94.85	111.00
1	F	160	GLU	N-CA-C	5.93	127.02	111.00
1	B	225	ASP	CB-CG-OD2	5.85	123.56	118.30
1	E	182	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	A	68	ARG	NE-CZ-NH1	-5.80	117.40	120.30
1	C	171	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	B	109	ASP	CB-CG-OD2	5.73	123.46	118.30
1	F	225	ASP	CB-CG-OD2	5.69	123.42	118.30
1	A	131	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	E	171	ARG	NE-CZ-NH2	-5.64	117.48	120.30
1	A	119	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	252	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	C	119	ASP	CB-CG-OD2	5.57	123.31	118.30
1	F	182	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	B	40	ASP	CB-CG-OD2	5.54	123.28	118.30
1	B	180	ASP	CB-CG-OD2	5.48	123.23	118.30
1	F	109	ASP	CB-CG-OD2	5.47	123.22	118.30
1	D	161	GLN	N-CA-C	-5.42	96.37	111.00
1	A	171	ARG	NE-CZ-NH1	5.37	122.99	120.30
1	C	192	ASP	CB-CG-OD2	5.36	123.13	118.30
1	F	40	ASP	CB-CG-OD2	5.36	123.13	118.30
1	D	375	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	D	171	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	C	258	ARG	NE-CZ-NH1	-5.31	117.65	120.30
1	E	119	ASP	CB-CG-OD2	5.25	123.02	118.30
1	A	180	ASP	CB-CG-OD2	5.24	123.02	118.30
1	C	333	ASP	CB-CG-OD2	5.23	123.01	118.30
1	D	68	ARG	NE-CZ-NH1	-5.20	117.70	120.30
1	B	68	ARG	NE-CZ-NH1	-5.09	117.75	120.30
1	F	361	ASP	CB-CG-OD2	5.08	122.87	118.30
1	C	180	ASP	CB-CG-OD2	5.06	122.85	118.30

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	283	THR	CB

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Mol	Chain	Res	Type	Atom
1	B	283	THR	CB
1	E	283	THR	CB
1	F	283	THR	CB

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	3138	0	3037	22	0
1	B	3137	0	3037	24	0
1	C	3138	0	3037	24	0
1	D	3137	0	3037	22	0
1	E	3137	0	3037	17	0
1	F	3137	0	3037	19	0
2	A	13	0	5	2	0
2	B	13	0	5	0	0
2	C	13	0	5	0	0
2	D	13	0	5	0	0
2	E	13	0	5	0	0
2	F	13	0	5	0	0
3	A	302	0	0	4	0
3	B	288	0	0	6	0
3	C	301	0	0	4	0
3	D	289	0	0	4	0
3	E	301	0	0	3	0
3	F	286	0	0	3	0
All	All	20669	0	18252	124	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 3.

All (124) 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:E:316:MET:HE2	1:E:316:MET:HA	1.55	0.88
1:D:283:THR:HG22	1:D:286:GLN:H	1.36	0.88
1:E:283:THR:HG21	1:E:335:ASP:OD2	1.81	0.81
1:A:283:THR:HG21	1:A:335:ASP:OD2	1.84	0.78
1:B:283:THR:HG21	1:B:335:ASP:OD2	1.85	0.76
1:A:68:ARG:HD3	3:D:2160:HOH:O	1.85	0.75
1:D:283:THR:HG21	1:D:335:ASP:OD2	1.91	0.69
1:B:283:THR:HG23	1:B:286:GLN:H	1.59	0.68
1:F:76:LYS:NZ	1:F:192:ASP:OD2	2.27	0.68
1:C:283:THR:HG22	1:C:286:GLN:H	1.59	0.68
1:F:283:THR:HG21	1:F:335:ASP:OD2	1.95	0.67
1:D:175:ALA:O	1:D:179:GLU:HG2	1.94	0.66
1:D:180:ASP:OD1	1:D:182:ARG:HD3	1.95	0.66
1:A:180:ASP:OD1	1:A:182:ARG:HD3	1.96	0.65
1:A:283:THR:HG23	1:A:286:GLN:H	1.61	0.64
1:C:11:GLU:CG	1:C:102:ASP:HB3	2.28	0.64
1:C:180:ASP:OD1	1:C:182:ARG:HD3	1.97	0.64
1:E:283:THR:HG23	1:E:286:GLN:H	1.63	0.63
1:E:180:ASP:OD1	1:E:182:ARG:HD3	1.97	0.63
1:F:283:THR:HG23	1:F:286:GLN:H	1.64	0.63
1:B:123:GLY:HA2	1:B:170:ARG:CD	2.29	0.63
1:B:180:ASP:OD1	1:B:182:ARG:HD3	1.99	0.63
1:C:11:GLU:HG2	1:C:102:ASP:HB3	1.81	0.62
1:F:123:GLY:HA2	1:F:170:ARG:CD	2.29	0.62
1:D:123:GLY:HA2	1:D:170:ARG:CD	2.29	0.62
1:D:43:TYR:CE2	1:D:65:LEU:HD13	2.34	0.62
1:F:123:GLY:HA2	1:F:170:ARG:HD3	1.82	0.61
1:E:316:MET:HE3	3:E:2254:HOH:O	2.00	0.61
1:B:123:GLY:HA2	1:B:170:ARG:HD3	1.83	0.60
1:A:226:LEU:HD11	2:A:1382:FLC:OA2	2.01	0.60
1:D:33:LEU:HD21	1:D:35:ILE:HD11	1.84	0.60
1:B:258:ARG:HG2	3:B:2188:HOH:O	2.03	0.58
1:C:113:LYS:NZ	3:C:2087:HOH:O	2.37	0.58
1:A:38:GLU:HB2	1:D:68:ARG:HD2	1.86	0.58
1:F:180:ASP:OD1	1:F:182:ARG:HD3	2.04	0.58
1:C:283:THR:HG21	1:C:335:ASP:OD2	2.04	0.56
1:F:33:LEU:HD21	1:F:35:ILE:HD11	1.89	0.55
1:B:259:VAL:HG21	1:B:360:TYR:CE1	2.42	0.55
1:B:43:TYR:CE2	1:B:65:LEU:HD13	2.43	0.54
1:F:43:TYR:CE2	1:F:65:LEU:HD13	2.43	0.54
1:B:330:GLN:NE2	3:B:2250:HOH:O	2.41	0.53
1:C:33:LEU:HD21	1:C:35:ILE:HD11	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:283:THR:HG23	3:C:2231:HOH:O	2.08	0.53
1:C:123:GLY:HA2	1:C:170:ARG:CD	2.39	0.53
1:E:9:LYS:HB2	3:E:2004:HOH:O	2.10	0.51
1:F:9:LYS:HE3	3:F:2006:HOH:O	2.09	0.51
1:D:123:GLY:HA2	1:D:170:ARG:HD3	1.92	0.51
1:F:194:LEU:O	1:F:198:VAL:HG13	2.11	0.51
1:B:76:LYS:NZ	1:B:192:ASP:OD2	2.42	0.50
1:A:43:TYR:CE2	1:A:65:LEU:HD13	2.46	0.50
1:C:381:GLU:HG2	1:C:382:ASN:N	2.26	0.50
3:B:2189:HOH:O	1:F:261:VAL:HG21	2.11	0.49
1:A:170:ARG:HD2	3:A:2088:HOH:O	2.11	0.49
1:A:304:CYS:HA	1:A:307:ARG:O	2.12	0.49
1:E:316:MET:HA	1:E:316:MET:CE	2.36	0.49
1:B:33:LEU:HD21	1:B:35:ILE:HD11	1.94	0.48
1:B:194:LEU:O	1:B:198:VAL:HG13	2.13	0.48
1:C:207:TYR:HE2	1:C:209:CYS:HB3	1.78	0.48
1:C:11:GLU:HG3	1:C:102:ASP:HB3	1.95	0.48
1:E:170:ARG:HD2	3:E:2100:HOH:O	2.14	0.48
1:B:21:ARG:HD2	3:B:2029:HOH:O	2.14	0.47
1:C:207:TYR:CE2	1:C:209:CYS:HB3	2.49	0.47
1:A:131:ARG:NH2	2:A:1382:FLC:OHB	2.48	0.47
1:C:123:GLY:HA2	1:C:170:ARG:HD3	1.96	0.47
1:B:259:VAL:HG21	1:B:360:TYR:HE1	1.78	0.47
1:A:97:HIS:CE1	1:A:168:LEU:HD12	2.50	0.47
1:B:247:GLN:N	1:B:248:PRO:CD	2.78	0.47
1:E:210:TYR:CE2	1:E:212:PRO:HA	2.51	0.46
1:D:328:HIS:HD2	3:D:2079:HOH:O	1.98	0.46
1:B:316:MET:HA	1:B:316:MET:CE	2.45	0.46
1:F:281:HIS:CD2	3:F:2207:HOH:O	2.68	0.46
1:C:213:GLU:OE2	1:C:262:HIS:CE1	2.69	0.46
1:D:255:HIS:HD2	1:D:256:PRO:O	1.99	0.46
1:E:123:GLY:HA2	1:E:170:ARG:HD3	1.98	0.45
1:F:304:CYS:HA	1:F:307:ARG:O	2.15	0.45
1:A:62:SER:OG	3:A:2032:HOH:O	2.21	0.45
1:A:33:LEU:HD21	1:A:35:ILE:HD11	1.98	0.45
1:C:129:PHE:CD1	1:C:214:GLY:HA3	2.51	0.45
1:D:168:LEU:C	1:D:168:LEU:HD23	2.37	0.45
1:F:255:HIS:HD2	1:F:256:PRO:O	1.99	0.45
1:F:316:MET:HA	1:F:316:MET:CE	2.47	0.45
1:D:316:MET:HA	1:D:316:MET:CE	2.48	0.44
1:A:207:TYR:HE2	1:A:209:CYS:HB3	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:THR:HG22	1:B:286:GLN:NE2	2.32	0.44
1:E:43:TYR:CE2	1:E:65:LEU:HD13	2.53	0.44
1:E:304:CYS:HA	1:E:307:ARG:O	2.18	0.44
1:C:182:ARG:HD2	3:C:2130:HOH:O	2.18	0.44
1:A:316:MET:HG2	3:A:2077:HOH:O	2.17	0.43
1:E:114:ALA:HA	1:E:125:VAL:HG11	1.99	0.43
1:A:252:ARG:HD3	3:F:2100:HOH:O	2.17	0.43
1:D:304:CYS:HA	1:D:307:ARG:O	2.17	0.43
1:D:92:GLN:HG2	3:D:2055:HOH:O	2.18	0.43
1:B:168:LEU:HD23	1:B:168:LEU:C	2.37	0.43
1:A:283:THR:HG22	1:A:286:GLN:NE2	2.34	0.43
1:F:247:GLN:N	1:F:248:PRO:CD	2.80	0.43
1:C:304:CYS:HA	1:C:307:ARG:O	2.18	0.43
1:D:114:ALA:HA	1:D:125:VAL:HG11	2.01	0.43
1:B:257:HIS:HD2	3:B:2188:HOH:O	2.02	0.43
1:E:103:ILE:HG21	1:E:110:TRP:CH2	2.54	0.43
1:F:168:LEU:HD23	1:F:168:LEU:C	2.39	0.42
1:C:213:GLU:OE2	1:C:262:HIS:NE2	2.52	0.42
1:D:316:MET:HE3	3:D:2242:HOH:O	2.18	0.42
1:A:328:HIS:HD2	3:A:2083:HOH:O	2.02	0.42
1:C:103:ILE:HG21	1:C:110:TRP:CH2	2.55	0.42
1:B:291:GLU:HA	1:B:298:ARG:HD3	2.00	0.42
1:A:207:TYR:CE2	1:A:209:CYS:HB3	2.55	0.42
1:B:283:THR:CG2	1:B:335:ASP:OD2	2.63	0.42
1:C:110:TRP:O	1:C:127:HIS:HE1	2.03	0.42
1:B:258:ARG:HD2	1:F:261:VAL:HG13	2.01	0.42
1:B:304:CYS:HA	1:B:307:ARG:O	2.19	0.42
1:C:168:LEU:HD23	1:C:168:LEU:C	2.41	0.41
1:B:316:MET:HE3	3:B:2241:HOH:O	2.20	0.41
1:E:207:TYR:HE2	1:E:209:CYS:HB3	1.85	0.41
1:A:114:ALA:HA	1:A:125:VAL:HG11	2.01	0.41
1:C:114:ALA:HA	1:C:125:VAL:HG11	2.03	0.41
1:D:362:TYR:O	1:D:365:GLN:HG3	2.20	0.41
1:A:38:GLU:OE1	1:D:68:ARG:HD3	2.20	0.41
1:A:38:GLU:OE1	1:D:68:ARG:CD	2.68	0.41
1:D:160:GLU:O	1:D:161:GLN:HG3	2.21	0.41
1:F:295:THR:HB	1:F:296:PRO:HD3	2.02	0.41
1:D:277:ARG:CZ	1:D:281:HIS:HD2	2.34	0.41
1:E:97:HIS:CE1	1:E:168:LEU:HD12	2.56	0.41
1:C:213:GLU:HG3	3:C:2158:HOH:O	2.21	0.40
1:E:295:THR:HB	1:E:296:PRO:HD3	2.04	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	379/397 (96%)	368 (97%)	9 (2%)	2 (0%)	34	21
1	B	378/397 (95%)	371 (98%)	7 (2%)	0	100	100
1	C	379/397 (96%)	366 (97%)	12 (3%)	1 (0%)	46	35
1	D	378/397 (95%)	370 (98%)	8 (2%)	0	100	100
1	E	378/397 (95%)	369 (98%)	9 (2%)	0	100	100
1	F	378/397 (95%)	369 (98%)	9 (2%)	0	100	100
All	All	2270/2382 (95%)	2213 (98%)	54 (2%)	3 (0%)	56	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	HIS
1	A	11	GLU
1	C	9	LYS

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	326/338 (96%)	309 (95%)	17 (5%)	29	13
1	B	326/338 (96%)	312 (96%)	14 (4%)	35	20

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	326/338 (96%)	308 (94%)	18 (6%)	27	12
1	D	326/338 (96%)	310 (95%)	16 (5%)	31	15
1	E	326/338 (96%)	308 (94%)	18 (6%)	27	12
1	F	326/338 (96%)	310 (95%)	16 (5%)	31	15
All	All	1956/2028 (96%)	1857 (95%)	99 (5%)	29	13

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	14	GLU
1	A	47	GLU
1	A	65	LEU
1	A	68	ARG
1	A	72	LEU
1	A	84	LEU
1	A	92	GLN
1	A	179	GLU
1	A	186	ARG
1	A	213	GLU
1	A	219	LEU
1	A	226	LEU
1	A	252	ARG
1	A	285	ARG
1	A	289	LEU
1	A	301	LEU
1	B	65	LEU
1	B	68	ARG
1	B	72	LEU
1	B	84	LEU
1	B	92	GLN
1	B	105	SER
1	B	124	LEU
1	B	179	GLU
1	B	186	ARG
1	B	215	LYS
1	B	257	HIS
1	B	258	ARG
1	B	339	LEU
1	B	381	GLU
1	C	9	LYS

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Mol	Chain	Res	Type
1	C	47	GLU
1	C	68	ARG
1	C	72	LEU
1	C	84	LEU
1	C	92	GLN
1	C	104	THR
1	C	105	SER
1	C	124	LEU
1	C	179	GLU
1	C	215	LYS
1	C	226	LEU
1	C	252	ARG
1	C	283	THR
1	C	285	ARG
1	C	288	GLU
1	C	289	LEU
1	C	381	GLU
1	D	39	ARG
1	D	62	SER
1	D	65	LEU
1	D	68	ARG
1	D	72	LEU
1	D	84	LEU
1	D	92	GLN
1	D	105	SER
1	D	124	LEU
1	D	186	ARG
1	D	215	LYS
1	D	226	LEU
1	D	283	THR
1	D	289	LEU
1	D	301	LEU
1	D	339	LEU
1	E	10	HIS
1	E	65	LEU
1	E	68	ARG
1	E	84	LEU
1	E	92	GLN
1	E	105	SER
1	E	109	ASP
1	E	124	LEU
1	E	179	GLU

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Mol	Chain	Res	Type
1	E	186	ARG
1	E	202	GLN
1	E	219	LEU
1	E	252	ARG
1	E	258	ARG
1	E	285	ARG
1	E	289	LEU
1	E	301	LEU
1	E	380	LEU
1	F	9	LYS
1	F	21	ARG
1	F	62	SER
1	F	65	LEU
1	F	68	ARG
1	F	72	LEU
1	F	84	LEU
1	F	92	GLN
1	F	105	SER
1	F	109	ASP
1	F	124	LEU
1	F	179	GLU
1	F	186	ARG
1	F	257	HIS
1	F	301	LEU
1	F	339	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	81	ASN
1	A	202	GLN
1	A	330	GLN
1	A	377	GLN
1	B	81	ASN
1	B	127	HIS
1	B	202	GLN
1	B	281	HIS
1	B	330	GLN
1	B	377	GLN
1	C	81	ASN
1	C	127	HIS

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Mol	Chain	Res	Type
1	C	330	GLN
1	C	377	GLN
1	D	81	ASN
1	D	127	HIS
1	D	255	HIS
1	D	330	GLN
1	D	377	GLN
1	E	81	ASN
1	E	127	HIS
1	E	330	GLN
1	E	377	GLN
1	F	66	GLN
1	F	81	ASN
1	F	127	HIS
1	F	161	GLN
1	F	255	HIS
1	F	377	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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	FLC	A	1382	-	3,12,12	0.39	0	3,17,17	1.11	0
2	FLC	B	1382	-	3,12,12	0.99	0	3,17,17	0.87	0
2	FLC	C	1382	-	3,12,12	0.68	0	3,17,17	0.55	0
2	FLC	D	1382	-	3,12,12	0.63	0	3,17,17	0.25	0
2	FLC	E	1382	-	3,12,12	0.66	0	3,17,17	0.44	0
2	FLC	F	1382	-	3,12,12	0.59	0	3,17,17	0.58	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
2	FLC	A	1382	-	-	0/6/16/16	0/0/0/0
2	FLC	B	1382	-	-	0/6/16/16	0/0/0/0
2	FLC	C	1382	-	-	0/6/16/16	0/0/0/0
2	FLC	D	1382	-	-	0/6/16/16	0/0/0/0
2	FLC	E	1382	-	-	0/6/16/16	0/0/0/0
2	FLC	F	1382	-	-	0/6/16/16	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1382	FLC	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	381/397 (95%)	0.26	7 (1%) 71 80	14, 19, 28, 42	0
1	B	380/397 (95%)	0.19	3 (0%) 87 92	14, 19, 27, 39	0
1	C	381/397 (95%)	0.26	9 (2%) 62 72	14, 19, 28, 42	0
1	D	380/397 (95%)	0.20	6 (1%) 74 83	14, 19, 27, 39	0
1	E	380/397 (95%)	0.24	9 (2%) 62 72	14, 19, 28, 44	0
1	F	380/397 (95%)	0.17	4 (1%) 82 88	14, 18, 27, 39	0
All	All	2282/2382 (95%)	0.22	38 (1%) 73 81	14, 19, 28, 44	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	ALA	5.5
1	C	216	ALA	5.0
1	C	10	HIS	4.8
1	C	219	LEU	4.5
1	B	258	ARG	4.4
1	F	258	ARG	4.4
1	D	258	ARG	4.4
1	A	10	HIS	4.2
1	A	219	LEU	3.5
1	E	216	ALA	3.5
1	F	247	GLN	3.2
1	E	12	HIS	3.1
1	D	247	GLN	2.9
1	B	247	GLN	2.9
1	F	259	VAL	2.6
1	E	380	LEU	2.6
1	A	9	LYS	2.5
1	E	10	HIS	2.5
1	C	217	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	257	HIS	2.4
1	E	219	LEU	2.4
1	D	246	GLY	2.4
1	A	217	HIS	2.4
1	E	9	LYS	2.4
1	C	214	GLY	2.3
1	C	381	GLU	2.3
1	D	181	GLU	2.2
1	E	214	GLY	2.2
1	C	213	GLU	2.2
1	D	259	VAL	2.2
1	A	380	LEU	2.2
1	B	257	HIS	2.2
1	C	103	ILE	2.2
1	E	55	ARG	2.1
1	C	9	LYS	2.1
1	D	257	HIS	2.1
1	A	381	GLU	2.1
1	E	104	THR	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 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
2	FLC	F	1382	13/13	0.89	0.18	3.32	30,36,40,41	0
2	FLC	D	1382	13/13	0.93	0.19	3.11	29,35,40,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	FLC	A	1382	13/13	0.82	0.21	2.19	29,38,47,51	0
2	FLC	B	1382	13/13	0.89	0.15	2.11	30,37,41,42	0
2	FLC	E	1382	13/13	0.90	0.20	1.24	26,34,41,42	0
2	FLC	C	1382	13/13	0.89	0.18	0.62	26,34,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.