



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:51 PM GMT

PDB ID : 4Q31
Title : The crystal structure of cystathione gamma lyase (CalE6) from *Micromonospora echinospora*
Authors : Tan, K.; Bigelow, L.; Jedrzejczak, R.; Babnigg, G.; Bingman, C.A.; Yen-namalli, R.M.; Singh, S.; Kharel, M.K.; Thorson, J.S.; Phillips Jr., G.N.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG); Enzyme Discovery for Natural Product Biosynthesis (NatPro)
Deposited on : 2014-04-10
Resolution : 2.10 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

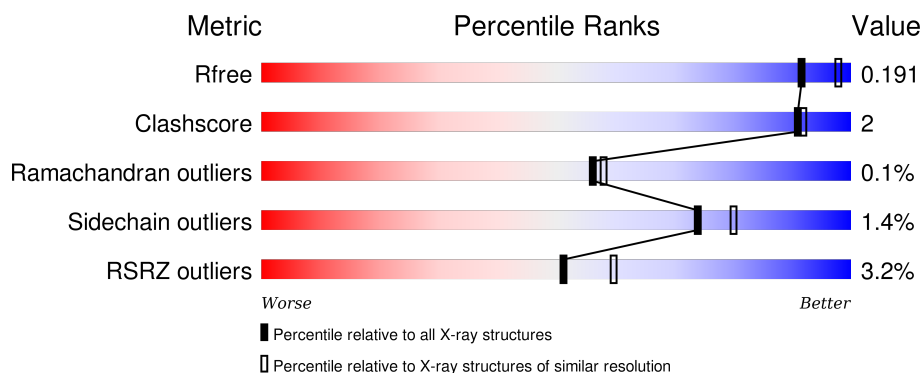
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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.



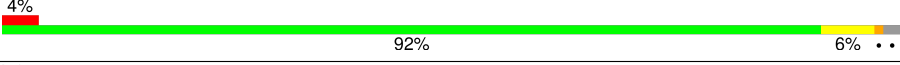
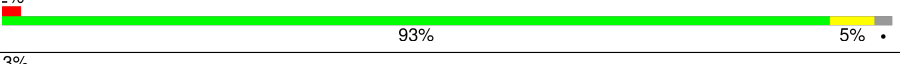
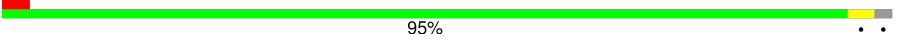
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	384	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>•</div> </div> </div>
1	B	384	<div> <div>2%</div> <div> <div></div> <div>93%</div> <div>5%</div> <div>••</div> </div> </div>
1	C	384	<div> <div>4%</div> <div> <div></div> <div>91%</div> <div>8%</div> <div>••</div> </div> </div>
1	D	384	<div> <div>4%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>•</div> </div> </div>
1	E	384	<div> <div>4%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>•</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	384	
1	G	384	
1	H	384	

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	A	407	-	-	-	X
4	GOL	B	403	-	-	-	X
4	GOL	B	404	-	-	-	X
4	GOL	B	405	-	-	-	X
4	GOL	B	406	-	-	-	X
4	GOL	B	407	-	-	-	X
4	GOL	D	404	-	-	-	X
4	GOL	D	405	-	-	-	X
4	GOL	D	407	-	-	X	-
4	GOL	E	403	-	-	-	X
4	GOL	E	404	-	-	-	X
4	GOL	E	405	-	-	-	X
4	GOL	E	406	-	-	-	X
4	GOL	F	403	-	-	-	X
4	GOL	F	404	-	-	-	X
4	GOL	F	405	-	-	-	X
4	GOL	G	405	-	-	-	X
5	FMT	F	406	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 24719 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 cystathione gamma lyase CalE6.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	P	S	Se	0	1	0
			2844	1778	526	530	1	5	4			
1	B	378	Total	C	N	O	P	S	Se	0	1	0
			2851	1781	528	532	1	5	4			
1	C	379	Total	C	N	O	P	S	Se	0	0	0
			2827	1769	523	526	1	5	3			
1	D	376	Total	C	N	O	P	S	Se	0	1	0
			2817	1764	519	524	1	5	4			
1	E	376	Total	C	N	O	P	S	Se	0	0	0
			2813	1764	512	528	1	5	3			
1	F	377	Total	C	N	O	P	S	Se	0	0	0
			2812	1760	518	525	1	5	3			
1	G	378	Total	C	N	O	P	S	Se	0	0	0
			2814	1761	516	528	1	5	3			
1	H	378	Total	C	N	O	P	S	Se	0	1	0
			2828	1770	520	529	1	5	3			

There are 32 discrepancies between the modelled and reference sequences:

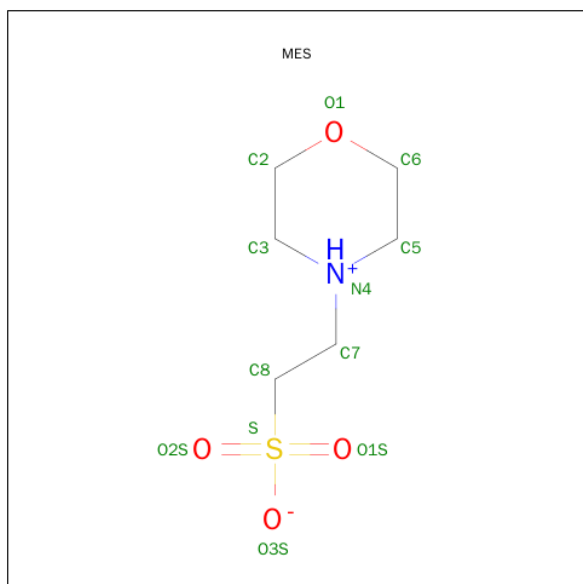
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
A	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
A	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
A	7	GLY	ASP	engineered mutation	UNP Q8KNG3
B	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
B	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
B	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
B	7	GLY	ASP	engineered mutation	UNP Q8KNG3
C	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
C	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
C	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
C	7	GLY	ASP	engineered mutation	UNP Q8KNG3
D	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
D	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
D	7	GLY	ASP	engineered mutation	UNP Q8KNG3
E	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
E	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
E	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
E	7	GLY	ASP	engineered mutation	UNP Q8KNG3
F	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
F	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
F	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
F	7	GLY	ASP	engineered mutation	UNP Q8KNG3
G	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
G	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
G	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
G	7	GLY	ASP	engineered mutation	UNP Q8KNG3
H	-2	SER	-	EXPRESSION TAG	UNP Q8KNG3
H	-1	ASN	-	EXPRESSION TAG	UNP Q8KNG3
H	0	ALA	-	EXPRESSION TAG	UNP Q8KNG3
H	7	GLY	ASP	engineered mutation	UNP Q8KNG3

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	E	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	F	1	Total	C	N	O	S	0	0
			11	6	1	3	1		
2	G	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	H	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Cl	0	0
			1	1		
3	D	2	Total	Cl	0	0
			2	2		
3	E	1	Total	Cl	0	0
			1	1		
3	H	2	Total	Cl	0	0
			2	2		
3	B	1	Total	Cl	0	0
			1	1		
3	C	2	Total	Cl	0	0
			2	2		
3	A	2	Total	Cl	0	0
			2	2		
3	F	1	Total	Cl	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



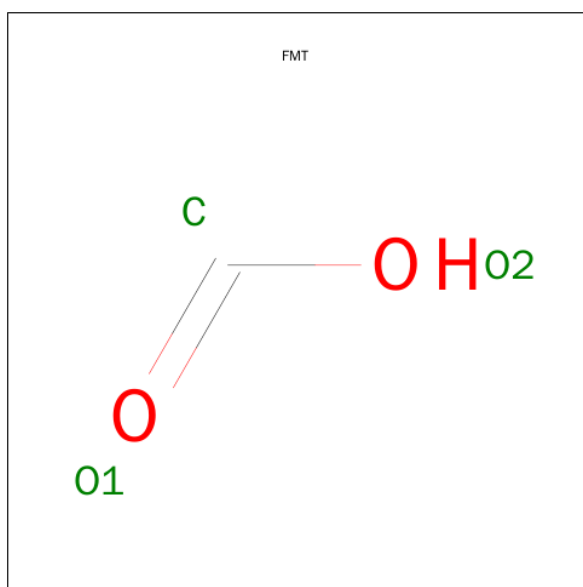
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			3	1	2		
5	C	1	Total	C	O	0	0
			3	1	2		
5	F	1	Total	C	O	0	0
			3	1	2		
5	G	1	Total	C	O	0	0
			3	1	2		
5	H	1	Total	C	O	0	0
			3	1	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	243	Total	O	0	0
			243	243		
6	B	228	Total	O	0	0
			228	228		
6	C	230	Total	O	0	0
			230	230		
6	D	215	Total	O	0	0
			215	215		
6	E	250	Total	O	0	0
			250	250		
6	F	204	Total	O	0	0
			204	204		
6	G	215	Total	O	0	0
			215	215		

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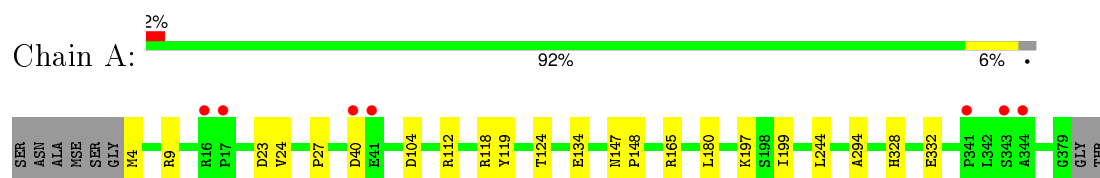
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	H	238	Total 238	O 238	0	0

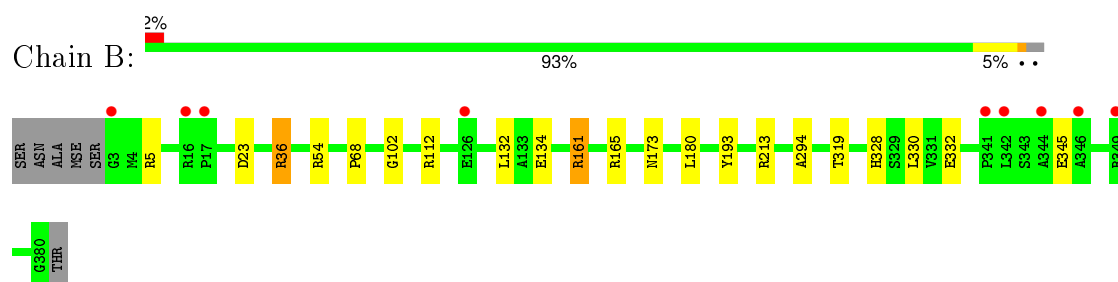
3 Residue-property plots [i](#)

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.

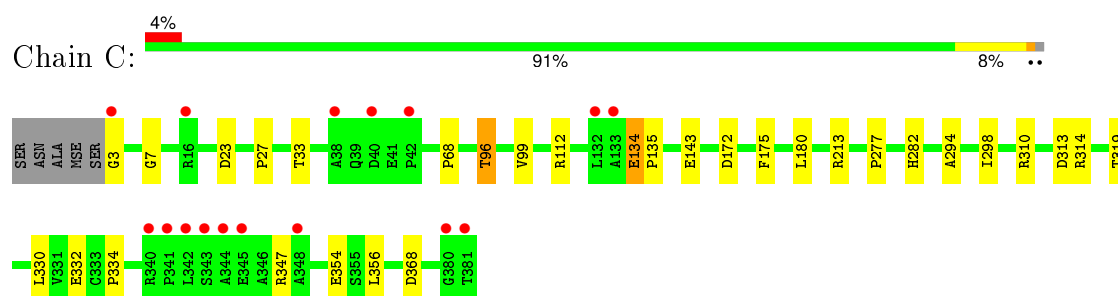
- Molecule 1: cystathione gamma lyase CalE6



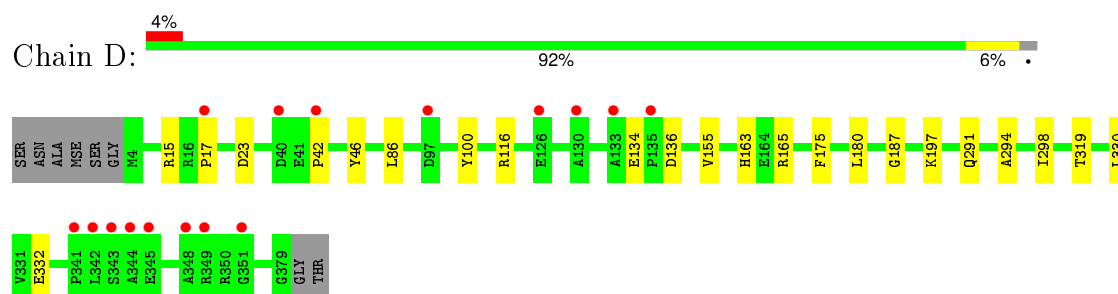
- Molecule 1: cystathione gamma lyase CalE6



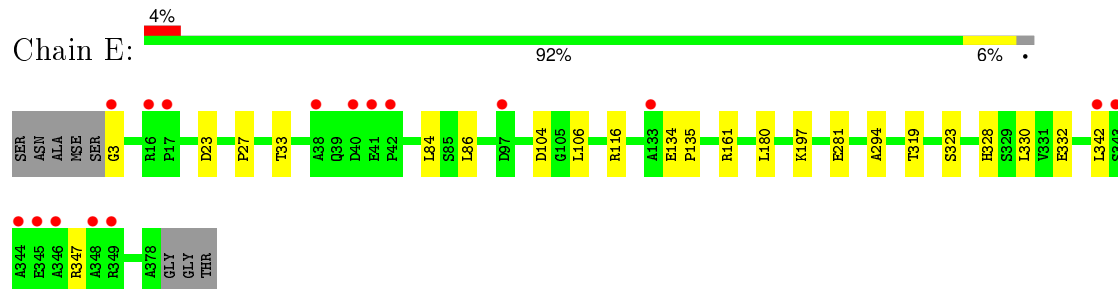
- Molecule 1: cystathione gamma lyase CalE6



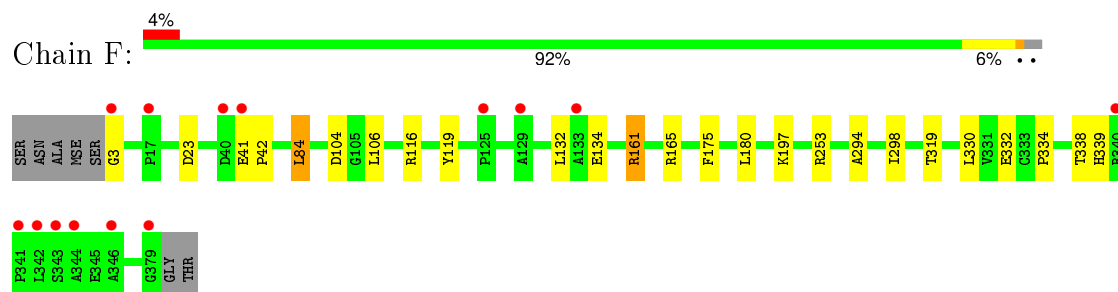
- Molecule 1: cystathione gamma lyase CalE6



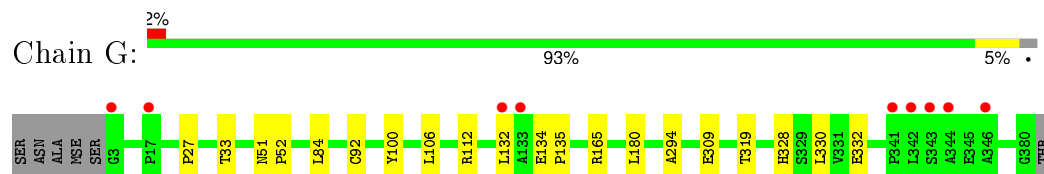
- Molecule 1: cystathione gamma lyase CalE6



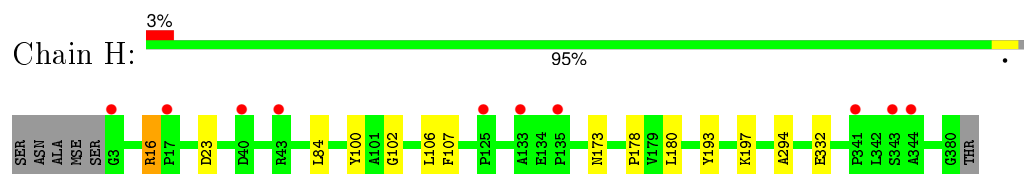
- Molecule 1: cystathione gamma lyase CalE6



- Molecule 1: cystathione gamma lyase CalE6



- Molecule 1: cystathione gamma lyase CalE6



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	146.85Å 146.98Å 349.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.97 – 2.10 33.97 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (33.97-2.10) 99.9 (33.97-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.152 , 0.191 0.155 , 0.191	Depositor DCC
R_{free} test set	11002 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.0	EDS
Estimated twinning fraction	0.055 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 219086 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	24719	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.26% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, FMT, LLP, MES, CL

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	0.36	0/2876	0.51	0/3916
1	B	0.37	0/2882	0.52	0/3922
1	C	0.37	0/2855	0.52	0/3890
1	D	0.35	0/2849	0.51	0/3884
1	E	0.37	0/2841	0.50	0/3871
1	F	0.36	0/2840	0.51	0/3871
1	G	0.35	0/2842	0.51	0/3874
1	H	0.36	0/2859	0.51	0/3894
All	All	0.36	0/22844	0.51	0/31122

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 [i](#)

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	2844	0	2843	10	0
1	B	2851	0	2848	14	0
1	C	2827	0	2812	20	0
1	D	2817	0	2800	15	0
1	E	2813	0	2791	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2812	0	2788	15	0
1	G	2814	0	2788	10	0
1	H	2828	0	2813	9	0
2	A	12	0	12	0	0
2	B	12	0	12	1	0
2	C	12	0	12	1	0
2	D	12	0	12	1	0
2	E	12	0	12	0	0
2	F	11	0	12	0	0
2	G	12	0	12	2	0
2	H	12	0	12	1	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	1	0
3	D	2	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	2	0	0	0	0
4	A	24	0	32	1	0
4	B	36	0	48	1	0
4	C	6	0	8	0	0
4	D	24	0	32	5	0
4	E	24	0	32	1	0
4	F	18	0	24	1	0
4	G	30	0	40	1	0
4	H	6	0	8	0	0
5	A	3	0	1	0	0
5	C	3	0	1	0	0
5	F	3	0	1	0	0
5	G	3	0	1	0	0
5	H	3	0	1	0	0
6	A	243	0	0	0	0
6	B	228	0	0	2	0
6	C	230	0	0	2	0
6	D	215	0	0	0	0
6	E	250	0	0	2	0
6	F	204	0	0	1	0
6	G	215	0	0	0	0
6	H	238	0	0	1	0
All	All	24719	0	22808	101	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 2.

All (101) 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:86:LEU:HB3	4:E:405:GOL:H11	1.75	0.67
1:F:134:GLU:O	1:F:165:ARG:NH1	2.28	0.67
1:C:3:GLY:N	6:C:644:HOH:O	2.29	0.65
1:F:338:THR:HG23	1:F:339:HIS:ND1	2.12	0.64
1:F:253:ARG:HD2	4:F:404:GOL:H11	1.83	0.61
1:B:132:LEU:HD12	1:B:161:ARG:HB3	1.82	0.60
1:A:4[B]:MSE:HE3	1:A:9:ARG:HG2	1.83	0.60
1:A:180:LEU:HD22	1:A:294:ALA:HB3	1.84	0.60
4:B:403:GOL:H2	1:C:7:GLY:HA3	1.84	0.59
1:D:17:PRO:HA	4:D:407:GOL:H2	1.84	0.59
1:F:116:ARG:NH1	1:F:134:GLU:OE1	2.38	0.57
4:G:405:GOL:H31	1:H:102:GLY:HA2	1.86	0.57
1:A:199:ILE:HG23	1:A:244:LEU:HD13	1.86	0.57
1:F:334:PRO:HA	1:F:338:THR:HG22	1.86	0.56
1:G:319:THR:HB	1:G:330:LEU:HD23	1.87	0.56
1:C:310:ARG:NH1	1:C:313:ASP:OD2	2.35	0.56
1:D:134:GLU:O	1:D:165:ARG:NH1	2.39	0.56
1:B:112:ARG:NH2	6:B:647:HOH:O	2.33	0.56
2:C:401:MES:H61	1:D:46:TYR:CD2	2.43	0.53
1:H:180:LEU:HD22	1:H:294:ALA:HB3	1.90	0.53
1:E:342:LEU:O	1:E:347:ARG:NH1	2.43	0.52
1:B:180:LEU:HD22	1:B:294:ALA:HB3	1.92	0.52
1:A:104:ASP:OD1	1:A:119:TYR:OH	2.23	0.52
1:D:15:ARG:NH2	4:D:407:GOL:H32	2.25	0.51
1:D:15:ARG:HH22	4:D:407:GOL:H32	1.74	0.51
1:B:134:GLU:O	1:B:165:ARG:NH1	2.43	0.51
1:F:132:LEU:O	1:F:165:ARG:NE	2.42	0.51
1:F:180:LEU:HD22	1:F:294:ALA:HB3	1.93	0.50
1:C:314:ARG:HD3	6:C:696:HOH:O	2.11	0.50
1:C:143:GLU:HG3	1:C:172:ASP:HB3	1.94	0.50
1:G:84:LEU:HD12	1:G:106:LEU:HB3	1.94	0.50
1:G:180:LEU:HD22	1:G:294:ALA:HB3	1.93	0.49
1:D:180:LEU:HD22	1:D:294:ALA:HB3	1.94	0.49
1:C:112:ARG:HD3	3:C:403:CL:CL	2.49	0.49
1:D:319:THR:HB	1:D:330:LEU:HD23	1.95	0.49
1:C:319:THR:HB	1:C:330:LEU:HD23	1.95	0.48
1:F:319:THR:HB	1:F:330:LEU:HD23	1.96	0.48
1:G:33:THR:HG23	6:H:626:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:319:THR:HB	1:E:330:LEU:HD23	1.96	0.48
1:E:3:GLY:N	6:E:664:HOH:O	2.47	0.47
1:H:178:PRO:HD3	1:H:193:TYR:CE1	2.50	0.47
1:H:100:TYR:CE1	2:H:401:MES:H81	2.48	0.47
1:C:180:LEU:HD22	1:C:294:ALA:HB3	1.95	0.47
1:A:24:VAL:HB	1:C:33:THR:HG23	1.96	0.47
1:G:100:TYR:CE1	2:G:401:MES:H81	2.49	0.47
4:A:407:GOL:H12	1:B:102:GLY:HA2	1.97	0.47
1:F:197:LLP:NZ	1:F:197:LLP:O3	2.47	0.47
1:B:345:GLU:OE1	1:B:345:GLU:N	2.46	0.47
1:B:5:ARG:HG2	1:C:368:ASP:CG	2.36	0.46
1:E:116:ARG:NE	6:E:722:HOH:O	2.37	0.46
1:D:175:PHE:CZ	1:D:298:ILE:HB	2.51	0.46
1:F:175:PHE:CZ	1:F:298:ILE:HB	2.51	0.46
1:B:54:ARG:NH1	6:B:526:HOH:O	2.46	0.45
1:C:175:PHE:CZ	1:C:298:ILE:HB	2.51	0.45
1:F:132:LEU:HD23	1:F:161:ARG:HB3	1.98	0.45
1:G:134:GLU:HG3	1:G:135:PRO:HD2	1.97	0.45
1:C:277:PRO:HA	1:C:282:HIS:CD2	2.52	0.44
1:B:36:ARG:HE	1:B:36:ARG:HB2	1.43	0.44
1:A:134:GLU:O	1:A:165:ARG:NH1	2.49	0.44
1:D:100:TYR:CE1	2:D:401:MES:H81	2.53	0.44
1:H:197:LLP:NZ	1:H:197:LLP:O3	2.51	0.44
1:B:5:ARG:HG2	1:C:368:ASP:OD2	2.18	0.44
1:D:17:PRO:HA	4:D:407:GOL:C2	2.46	0.43
1:F:3:GLY:N	6:F:666:HOH:O	2.49	0.43
1:C:68:PRO:HG2	1:C:213:ARG:HA	2.00	0.43
1:D:163:HIS:HE1	1:D:187:GLY:O	2.02	0.43
1:E:84:LEU:HD12	1:E:106:LEU:HB3	2.01	0.43
1:H:84:LEU:HD11	1:H:107:PHE:HD1	1.83	0.43
1:D:116:ARG:HH12	1:D:136:ASP:H	1.66	0.43
1:H:84:LEU:HD12	1:H:106:LEU:HB3	2.01	0.43
1:E:27:PRO:HB2	1:G:27:PRO:HB2	2.01	0.43
1:H:173:ASN:HB3	1:H:193:TYR:CE1	2.54	0.43
1:C:134:GLU:HG3	1:C:135:PRO:HD2	2.01	0.43
1:F:84:LEU:HD12	1:F:106:LEU:HB3	2.00	0.43
1:E:180:LEU:HD22	1:E:294:ALA:HB3	2.00	0.42
1:E:161:ARG:HA	1:E:161:ARG:HD3	1.90	0.42
1:D:197:LLP:O3	1:D:197:LLP:NZ	2.52	0.42
1:C:96:THR:HG23	1:C:99:VAL:HG13	2.01	0.42
1:D:86:LEU:HD13	4:D:406:GOL:H32	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:334:PRO:HD2	1:C:356:LEU:O	2.19	0.42
1:F:104:ASP:OD1	1:F:119:TYR:OH	2.30	0.42
1:G:51:ASN:HA	1:G:52:PRO:HD3	1.93	0.42
1:D:155:VAL:N	1:D:291:GLN:OE1	2.39	0.41
2:G:401:MES:H82	2:G:401:MES:H31	1.71	0.41
1:G:92:CYS:SG	1:G:134:GLU:HG2	2.60	0.41
1:C:332:GLU:OE2	1:C:334:PRO:HA	2.20	0.41
1:B:132:LEU:CD1	1:B:161:ARG:HB3	2.50	0.41
1:G:132:LEU:O	1:G:165:ARG:NH1	2.53	0.41
1:E:134:GLU:HA	1:E:135:PRO:HD3	1.95	0.41
1:B:68:PRO:HG2	1:B:213:ARG:HA	2.01	0.41
1:A:197:LLP:O3	1:A:197:LLP:NZ	2.53	0.41
1:A:112:ARG:NH2	1:E:135:PRO:HG2	2.35	0.41
2:B:401:MES:H82	2:B:401:MES:H32	1.79	0.41
1:E:197:LLP:O3	1:E:197:LLP:NZ	2.53	0.41
1:C:347:ARG:NH2	1:C:354:GLU:OE2	2.53	0.41
1:B:319:THR:HB	1:B:330:LEU:HD23	2.03	0.40
1:F:41:GLU:HA	1:F:42:PRO:HD3	1.85	0.40
1:H:16:ARG:HH11	1:H:16:ARG:HA	1.86	0.40
1:A:27:PRO:HB2	1:C:27:PRO:HB2	2.02	0.40
1:A:147:ASN:HA	1:A:148:PRO:HA	1.88	0.40
1:B:173:ASN:HB3	1:B:193:TYR:CE1	2.56	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	373/384 (97%)	366 (98%)	7 (2%)	0	100	100
1	B	376/384 (98%)	364 (97%)	12 (3%)	0	100	100
1	C	376/384 (98%)	367 (98%)	9 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	373/384 (97%)	363 (97%)	9 (2%)	1 (0%)	46	45
1	E	373/384 (97%)	366 (98%)	6 (2%)	1 (0%)	46	45
1	F	374/384 (97%)	367 (98%)	7 (2%)	0	100	100
1	G	375/384 (98%)	364 (97%)	11 (3%)	0	100	100
1	H	376/384 (98%)	365 (97%)	11 (3%)	0	100	100
All	All	2996/3072 (98%)	2922 (98%)	72 (2%)	2 (0%)	56	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	323	SER
1	D	42	PRO

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	292/293 (100%)	286 (98%)	6 (2%)	61	66
1	B	292/293 (100%)	287 (98%)	5 (2%)	68	74
1	C	286/293 (98%)	283 (99%)	3 (1%)	82	87
1	D	286/293 (98%)	284 (99%)	2 (1%)	88	92
1	E	286/293 (98%)	280 (98%)	6 (2%)	61	66
1	F	284/293 (97%)	280 (99%)	4 (1%)	74	80
1	G	285/293 (97%)	281 (99%)	4 (1%)	74	80
1	H	287/293 (98%)	284 (99%)	3 (1%)	82	87
All	All	2298/2344 (98%)	2265 (99%)	33 (1%)	74	80

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ASP

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Mol	Chain	Res	Type
1	A	40	ASP
1	A	118	ARG
1	A	124	THR
1	A	328	HIS
1	A	332	GLU
1	B	23	ASP
1	B	36	ARG
1	B	161	ARG
1	B	328	HIS
1	B	332	GLU
1	C	23	ASP
1	C	96	THR
1	C	134	GLU
1	D	23	ASP
1	D	332	GLU
1	E	23	ASP
1	E	33	THR
1	E	104	ASP
1	E	281	GLU
1	E	328	HIS
1	E	332	GLU
1	F	23	ASP
1	F	84	LEU
1	F	161	ARG
1	F	332	GLU
1	G	112	ARG
1	G	309	GLU
1	G	328	HIS
1	G	332	GLU
1	H	16	ARG
1	H	23	ASP
1	H	332	GLU

Some 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 ⓘ

8 non-standard protein/DNA/RNA residues 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$
1	LLP	A	197	1	23,24,25	3.36	7 (30%)	28,32,34	1.38	6 (21%)
1	LLP	B	197	1	23,24,25	3.40	7 (30%)	28,32,34	1.31	6 (21%)
1	LLP	C	197	1	23,24,25	3.49	7 (30%)	28,32,34	1.30	4 (14%)
1	LLP	D	197	1	23,24,25	3.35	7 (30%)	28,32,34	1.27	6 (21%)
1	LLP	E	197	1	23,24,25	3.30	7 (30%)	28,32,34	1.25	4 (14%)
1	LLP	F	197	1	23,24,25	3.43	7 (30%)	28,32,34	1.30	5 (17%)
1	LLP	G	197	1	23,24,25	3.35	7 (30%)	28,32,34	1.33	6 (21%)
1	LLP	H	197	1	23,24,25	3.35	7 (30%)	28,32,34	1.52	6 (21%)

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
1	LLP	A	197	1	-	0/15/17/19	0/1/1/1
1	LLP	B	197	1	-	0/15/17/19	0/1/1/1
1	LLP	C	197	1	-	0/15/17/19	0/1/1/1
1	LLP	D	197	1	-	0/15/17/19	0/1/1/1
1	LLP	E	197	1	-	0/15/17/19	0/1/1/1
1	LLP	F	197	1	-	0/15/17/19	0/1/1/1
1	LLP	G	197	1	-	0/15/17/19	0/1/1/1
1	LLP	H	197	1	-	0/15/17/19	0/1/1/1

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	197	LLP	C4-C5	-6.32	1.33	1.42
1	F	197	LLP	C4-C5	-5.99	1.33	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	197	LLP	C4-C5	-5.95	1.33	1.42
1	D	197	LLP	C4-C5	-5.85	1.34	1.42
1	B	197	LLP	C4-C5	-5.76	1.34	1.42
1	B	197	LLP	C3-C2	-5.61	1.36	1.40
1	D	197	LLP	C3-C2	-5.55	1.36	1.40
1	F	197	LLP	C3-C2	-5.47	1.37	1.40
1	E	197	LLP	C4-C5	-5.37	1.34	1.42
1	G	197	LLP	C4-C5	-5.24	1.34	1.42
1	A	197	LLP	C3-C2	-5.16	1.37	1.40
1	H	197	LLP	C4-C5	-4.93	1.35	1.42
1	G	197	LLP	C3-C2	-4.81	1.37	1.40
1	C	197	LLP	C3-C2	-4.69	1.37	1.40
1	H	197	LLP	C3-C2	-4.49	1.37	1.40
1	E	197	LLP	C3-C2	-4.30	1.37	1.40
1	A	197	LLP	C4-C4'	5.28	1.55	1.46
1	E	197	LLP	C2-N1	5.32	1.45	1.34
1	B	197	LLP	C4-C4'	5.33	1.56	1.46
1	D	197	LLP	C4-C4'	5.34	1.56	1.46
1	C	197	LLP	C2-N1	5.40	1.45	1.34
1	G	197	LLP	C2-N1	5.42	1.45	1.34
1	F	197	LLP	C4-C4'	5.43	1.56	1.46
1	A	197	LLP	C2-N1	5.50	1.45	1.34
1	H	197	LLP	C2-N1	5.52	1.45	1.34
1	B	197	LLP	C2-N1	5.53	1.45	1.34
1	E	197	LLP	C4-C4'	5.60	1.56	1.46
1	F	197	LLP	C2-N1	5.61	1.46	1.34
1	A	197	LLP	C6-C5	5.61	1.50	1.37
1	D	197	LLP	C2-N1	5.62	1.46	1.34
1	G	197	LLP	C4-C4'	5.64	1.56	1.46
1	C	197	LLP	C4-C4'	5.68	1.56	1.46
1	E	197	LLP	C6-C5	5.69	1.50	1.37
1	H	197	LLP	C6-C5	5.71	1.50	1.37
1	B	197	LLP	C6-C5	5.73	1.50	1.37
1	H	197	LLP	C4-C4'	5.79	1.56	1.46
1	G	197	LLP	C6-C5	5.82	1.50	1.37
1	C	197	LLP	C6-C5	5.84	1.50	1.37
1	D	197	LLP	C6-C5	5.87	1.50	1.37
1	F	197	LLP	C6-C5	5.93	1.50	1.37
1	D	197	LLP	C4'-NZ	6.01	1.45	1.27
1	A	197	LLP	C4'-NZ	6.18	1.46	1.27
1	E	197	LLP	C4'-NZ	6.21	1.46	1.27
1	F	197	LLP	C4'-NZ	6.25	1.46	1.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	197	LLP	C4'-NZ	6.27	1.46	1.27
1	C	197	LLP	C4'-NZ	6.28	1.46	1.27
1	G	197	LLP	C4'-NZ	6.43	1.46	1.27
1	H	197	LLP	C4'-NZ	6.55	1.47	1.27
1	D	197	LLP	C4-C3	6.90	1.49	1.40
1	F	197	LLP	C4-C3	7.21	1.49	1.40
1	B	197	LLP	C4-C3	7.28	1.49	1.40
1	H	197	LLP	C4-C3	7.30	1.49	1.40
1	A	197	LLP	C4-C3	7.34	1.50	1.40
1	G	197	LLP	C4-C3	7.34	1.50	1.40
1	E	197	LLP	C4-C3	7.41	1.50	1.40
1	C	197	LLP	C4-C3	8.02	1.50	1.40

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	LLP	O-C-CA	-2.75	118.32	125.49
1	A	197	LLP	CE-NZ-C4'	-2.61	111.42	118.97
1	D	197	LLP	C4-C4'-NZ	-2.59	110.67	125.06
1	B	197	LLP	CE-NZ-C4'	-2.55	111.59	118.97
1	C	197	LLP	C4-C4'-NZ	-2.54	110.93	125.06
1	F	197	LLP	O-C-CA	-2.51	118.95	125.49
1	F	197	LLP	C4-C4'-NZ	-2.51	111.11	125.06
1	E	197	LLP	C4-C4'-NZ	-2.45	111.44	125.06
1	G	197	LLP	O-C-CA	-2.44	119.14	125.49
1	E	197	LLP	O-C-CA	-2.44	119.15	125.49
1	H	197	LLP	O-C-CA	-2.43	119.16	125.49
1	B	197	LLP	O-C-CA	-2.43	119.16	125.49
1	D	197	LLP	C5-C6-N1	-2.43	119.65	123.86
1	F	197	LLP	C5-C6-N1	-2.42	119.66	123.86
1	C	197	LLP	O-C-CA	-2.40	119.23	125.49
1	G	197	LLP	CE-NZ-C4'	-2.38	112.09	118.97
1	D	197	LLP	O-C-CA	-2.37	119.31	125.49
1	A	197	LLP	C5-C6-N1	-2.32	119.83	123.86
1	H	197	LLP	C5-C6-N1	-2.30	119.87	123.86
1	H	197	LLP	CE-NZ-C4'	-2.24	112.51	118.97
1	A	197	LLP	C4-C4'-NZ	-2.19	112.85	125.06
1	B	197	LLP	C4-C4'-NZ	-2.19	112.85	125.06
1	B	197	LLP	C5-C6-N1	-2.19	120.06	123.86
1	D	197	LLP	CE-NZ-C4'	-2.19	112.65	118.97
1	G	197	LLP	C4-C4'-NZ	-2.17	112.98	125.06
1	H	197	LLP	C5'-C5-C6	-2.14	115.23	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	197	LLP	C5-C6-N1	-2.01	120.37	123.86
1	B	197	LLP	OP2-P-OP4	2.06	112.51	106.56
1	C	197	LLP	OP2-P-OP4	2.21	112.93	106.56
1	D	197	LLP	OP3-P-OP4	2.22	112.95	106.56
1	D	197	LLP	OP2-P-OP4	2.28	113.14	106.56
1	E	197	LLP	OP2-P-OP4	2.28	113.14	106.56
1	G	197	LLP	OP2-P-OP4	2.30	113.18	106.56
1	F	197	LLP	OP2-P-OP4	2.40	113.49	106.56
1	A	197	LLP	OP3-P-OP4	2.54	113.89	106.56
1	G	197	LLP	OP3-P-OP4	2.64	114.16	106.56
1	A	197	LLP	OP2-P-OP4	2.70	114.33	106.56
1	E	197	LLP	OP3-P-OP4	2.81	114.66	106.56
1	F	197	LLP	OP3-P-OP4	2.86	114.80	106.56
1	H	197	LLP	OP3-P-OP4	2.92	114.96	106.56
1	C	197	LLP	OP3-P-OP4	2.94	115.04	106.56
1	H	197	LLP	OP2-P-OP4	3.02	115.25	106.56
1	B	197	LLP	OP3-P-OP4	3.13	115.58	106.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	197	LLP	1	0
1	D	197	LLP	1	0
1	E	197	LLP	1	0
1	F	197	LLP	1	0
1	H	197	LLP	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 53 ligands modelled in this entry, 12 are monoatomic - leaving 41 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
2	MES	A	401	-	11,12,12	0.67	0	14,16,16	2.62	3 (21%)
4	GOL	A	404	-	5,5,5	0.30	0	5,5,5	0.24	0
4	GOL	A	405	-	5,5,5	0.29	0	5,5,5	0.22	0
4	GOL	A	406	-	5,5,5	0.29	0	5,5,5	0.38	0
4	GOL	A	407	-	5,5,5	0.42	0	5,5,5	0.46	0
5	FMT	A	408	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	B	401	-	11,12,12	0.73	0	14,16,16	2.36	6 (42%)
4	GOL	B	403	-	5,5,5	0.34	0	5,5,5	0.24	0
4	GOL	B	404	-	5,5,5	0.33	0	5,5,5	0.43	0
4	GOL	B	405	-	5,5,5	0.38	0	5,5,5	0.33	0
4	GOL	B	406	-	5,5,5	0.34	0	5,5,5	0.67	0
4	GOL	B	407	-	5,5,5	0.33	0	5,5,5	0.42	0
4	GOL	B	408	-	5,5,5	0.34	0	5,5,5	0.32	0
2	MES	C	401	-	11,12,12	0.73	0	14,16,16	2.42	4 (28%)
4	GOL	C	404	-	5,5,5	0.29	0	5,5,5	0.36	0
5	FMT	C	405	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	D	401	-	11,12,12	0.76	0	14,16,16	2.12	3 (21%)
4	GOL	D	404	-	5,5,5	0.35	0	5,5,5	0.32	0
4	GOL	D	405	-	5,5,5	0.35	0	5,5,5	0.32	0
4	GOL	D	406	-	5,5,5	0.34	0	5,5,5	0.28	0
4	GOL	D	407	-	5,5,5	0.41	0	5,5,5	0.45	0
2	MES	E	401	-	11,12,12	0.72	0	14,16,16	2.31	3 (21%)
4	GOL	E	403	-	5,5,5	0.35	0	5,5,5	0.25	0
4	GOL	E	404	-	5,5,5	0.36	0	5,5,5	0.15	0
4	GOL	E	405	-	5,5,5	0.41	0	5,5,5	0.25	0
4	GOL	E	406	-	5,5,5	0.38	0	5,5,5	0.15	0
2	MES	F	401	-	9,11,12	0.70	0	11,13,16	2.03	2 (18%)
4	GOL	F	403	-	5,5,5	0.37	0	5,5,5	0.57	0
4	GOL	F	404	-	5,5,5	0.41	0	5,5,5	0.35	0
4	GOL	F	405	-	5,5,5	0.31	0	5,5,5	0.36	0
5	FMT	F	406	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	G	401	-	11,12,12	0.70	0	14,16,16	2.51	3 (21%)
4	GOL	G	403	-	5,5,5	0.38	0	5,5,5	0.35	0
4	GOL	G	404	-	5,5,5	0.29	0	5,5,5	0.37	0
4	GOL	G	405	-	5,5,5	0.39	0	5,5,5	0.09	0
4	GOL	G	406	-	5,5,5	0.32	0	5,5,5	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	G	407	-	5,5,5	0.30	0	5,5,5	0.33	0
5	FMT	G	408	-	0,2,2	0.00	-	0,1,1	0.00	-
2	MES	H	401	-	11,12,12	0.64	0	14,16,16	2.18	3 (21%)
4	GOL	H	404	-	5,5,5	0.35	0	5,5,5	0.31	0
5	FMT	H	405	-	0,2,2	0.00	-	0,1,1	0.00	-

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	MES	A	401	-	-	0/6/14/14	0/1/1/1
4	GOL	A	404	-	-	0/4/4/4	0/0/0/0
4	GOL	A	405	-	-	0/4/4/4	0/0/0/0
4	GOL	A	406	-	-	0/4/4/4	0/0/0/0
4	GOL	A	407	-	-	0/4/4/4	0/0/0/0
5	FMT	A	408	-	-	0/0/0/0	0/0/0/0
2	MES	B	401	-	-	0/6/14/14	0/1/1/1
4	GOL	B	403	-	-	0/4/4/4	0/0/0/0
4	GOL	B	404	-	-	0/4/4/4	0/0/0/0
4	GOL	B	405	-	-	0/4/4/4	0/0/0/0
4	GOL	B	406	-	-	0/4/4/4	0/0/0/0
4	GOL	B	407	-	-	0/4/4/4	0/0/0/0
4	GOL	B	408	-	-	0/4/4/4	0/0/0/0
2	MES	C	401	-	-	0/6/14/14	0/1/1/1
4	GOL	C	404	-	-	0/4/4/4	0/0/0/0
5	FMT	C	405	-	-	0/0/0/0	0/0/0/0
2	MES	D	401	-	-	0/6/14/14	0/1/1/1
4	GOL	D	404	-	-	0/4/4/4	0/0/0/0
4	GOL	D	405	-	-	0/4/4/4	0/0/0/0
4	GOL	D	406	-	-	0/4/4/4	0/0/0/0
4	GOL	D	407	-	-	0/4/4/4	0/0/0/0
2	MES	E	401	-	-	0/6/14/14	0/1/1/1
4	GOL	E	403	-	-	0/4/4/4	0/0/0/0
4	GOL	E	404	-	-	0/4/4/4	0/0/0/0
4	GOL	E	405	-	-	0/4/4/4	0/0/0/0
4	GOL	E	406	-	-	0/4/4/4	0/0/0/0
2	MES	F	401	-	-	0/5/13/14	0/1/1/1
4	GOL	F	403	-	-	0/4/4/4	0/0/0/0
4	GOL	F	404	-	-	0/4/4/4	0/0/0/0
4	GOL	F	405	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	FMT	F	406	-	-	0/0/0/0	0/0/0/0
2	MES	G	401	-	-	0/6/14/14	0/1/1/1
4	GOL	G	403	-	-	0/4/4/4	0/0/0/0
4	GOL	G	404	-	-	0/4/4/4	0/0/0/0
4	GOL	G	405	-	-	0/4/4/4	0/0/0/0
4	GOL	G	406	-	-	0/4/4/4	0/0/0/0
4	GOL	G	407	-	-	0/4/4/4	0/0/0/0
5	FMT	G	408	-	-	0/0/0/0	0/0/0/0
2	MES	H	401	-	-	0/6/14/14	0/1/1/1
4	GOL	H	404	-	-	0/4/4/4	0/0/0/0
5	FMT	H	405	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	MES	C7-N4-C5	2.04	116.50	111.27
2	B	401	MES	O1-C2-C3	2.12	116.69	111.84
2	C	401	MES	C7-N4-C5	2.31	117.20	111.27
2	D	401	MES	C7-N4-C3	2.52	117.74	111.27
2	G	401	MES	C7-N4-C3	2.82	118.51	111.27
2	H	401	MES	C7-N4-C3	2.88	118.65	111.27
2	F	401	MES	O2S-S-C8	2.97	110.49	105.81
2	E	401	MES	O2S-S-C8	3.20	109.64	106.91
2	B	401	MES	C7-N4-C3	3.30	119.72	111.27
2	B	401	MES	O1S-S-C8	3.40	109.81	106.91
2	B	401	MES	O2S-S-C8	3.48	109.88	106.91
2	H	401	MES	C5-N4-C3	3.48	116.45	108.90
2	G	401	MES	C5-N4-C3	3.64	116.79	108.90
2	C	401	MES	O2S-S-C8	3.79	110.14	106.91
2	C	401	MES	C5-N4-C3	4.51	118.66	108.90
2	D	401	MES	C5-N4-C3	4.68	119.03	108.90
2	A	401	MES	O2S-S-C8	4.73	110.94	106.91
2	A	401	MES	C5-N4-C3	4.74	119.17	108.90
2	B	401	MES	C5-N4-C3	4.88	119.47	108.90
2	E	401	MES	O1S-S-C8	4.92	111.10	106.91
2	D	401	MES	O1S-S-C8	5.16	111.31	106.91
2	F	401	MES	C5-N4-C3	5.39	120.57	108.90
2	E	401	MES	C5-N4-C3	5.56	120.93	108.90
2	C	401	MES	O1S-S-C8	5.62	111.70	106.91
2	H	401	MES	O1S-S-C8	5.78	111.84	106.91
2	A	401	MES	O1S-S-C8	6.37	112.34	106.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	MES	O1S-S-C8	7.42	113.23	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	407	GOL	1	0
2	B	401	MES	1	0
4	B	403	GOL	1	0
2	C	401	MES	1	0
2	D	401	MES	1	0
4	D	406	GOL	1	0
4	D	407	GOL	4	0
4	E	405	GOL	1	0
4	F	404	GOL	1	0
2	G	401	MES	2	0
4	G	405	GOL	1	0
2	H	401	MES	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	372/384 (96%)	-0.36	7 (1%) 70 75	13, 24, 57, 90	0
1	B	374/384 (97%)	-0.37	9 (2%) 62 68	13, 24, 56, 72	0
1	C	375/384 (97%)	-0.32	16 (4%) 39 48	13, 24, 58, 81	0
1	D	372/384 (96%)	-0.31	16 (4%) 39 48	13, 23, 62, 82	0
1	E	372/384 (96%)	-0.38	16 (4%) 39 48	14, 23, 56, 88	0
1	F	373/384 (97%)	-0.28	14 (3%) 44 53	13, 24, 57, 74	0
1	G	374/384 (97%)	-0.36	9 (2%) 62 68	13, 26, 57, 71	0
1	H	374/384 (97%)	-0.37	10 (2%) 58 65	13, 24, 57, 76	0
All	All	2986/3072 (97%)	-0.34	97 (3%) 51 60	13, 24, 58, 90	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	3	GLY	4.7
1	F	344	ALA	4.6
1	H	341	PRO	4.6
1	G	3	GLY	4.3
1	C	3	GLY	4.3
1	D	40	ASP	4.2
1	B	341	PRO	4.2
1	E	40	ASP	4.2
1	G	344	ALA	4.1
1	D	135	PRO	4.0
1	G	341	PRO	4.0
1	E	345	GLU	3.9
1	E	342	LEU	3.8
1	G	343	SER	3.8
1	C	344	ALA	3.7
1	E	344	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	343	SER	3.7
1	E	343	SER	3.7
1	D	133	ALA	3.7
1	C	341	PRO	3.6
1	B	3	GLY	3.6
1	H	343[A]	SER	3.6
1	A	40	ASP	3.5
1	A	17	PRO	3.5
1	G	133	ALA	3.5
1	D	344	ALA	3.5
1	C	42	PRO	3.4
1	D	42	PRO	3.4
1	C	342	LEU	3.3
1	A	344	ALA	3.3
1	D	348	ALA	3.3
1	F	343	SER	3.3
1	E	42	PRO	3.3
1	F	3	GLY	3.2
1	D	349	ARG	3.2
1	F	41	GLU	3.1
1	C	40	ASP	3.0
1	C	381	THR	2.9
1	E	17	PRO	2.9
1	H	133	ALA	2.9
1	C	38	ALA	2.8
1	D	17	PRO	2.8
1	D	345	GLU	2.8
1	F	341	PRO	2.8
1	C	380	GLY	2.8
1	F	17	PRO	2.7
1	E	41	GLU	2.7
1	H	344	ALA	2.7
1	H	40	ASP	2.7
1	B	349	ARG	2.7
1	D	341	PRO	2.7
1	D	343	SER	2.6
1	H	135	PRO	2.6
1	C	133	ALA	2.6
1	D	342	LEU	2.6
1	G	342	LEU	2.6
1	A	341	PRO	2.5
1	D	97	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	16	ARG	2.5
1	E	16	ARG	2.4
1	D	351	GLY	2.4
1	C	16	ARG	2.4
1	D	126	GLU	2.4
1	G	132	LEU	2.3
1	E	346	ALA	2.3
1	C	345	GLU	2.3
1	A	343	SER	2.3
1	E	348	ALA	2.3
1	B	16	ARG	2.3
1	F	379	GLY	2.3
1	B	17	PRO	2.3
1	H	3	GLY	2.2
1	E	97	ASP	2.2
1	E	349	ARG	2.2
1	E	38	ALA	2.2
1	C	340	ARG	2.2
1	H	43	ARG	2.1
1	B	342	LEU	2.1
1	C	348	ALA	2.1
1	G	346	ALA	2.1
1	B	126	GLU	2.1
1	E	133	ALA	2.1
1	F	133	ALA	2.1
1	F	40	ASP	2.1
1	H	125	PRO	2.1
1	B	344	ALA	2.1
1	H	17	PRO	2.0
1	B	346	ALA	2.0
1	F	346	ALA	2.0
1	A	41	GLU	2.0
1	F	125	PRO	2.0
1	G	17	PRO	2.0
1	C	132	LEU	2.0
1	F	342	LEU	2.0
1	D	130	ALA	2.0
1	F	129	ALA	2.0
1	F	340	ARG	2.0

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

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
1	LLP	E	197	24/25	0.98	0.14	-	13,16,20,21	0
1	LLP	F	197	24/25	0.98	0.15	-	15,18,25,29	0
1	LLP	G	197	24/25	0.99	0.16	-	13,19,22,23	0
1	LLP	H	197	24/25	0.99	0.13	-	13,18,22,23	0
1	LLP	A	197	24/25	0.99	0.13	-	13,18,21,26	0
1	LLP	B	197	24/25	0.98	0.13	-	10,17,22,23	0
1	LLP	C	197	24/25	0.98	0.16	-	13,18,22,25	0
1	LLP	D	197	24/25	0.99	0.12	-	12,19,22,22	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
4	GOL	B	403	6/6	0.84	0.22	8.54	55,59,60,62	0
4	GOL	F	404	6/6	0.83	0.21	8.15	41,50,53,54	0
4	GOL	B	405	6/6	0.87	0.26	7.46	46,51,56,57	0
4	GOL	B	407	6/6	0.76	0.20	6.39	36,58,61,63	0
4	GOL	D	405	6/6	0.82	0.18	4.45	55,58,65,66	0
4	GOL	E	405	6/6	0.91	0.17	4.42	54,58,60,63	0
4	GOL	A	407	6/6	0.85	0.20	4.35	39,51,53,54	0
4	GOL	F	403	6/6	0.89	0.24	4.31	42,56,57,61	0
4	GOL	B	404	6/6	0.94	0.20	3.85	43,45,46,48	0
4	GOL	E	406	6/6	0.88	0.19	3.49	44,50,54,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	GOL	D	404	6/6	0.92	0.18	3.32	41,47,48,50	0
4	GOL	E	403	6/6	0.94	0.22	3.14	42,43,45,51	0
4	GOL	E	404	6/6	0.87	0.17	2.98	48,55,57,57	0
5	FMT	F	406	3/3	0.86	0.15	2.90	29,29,39,44	0
4	GOL	G	405	6/6	0.88	0.14	2.80	39,47,50,50	0
4	GOL	F	405	6/6	0.89	0.19	2.65	44,50,51,52	0
4	GOL	B	406	6/6	0.90	0.33	2.63	37,59,63,64	0
4	GOL	A	404	6/6	0.92	0.17	1.92	43,51,53,57	0
4	GOL	G	403	6/6	0.94	0.15	1.75	39,47,50,56	0
4	GOL	C	404	6/6	0.95	0.16	1.61	39,46,49,51	0
4	GOL	A	406	6/6	0.90	0.15	1.49	45,47,52,52	0
4	GOL	D	406	6/6	0.93	0.14	1.15	46,56,61,64	0
4	GOL	D	407	6/6	0.85	0.20	1.11	49,55,57,57	0
4	GOL	G	406	6/6	0.88	0.17	1.11	65,67,67,69	0
4	GOL	A	405	6/6	0.85	0.23	1.00	33,53,57,58	0
4	GOL	H	404	6/6	0.95	0.12	0.79	40,49,53,54	0
2	MES	F	401	11/12	0.96	0.11	-0.03	28,44,53,54	0
5	FMT	A	408	3/3	0.75	0.13	-0.24	57,57,60,61	0
2	MES	D	401	12/12	0.97	0.09	-0.25	31,38,45,45	0
2	MES	C	401	12/12	0.95	0.10	-0.45	31,37,41,41	0
2	MES	B	401	12/12	0.97	0.08	-0.92	29,32,36,39	0
2	MES	H	401	12/12	0.99	0.08	-1.01	26,29,38,43	0
2	MES	G	401	12/12	0.97	0.09	-1.24	32,38,39,41	0
2	MES	E	401	12/12	0.97	0.09	-1.40	31,39,50,52	0
2	MES	A	401	12/12	0.97	0.09	-1.67	25,28,32,35	0
3	CL	A	402	1/1	0.90	0.08	-	66,66,66,66	0
5	FMT	H	405	3/3	0.91	0.09	-	47,47,47,48	0
3	CL	C	403	1/1	0.94	0.15	-	66,66,66,66	0
3	CL	E	402	1/1	0.95	0.06	-	56,56,56,56	0
3	CL	A	403	1/1	0.92	0.11	-	53,53,53,53	0
4	GOL	G	404	6/6	0.61	0.25	-	48,56,63,64	0
3	CL	G	402	1/1	0.94	0.06	-	66,66,66,66	0
4	GOL	G	407	6/6	0.81	0.25	-	61,64,65,66	0
4	GOL	B	408	6/6	0.83	0.20	-	71,74,75,76	0
3	CL	F	402	1/1	0.95	0.05	-	60,60,60,60	0
3	CL	D	403	1/1	0.95	0.05	-	64,64,64,64	0
3	CL	D	402	1/1	0.90	0.10	-	61,61,61,61	0
3	CL	B	402	1/1	0.96	0.06	-	67,67,67,67	0
5	FMT	G	408	3/3	0.89	0.07	-	43,43,46,46	0
5	FMT	C	405	3/3	0.92	0.07	-	51,51,52,53	0
3	CL	H	402	1/1	0.95	0.07	-	64,64,64,64	0
3	CL	C	402	1/1	0.94	0.07	-	64,64,64,64	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	CL	H	403	1/1	0.93	0.08	-	52,52,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.