



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 09:39 PM GMT

PDB ID : 3HSA  
Title : Crystal structure of pleckstrin homology domain (YP\_926556.1) from SHE-  
WANELLA AMAZONENSIS SB2B at 1.99 Å resolution  
Authors : Joint Center for Structural Genomics (JCSG)  
Deposited on : 2009-06-10  
Resolution : 1.99 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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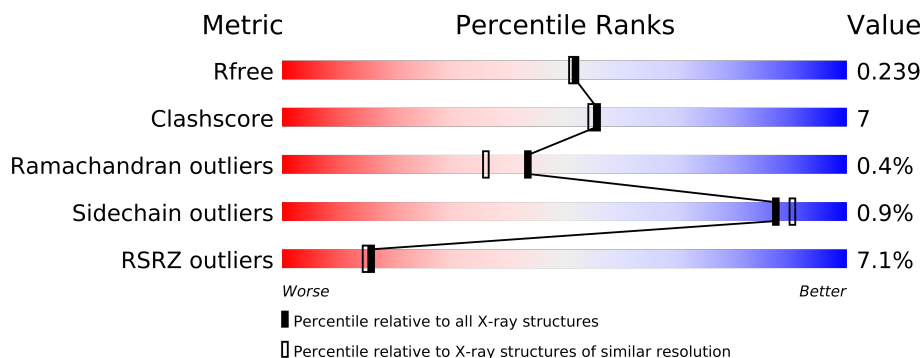
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.99 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))
$R_{free}$	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	126	
1	B	126	
1	C	126	
1	D	126	
1	E	126	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	GOL	A	128	-	X
2	GOL	B	1126	-	X
2	GOL	C	2126	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	GOL	D	3126	-	X
3	PEG	D	3127	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4967 atoms, of which 0 are hydrogen and 0 are deuterium.

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 pleckstrin homology domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	Se	0	4	1
			1015	667	164	179	5			
1	B	124	Total	C	N	O	Se	0	1	1
			974	636	158	177	3			
1	C	112	Total	C	N	O	Se	0	0	1
			896	587	147	160	2			
1	D	113	Total	C	N	O	Se	0	2	0
			914	599	149	164	2			
1	E	111	Total	C	N	O	Se	0	1	0
			876	571	144	159	2			

There are 5 discrepancies between the modelled and reference sequences:

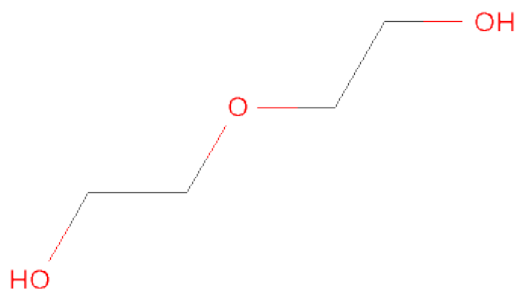
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP A1S3D0
B	0	GLY	-	leader sequence	UNP A1S3D0
C	0	GLY	-	leader sequence	UNP A1S3D0
D	0	GLY	-	leader sequence	UNP A1S3D0
E	0	GLY	-	leader sequence	UNP A1S3D0

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			7	4	3		
3	E	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	63	Total	O	0	0
			63	63		
4	B	49	Total	O	0	0
			49	49		
4	C	36	Total	O	0	0
			36	36		
4	D	61	Total	O	0	0
			61	61		
4	E	33	Total	O	0	0
			33	33		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	33.23Å 129.49Å 138.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.35 – 1.99 47.33 – 1.99	Depositor EDS
% Data completeness (in resolution range)	98.5 (47.35-1.99) 98.5 (47.33-1.99)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.96 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.5.0092	Depositor
R, $R_{free}$	0.189 , 0.233 0.194 , 0.239	Depositor DCC
$R_{free}$ test set	2115 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.7	Xtriage
Anisotropy	0.380	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 45.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 41644 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4967	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PEG, MLY

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	0.74	0/920	0.78	0/1247
1	B	0.69	0/880	0.82	0/1195
1	C	0.73	0/794	0.84	2/1080 (0.2%)
1	D	0.71	0/816	0.78	1/1109 (0.1%)
1	E	0.63	0/793	0.76	0/1076
All	All	0.70	0/4203	0.80	3/5707 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	121	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	C	108	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	C	108	ARG	NE-CZ-NH1	5.22	122.91	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1015	0	1068	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	974	0	987	11	0
1	C	896	0	928	17	0
1	D	914	0	953	15	0
1	E	876	0	885	13	0
2	A	18	0	24	0	0
2	B	6	0	8	0	0
2	C	6	0	8	1	0
2	D	6	0	8	0	0
3	D	7	0	10	0	0
3	E	7	0	10	1	0
4	A	63	0	0	1	0
4	B	49	0	0	2	0
4	C	36	0	0	1	0
4	D	61	0	0	0	0
4	E	33	0	0	0	0
All	All	4967	0	4889	63	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (63) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:40:VAL:HG11	1:A:107:MLY:HD2	1.21	1.19
1:C:39:MSE:HE2	1:C:110:THR:HG23	1.32	1.08
1:C:39:MSE:CE	1:C:110:THR:HG23	1.98	0.93
1:C:39:MSE:HE1	4:C:2167:HOH:O	1.73	0.89
1:E:93[B]:LEU:HD21	1:E:119:ILE:HD13	1.62	0.81
1:A:63:MLY:CH1	1:D:87:MSE:HE2	2.11	0.80
1:E:42:ASP:OD2	1:E:104:MLY:NZ	2.15	0.79
1:A:40:VAL:CG1	1:A:107:MLY:HD2	2.09	0.79
1:C:38:MLY:O	1:C:39:MSE:HE3	1.86	0.74
1:A:63:MLY:HH11	1:D:87:MSE:HE2	1.75	0.69
1:E:93[A]:LEU:CD2	1:E:95:ILE:HD11	2.25	0.66
1:D:72:MLY:HH22	1:D:72:MLY:HG2	1.78	0.65
1:B:8:MSE:N	1:B:8:MSE:HE2	2.12	0.65
1:C:39:MSE:HE2	1:C:110:THR:CG2	2.19	0.64
1:E:79:VAL:HG21	1:E:116:GLN:HG2	1.80	0.62
1:C:98:GLN:HG3	2:C:2126:GOL:O2	2.00	0.61
1:C:79:VAL:HG21	1:C:116:GLN:HG2	1.83	0.59
1:B:7:LEU:O	1:B:8:MSE:HB2	2.04	0.57
1:D:72:MLY:CG	1:D:72:MLY:HH22	2.34	0.57
1:A:87:MSE:HA	1:A:87:MSE:HE2	1.87	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:27:LEU:O	1:D:117:MLY:HH12	2.07	0.54
1:C:75:VAL:HA	1:D:72:MLY:HH21	1.90	0.54
1:E:93[A]:LEU:CD2	1:E:95:ILE:CD1	2.87	0.53
1:E:74:ILE:HG12	1:E:93[A]:LEU:HD21	1.90	0.52
1:A:40:VAL:HG23	1:A:41:ARG:HG2	1.91	0.52
1:E:93[A]:LEU:HD22	1:E:95:ILE:HD11	1.92	0.51
1:C:79:VAL:HG21	1:C:116:GLN:CG	2.40	0.50
1:A:8[B]:MSE:SE	4:A:151:HOH:O	2.80	0.49
4:B:1240:HOH:O	1:C:72:MLY:HE3	2.10	0.49
1:E:79:VAL:HG21	1:E:116:GLN:CG	2.43	0.47
1:E:37:TYR:HB3	1:E:39:MSE:HE3	1.97	0.47
1:E:55:ASP:OD1	1:E:56:MLY:N	2.47	0.47
1:C:16:LEU:HD22	1:C:33:LEU:CD2	2.44	0.47
1:B:8:MSE:H	1:B:8:MSE:HE2	1.79	0.46
1:D:14:VAL:CG1	1:D:14:VAL:O	2.63	0.46
1:C:62:MLY:HH22	1:C:62:MLY:HG2	1.97	0.46
1:C:38:MLY:C	1:C:39:MSE:HE3	2.46	0.45
1:B:56:MLY:CG	1:B:56:MLY:HH22	2.46	0.45
1:E:107:MLY:O	1:E:109:GLY:N	2.51	0.44
1:B:55:ASP:OD2	1:B:104:MLY:NZ	2.51	0.44
1:C:74:ILE:O	1:D:72:MLY:HH11	2.17	0.44
1:B:121:ARG:NH1	4:B:1229:HOH:O	2.51	0.43
1:C:123:ALA:O	1:D:72:MLY:HH12	2.18	0.43
1:A:1:MSE:HE1	1:A:6:ALA:HA	2.00	0.43
1:D:106:LEU:HB3	1:D:112:VAL:HG22	1.99	0.43
1:C:85:PHE:CD1	1:D:62:MLY:NZ	2.87	0.42
1:D:100:GLU:OE1	1:D:100:GLU:HA	2.19	0.42
1:E:39:MSE:CE	1:E:110:THR:HG23	2.49	0.42
1:A:79:VAL:HG21	1:A:116:GLN:HG2	2.02	0.42
1:B:72:MLY:HH13	1:E:123:ALA:O	2.20	0.42
1:D:16:LEU:HD22	1:D:33:LEU:HB3	2.01	0.42
1:D:80:GLU:OE2	1:D:92:MLY:HH13	2.20	0.42
1:A:40:VAL:HG23	1:A:41:ARG:N	2.36	0.41
1:B:30:ASN:ND2	1:B:49:MLY:HH22	2.36	0.41
1:A:95[B]:ILE:HD11	1:A:102:LEU:CD1	2.51	0.41
1:C:39:MSE:SE	1:C:110:THR:HG23	2.69	0.41
1:D:70:PRO:HB2	1:D:72:MLY:HH23	2.02	0.41
1:B:6:ALA:C	1:B:7:LEU:O	2.58	0.41
1:B:56:MLY:HH22	1:B:56:MLY:HG2	2.03	0.41
1:A:1:MSE:HE2	1:A:5:ASP:HB2	2.03	0.41
1:A:79:VAL:HG12	3:E:4126:PEG:H22	2.02	0.41
1:A:54:ILE:HA	1:A:65:SER:O	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:6:ALA:HA	1:B:11:ALA:HB3	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	116/126 (92%)	114 (98%)	2 (2%)	0	100	100
1	B	112/126 (89%)	108 (96%)	3 (3%)	1 (1%)	25	14
1	C	99/126 (79%)	97 (98%)	2 (2%)	0	100	100
1	D	102/126 (81%)	101 (99%)	1 (1%)	0	100	100
1	E	99/126 (79%)	94 (95%)	4 (4%)	1 (1%)	22	12
All	All	528/630 (84%)	514 (97%)	12 (2%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	108	ARG
1	B	40	VAL

### 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	97/89 (109%)	97 (100%)	0	100	100
1	B	91/89 (102%)	90 (99%)	1 (1%)	84	86
1	C	84/89 (94%)	82 (98%)	2 (2%)	61	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	86/89 (97%)	86 (100%)	0	100	100
1	E	82/89 (92%)	81 (99%)	1 (1%)	82	84
All	All	440/445 (99%)	436 (99%)	4 (1%)	87	90

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	4	LEU
1	C	29	ASP
1	C	39	MSE
1	E	57	GLN

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 chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

55 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	MLY	A	104	1	10,10,11	6.23	1 (10%)	9,11,13	0.69	0
1	MLY	A	107	1	10,10,11	5.92	2 (20%)	9,11,13	0.92	1 (11%)
1	MLY	A	117	1	10,10,11	7.13	2 (20%)	9,11,13	1.49	2 (22%)
1	MLY	A	18	1	10,10,11	5.59	2 (20%)	9,11,13	1.50	2 (22%)
1	MLY	A	38	1	10,10,11	5.46	2 (20%)	9,11,13	3.07	1 (11%)
1	MLY	A	49	1	10,10,11	5.84	2 (20%)	9,11,13	1.23	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	MLY	A	56	1	10,10,11	5.08	2 (20%)	9,11,13	1.81	1 (11%)
1	MLY	A	62	1	10,10,11	6.00	2 (20%)	9,11,13	0.59	0
1	MLY	A	63	1	10,10,11	5.84	2 (20%)	9,11,13	2.02	1 (11%)
1	MLY	A	72	1	10,10,11	5.34	2 (20%)	9,11,13	0.49	0
1	MLY	A	92	1	10,10,11	5.72	1 (10%)	9,11,13	2.17	1 (11%)
1	MLY	B	104	1	7,8,11	8.70	1 (14%)	5,8,13	2.22	1 (20%)
1	MLY	B	107	1	10,10,11	6.22	2 (20%)	9,11,13	1.48	1 (11%)
1	MLY	B	117	1	7,8,11	6.13	2 (28%)	5,8,13	1.46	1 (20%)
1	MLY	B	18	1	10,10,11	4.34	2 (20%)	9,11,13	1.50	1 (11%)
1	MLY	B	38	1	10,10,11	5.10	3 (30%)	9,11,13	1.37	1 (11%)
1	MLY	B	49	1	10,10,11	5.34	3 (30%)	9,11,13	1.04	1 (11%)
1	MLY	B	56	1	10,10,11	5.28	1 (10%)	9,11,13	3.17	1 (11%)
1	MLY	B	62	1	3,4,11	10.46	2 (66%)	1,4,13	0.27	0
1	MLY	B	63	1	10,10,11	6.71	2 (20%)	9,11,13	1.20	1 (11%)
1	MLY	B	72	1	10,10,11	5.29	3 (30%)	9,11,13	0.84	0
1	MLY	B	92	1	10,10,11	5.42	1 (10%)	9,11,13	1.50	1 (11%)
1	MLY	C	104	1	10,10,11	5.18	1 (10%)	9,11,13	1.04	1 (11%)
1	MLY	C	107	1	10,10,11	6.13	3 (30%)	9,11,13	1.42	2 (22%)
1	MLY	C	117	1	10,10,11	5.24	2 (20%)	9,11,13	0.89	1 (11%)
1	MLY	C	18	1	3,4,11	10.57	2 (66%)	1,4,13	2.20	1 (100%)
1	MLY	C	38	1	10,10,11	6.16	2 (20%)	9,11,13	1.50	1 (11%)
1	MLY	C	49	1	10,10,11	5.91	2 (20%)	9,11,13	0.99	1 (11%)
1	MLY	C	56	1	10,10,11	5.55	2 (20%)	9,11,13	1.82	1 (11%)
1	MLY	C	62	1	10,10,11	5.18	2 (20%)	9,11,13	0.95	0
1	MLY	C	63	1	10,10,11	6.54	2 (20%)	9,11,13	1.77	1 (11%)
1	MLY	C	72	1	10,10,11	5.21	3 (30%)	9,11,13	0.58	0
1	MLY	C	92	1	10,10,11	5.90	2 (20%)	9,11,13	1.91	1 (11%)
1	MLY	D	104	1	10,10,11	5.45	3 (30%)	9,11,13	0.76	0
1	MLY	D	107	1	10,10,11	5.61	2 (20%)	9,11,13	0.47	0
1	MLY	D	117	1	10,10,11	4.60	2 (20%)	9,11,13	1.88	1 (11%)
1	MLY	D	18	1	10,10,11	5.55	2 (20%)	9,11,13	1.14	1 (11%)
1	MLY	D	38	1	10,10,11	5.80	2 (20%)	9,11,13	1.76	1 (11%)
1	MLY	D	49	1	10,10,11	6.28	2 (20%)	9,11,13	1.53	1 (11%)
1	MLY	D	56	1	10,10,11	6.08	1 (10%)	9,11,13	3.05	1 (11%)
1	MLY	D	62	1	7,8,11	6.27	1 (14%)	5,8,13	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	MLY	D	63	1	7,8,11	7.54	2 (28%)	5,8,13	2.02	1 (20%)
1	MLY	D	72	1	10,10,11	4.77	3 (30%)	9,11,13	0.86	0
1	MLY	D	92	1	10,10,11	5.09	3 (30%)	9,11,13	2.40	1 (11%)
1	MLY	E	104	1	10,10,11	5.54	1 (10%)	9,11,13	1.17	1 (11%)
1	MLY	E	107	1	7,8,11	6.82	1 (14%)	5,8,13	0.67	0
1	MLY	E	117	1	10,10,11	6.22	2 (20%)	9,11,13	0.98	0
1	MLY	E	18	1	3,4,11	10.48	2 (66%)	1,4,13	5.35	1 (100%)
1	MLY	E	38	1	10,10,11	5.37	2 (20%)	9,11,13	1.65	1 (11%)
1	MLY	E	49	1	7,8,11	7.27	2 (28%)	5,8,13	1.08	1 (20%)
1	MLY	E	56	1	10,10,11	5.79	2 (20%)	9,11,13	2.27	1 (11%)
1	MLY	E	62	1	3,4,11	10.87	2 (66%)	1,4,13	1.07	0
1	MLY	E	63	1	3,4,11	10.31	2 (66%)	1,4,13	1.99	0
1	MLY	E	72	1	10,10,11	5.37	2 (20%)	9,11,13	0.54	0
1	MLY	E	92	1	10,10,11	6.10	2 (20%)	9,11,13	1.31	1 (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
1	MLY	A	104	1	-	0/7/9/11	0/0/0/0
1	MLY	A	107	1	-	0/7/9/11	0/0/0/0
1	MLY	A	117	1	-	0/7/9/11	0/0/0/0
1	MLY	A	18	1	-	0/7/9/11	0/0/0/0
1	MLY	A	38	1	-	0/7/9/11	0/0/0/0
1	MLY	A	49	1	-	0/7/9/11	0/0/0/0
1	MLY	A	56	1	-	0/7/9/11	0/0/0/0
1	MLY	A	62	1	-	0/7/9/11	0/0/0/0
1	MLY	A	63	1	-	0/7/9/11	0/0/0/0
1	MLY	A	72	1	-	0/7/9/11	0/0/0/0
1	MLY	A	92	1	-	0/7/9/11	0/0/0/0
1	MLY	B	104	1	-	0/5/7/11	0/0/0/0
1	MLY	B	107	1	-	0/7/9/11	0/0/0/0
1	MLY	B	117	1	-	0/5/7/11	0/0/0/0
1	MLY	B	18	1	-	0/7/9/11	0/0/0/0
1	MLY	B	38	1	-	0/7/9/11	0/0/0/0
1	MLY	B	49	1	-	0/7/9/11	0/0/0/0
1	MLY	B	56	1	-	0/7/9/11	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	B	62	1	-	0/0/2/11	0/0/0/0
1	MLY	B	63	1	-	0/7/9/11	0/0/0/0
1	MLY	B	72	1	-	0/7/9/11	0/0/0/0
1	MLY	B	92	1	-	0/7/9/11	0/0/0/0
1	MLY	C	104	1	-	0/7/9/11	0/0/0/0
1	MLY	C	107	1	-	0/7/9/11	0/0/0/0
1	MLY	C	117	1	-	0/7/9/11	0/0/0/0
1	MLY	C	18	1	-	0/0/2/11	0/0/0/0
1	MLY	C	38	1	-	0/7/9/11	0/0/0/0
1	MLY	C	49	1	-	0/7/9/11	0/0/0/0
1	MLY	C	56	1	-	0/7/9/11	0/0/0/0
1	MLY	C	62	1	-	0/7/9/11	0/0/0/0
1	MLY	C	63	1	-	0/7/9/11	0/0/0/0
1	MLY	C	72	1	-	0/7/9/11	0/0/0/0
1	MLY	C	92	1	-	0/7/9/11	0/0/0/0
1	MLY	D	104	1	-	0/7/9/11	0/0/0/0
1	MLY	D	107	1	-	0/7/9/11	0/0/0/0
1	MLY	D	117	1	-	0/7/9/11	0/0/0/0
1	MLY	D	18	1	-	0/7/9/11	0/0/0/0
1	MLY	D	38	1	-	0/7/9/11	0/0/0/0
1	MLY	D	49	1	-	0/7/9/11	0/0/0/0
1	MLY	D	56	1	-	0/7/9/11	0/0/0/0
1	MLY	D	62	1	-	0/5/7/11	0/0/0/0
1	MLY	D	63	1	-	0/5/7/11	0/0/0/0
1	MLY	D	72	1	-	0/7/9/11	0/0/0/0
1	MLY	D	92	1	-	0/7/9/11	0/0/0/0
1	MLY	E	104	1	-	0/7/9/11	0/0/0/0
1	MLY	E	107	1	-	0/5/7/11	0/0/0/0
1	MLY	E	117	1	-	0/7/9/11	0/0/0/0
1	MLY	E	18	1	-	0/0/2/11	0/0/0/0
1	MLY	E	38	1	-	0/7/9/11	0/0/0/0
1	MLY	E	49	1	-	0/5/7/11	0/0/0/0
1	MLY	E	56	1	-	0/7/9/11	0/0/0/0
1	MLY	E	62	1	-	0/0/2/11	0/0/0/0
1	MLY	E	63	1	-	0/0/2/11	0/0/0/0
1	MLY	E	72	1	-	0/7/9/11	0/0/0/0
1	MLY	E	92	1	-	0/7/9/11	0/0/0/0

All (108) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	104	MLY	O-C	22.93	1.27	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	MLY	O-C	22.27	1.26	1.11
1	B	63	MLY	O-C	20.80	1.25	1.11
1	C	63	MLY	O-C	20.43	1.25	1.11
1	D	49	MLY	O-C	19.66	1.25	1.11
1	D	63	MLY	O-C	19.63	1.24	1.11
1	A	104	MLY	O-C	19.50	1.24	1.11
1	B	107	MLY	O-C	19.45	1.24	1.11
1	E	117	MLY	O-C	19.44	1.24	1.11
1	C	38	MLY	O-C	19.15	1.24	1.11
1	E	92	MLY	O-C	19.07	1.24	1.11
1	D	56	MLY	O-C	19.04	1.24	1.11
1	E	49	MLY	O-C	19.02	1.24	1.11
1	C	107	MLY	O-C	18.74	1.24	1.11
1	A	62	MLY	O-C	18.71	1.24	1.11
1	E	62	MLY	O-C	18.58	1.24	1.11
1	A	107	MLY	O-C	18.41	1.24	1.11
1	C	92	MLY	O-C	18.40	1.24	1.11
1	A	49	MLY	O-C	18.16	1.23	1.11
1	A	63	MLY	O-C	18.15	1.23	1.11
1	C	49	MLY	O-C	18.13	1.23	1.11
1	C	18	MLY	O-C	18.10	1.23	1.11
1	E	107	MLY	O-C	17.98	1.23	1.11
1	E	56	MLY	O-C	17.96	1.23	1.11
1	A	92	MLY	O-C	17.94	1.23	1.11
1	D	38	MLY	O-C	17.94	1.23	1.11
1	E	18	MLY	O-C	17.84	1.23	1.11
1	B	62	MLY	O-C	17.84	1.23	1.11
1	E	63	MLY	O-C	17.63	1.23	1.11
1	D	107	MLY	O-C	17.38	1.23	1.11
1	E	104	MLY	O-C	17.35	1.23	1.11
1	D	18	MLY	O-C	17.34	1.23	1.11
1	A	18	MLY	O-C	17.26	1.23	1.11
1	C	56	MLY	O-C	17.26	1.23	1.11
1	A	38	MLY	O-C	16.99	1.23	1.11
1	B	92	MLY	O-C	16.97	1.23	1.11
1	D	104	MLY	O-C	16.81	1.23	1.11
1	E	38	MLY	O-C	16.69	1.22	1.11
1	E	72	MLY	O-C	16.50	1.22	1.11
1	D	62	MLY	O-C	16.49	1.22	1.11
1	B	56	MLY	O-C	16.42	1.22	1.11
1	B	49	MLY	O-C	16.38	1.22	1.11
1	A	72	MLY	O-C	16.24	1.22	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	117	MLY	O-C	16.23	1.22	1.11
1	C	104	MLY	O-C	16.21	1.22	1.11
1	B	72	MLY	O-C	16.09	1.22	1.11
1	C	62	MLY	O-C	16.04	1.22	1.11
1	C	72	MLY	O-C	15.94	1.22	1.11
1	B	117	MLY	O-C	15.87	1.22	1.11
1	A	56	MLY	O-C	15.76	1.22	1.11
1	D	92	MLY	O-C	15.73	1.22	1.11
1	B	38	MLY	O-C	15.70	1.22	1.11
1	D	72	MLY	O-C	14.44	1.21	1.11
1	D	117	MLY	O-C	14.11	1.21	1.11
1	B	18	MLY	O-C	13.34	1.20	1.11
1	A	72	MLY	CA-C	3.94	1.55	1.48
1	B	63	MLY	CA-C	3.84	1.55	1.48
1	C	49	MLY	CA-C	3.81	1.55	1.48
1	C	107	MLY	CA-C	3.76	1.55	1.48
1	D	63	MLY	CA-C	3.34	1.54	1.48
1	D	38	MLY	CA-C	3.29	1.54	1.48
1	E	18	MLY	CA-C	3.28	1.54	1.48
1	B	62	MLY	CA-C	3.21	1.54	1.48
1	C	117	MLY	CA-C	3.19	1.54	1.48
1	B	117	MLY	CA-C	3.19	1.54	1.48
1	A	117	MLY	CA-C	3.16	1.54	1.48
1	B	49	MLY	CA-C	3.13	1.54	1.48
1	E	72	MLY	CA-C	3.10	1.54	1.48
1	C	72	MLY	CA-C	3.07	1.54	1.48
1	A	18	MLY	CA-C	2.96	1.53	1.48
1	E	56	MLY	CA-C	2.96	1.53	1.48
1	E	62	MLY	CA-C	2.89	1.53	1.48
1	A	49	MLY	CA-C	2.87	1.53	1.48
1	D	72	MLY	CA-C	2.87	1.53	1.48
1	D	104	MLY	CA-C	2.86	1.53	1.48
1	D	107	MLY	CA-C	2.83	1.53	1.48
1	C	38	MLY	CA-C	2.80	1.53	1.48
1	E	49	MLY	CA-C	2.80	1.53	1.48
1	A	63	MLY	CA-C	2.78	1.53	1.48
1	E	63	MLY	CA-C	2.76	1.53	1.48
1	B	38	MLY	CA-C	2.70	1.53	1.48
1	B	72	MLY	CB-CA	2.67	1.55	1.53
1	E	38	MLY	CA-C	2.65	1.53	1.48
1	B	72	MLY	CE-NZ	2.57	1.56	1.46
1	C	63	MLY	CA-C	2.50	1.53	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	18	MLY	CA-C	2.46	1.52	1.48
1	C	18	MLY	CA-C	2.45	1.52	1.48
1	E	117	MLY	CA-C	2.44	1.52	1.48
1	D	49	MLY	CA-C	2.40	1.52	1.48
1	B	49	MLY	CB-CA	2.37	1.55	1.53
1	C	56	MLY	CB-CA	2.31	1.55	1.53
1	A	62	MLY	CA-C	2.29	1.52	1.48
1	C	92	MLY	CA-C	2.27	1.52	1.48
1	E	92	MLY	CA-C	2.26	1.52	1.48
1	A	38	MLY	CA-C	2.23	1.52	1.48
1	D	72	MLY	CB-CA	2.21	1.55	1.53
1	D	92	MLY	CA-C	2.20	1.52	1.48
1	C	72	MLY	CE-NZ	2.19	1.55	1.46
1	D	18	MLY	CA-C	2.18	1.52	1.48
1	B	38	MLY	CB-CA	-2.18	1.51	1.53
1	C	62	MLY	CA-C	2.16	1.52	1.48
1	D	104	MLY	CB-CA	2.16	1.55	1.53
1	D	92	MLY	CE-NZ	2.13	1.54	1.46
1	A	107	MLY	CE-NZ	2.12	1.54	1.46
1	C	107	MLY	CE-NZ	2.11	1.54	1.46
1	A	56	MLY	CE-NZ	2.10	1.54	1.46
1	D	117	MLY	CA-C	2.06	1.52	1.48
1	B	107	MLY	CA-C	2.01	1.52	1.48

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	56	MLY	C-CA-N	-9.36	104.48	113.83
1	D	56	MLY	C-CA-N	-9.08	104.76	113.83
1	A	38	MLY	C-CA-N	-8.94	104.90	113.83
1	D	92	MLY	C-CA-N	-6.98	106.86	113.83
1	E	56	MLY	C-CA-N	-6.73	107.11	113.83
1	A	92	MLY	C-CA-N	-6.11	107.72	113.83
1	A	63	MLY	C-CA-N	-5.75	108.09	113.83
1	C	92	MLY	C-CA-N	-5.61	108.23	113.83
1	D	117	MLY	C-CA-N	-5.42	108.42	113.83
1	E	18	MLY	C-CA-N	-5.35	108.48	113.83
1	A	56	MLY	C-CA-N	-5.33	108.50	113.83
1	C	56	MLY	C-CA-N	-5.28	108.55	113.83
1	C	63	MLY	C-CA-N	-5.02	108.82	113.83
1	B	104	MLY	C-CA-N	-4.82	109.02	113.83
1	E	38	MLY	C-CA-N	-4.71	109.12	113.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	63	MLY	C-CA-N	-4.47	109.37	113.83
1	D	38	MLY	C-CA-N	-4.41	109.42	113.83
1	B	92	MLY	C-CA-N	-4.26	109.57	113.83
1	C	38	MLY	C-CA-N	-4.25	109.59	113.83
1	B	18	MLY	C-CA-N	-4.14	109.69	113.83
1	D	49	MLY	C-CA-N	4.07	117.90	113.83
1	B	107	MLY	C-CA-N	-3.88	109.95	113.83
1	E	92	MLY	C-CA-N	-3.73	110.11	113.83
1	A	117	MLY	CD-CE-NZ	-3.29	104.98	113.59
1	E	104	MLY	C-CA-N	-3.25	110.58	113.83
1	B	117	MLY	C-CA-N	-3.24	110.59	113.83
1	A	18	MLY	C-CA-N	-3.17	110.66	113.83
1	A	49	MLY	C-CA-N	3.14	116.97	113.83
1	B	38	MLY	C-CA-N	-3.08	110.75	113.83
1	B	63	MLY	C-CA-N	-3.07	110.76	113.83
1	C	107	MLY	C-CA-N	-3.06	110.77	113.83
1	D	18	MLY	C-CA-N	-2.82	111.01	113.83
1	C	107	MLY	CD-CE-NZ	2.66	120.56	113.59
1	C	49	MLY	CD-CE-NZ	-2.37	107.39	113.59
1	C	117	MLY	C-CA-N	-2.35	111.48	113.83
1	B	49	MLY	C-CA-N	2.35	116.18	113.83
1	A	117	MLY	CH2-NZ-CH1	2.32	116.07	109.75
1	A	18	MLY	CH2-NZ-CH1	2.23	115.82	109.75
1	C	18	MLY	C-CA-N	-2.20	111.63	113.83
1	E	49	MLY	C-CA-N	2.15	115.98	113.83
1	A	107	MLY	C-CA-N	2.12	115.95	113.83
1	C	104	MLY	C-CA-N	-2.11	111.72	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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	$\# Z  > 2$
2	GOL	A	126	-	5,5,5	0.34	0	5,5,5	0.69	0
2	GOL	A	127	-	5,5,5	0.39	0	5,5,5	0.52	0
2	GOL	A	128	-	5,5,5	0.49	0	5,5,5	0.55	0
2	GOL	B	1126	-	5,5,5	0.44	0	5,5,5	0.56	0
2	GOL	C	2126	-	5,5,5	0.56	0	5,5,5	0.45	0
2	GOL	D	3126	-	5,5,5	0.30	0	5,5,5	0.38	0
3	PEG	D	3127	-	6,6,6	0.45	0	5,5,5	0.37	0
3	PEG	E	4126	-	6,6,6	0.33	0	5,5,5	0.39	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	GOL	A	126	-	-	0/4/4/4	0/0/0/0
2	GOL	A	127	-	-	0/4/4/4	0/0/0/0
2	GOL	A	128	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1126	-	-	0/4/4/4	0/0/0/0
2	GOL	C	2126	-	-	0/4/4/4	0/0/0/0
2	GOL	D	3126	-	-	0/4/4/4	0/0/0/0
3	PEG	D	3127	-	-	0/4/4/4	0/0/0/0
3	PEG	E	4126	-	-	0/4/4/4	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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	125/126 (99%)	0.08	6 (4%)	29	28	9, 17, 33, 48	0
1	B	124/126 (98%)	0.31	6 (4%)	29	28	10, 19, 37, 46	0
1	C	112/126 (88%)	0.28	8 (7%)	16	15	11, 20, 33, 38	0
1	D	113/126 (89%)	0.04	4 (3%)	42	41	12, 21, 36, 43	0
1	E	111/126 (88%)	0.75	18 (16%)	2	3	11, 23, 47, 55	0
All	All	585/630 (92%)	0.29	42 (7%)	16	14	9, 20, 39, 55	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	59	VAL	6.9
1	B	10	ASN	5.0
1	E	58	GLY	4.8
1	E	109	GLY	4.7
1	B	85	PHE	4.5
1	A	84	THR	4.5
1	E	60	THR	4.2
1	E	56	MLY	4.2
1	E	62	MLY	4.1
1	A	85	PHE	3.9
1	B	9	GLY	3.8
1	E	61	GLY	3.8
1	B	11	ALA	3.7
1	A	82	ALA	3.6
1	E	21	ALA	3.5
1	E	29	ASP	3.5
1	C	14	VAL	3.3
1	E	16	LEU	3.1
1	E	15	ASP	3.1
1	A	83	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	107	MLY	2.8
1	D	14	VAL	2.7
1	E	110	THR	2.6
1	C	110	THR	2.5
1	C	85	PHE	2.4
1	C	97	GLY	2.4
1	A	86	ASP	2.4
1	C	59	VAL	2.4
1	B	99[A]	HIS	2.3
1	E	57	GLN	2.3
1	B	86	ASP	2.3
1	D	125	GLY	2.3
1	D	30	ASN	2.3
1	C	29	ASP	2.2
1	A	30	ASN	2.2
1	C	107	MLY	2.2
1	C	84	THR	2.1
1	D	29	ASP	2.1
1	E	64	VAL	2.1
1	E	125	GLY	2.0
1	E	87	MSE	2.0
1	E	88	ASP	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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	MLY	E	72	11/12	0.20	9.49	14,16,33,34	0
1	MLY	C	72	11/12	0.21	8.38	13,18,36,37	0
1	MLY	A	62	11/12	0.18	4.58	19,21,40,41	0
1	MLY	B	18	11/12	0.18	3.46	12,16,32,35	0
1	MLY	B	62	5/12	0.23	2.71	34,34,34,36	0
1	MLY	B	38	11/12	0.14	2.45	17,22,33,33	0
1	MLY	D	72	11/12	0.15	2.45	14,19,34,36	0
1	MLY	C	107	11/12	0.27	2.43	25,30,43,44	0
1	MLY	E	62	5/12	0.39	1.97	45,46,46,48	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	MLY	B	92	11/12	0.14	1.86	12,13,26,29	0
1	MLY	B	104	9/12	0.16	1.82	11,16,30,33	0
1	MLY	E	56	11/12	0.39	1.79	44,49,51,51	0
1	MLY	B	56	11/12	0.17	1.71	21,26,41,42	0
1	MLY	C	104	11/12	0.17	1.67	16,20,28,29	0
1	MLY	B	63	11/12	0.20	1.54	31,33,41,41	0
1	MLY	E	107	9/12	0.31	1.41	33,35,41,42	0
1	MLY	C	49	11/12	0.15	1.28	17,21,36,37	0
1	MLY	C	38	11/12	0.21	1.25	22,28,40,41	0
1	MLY	D	49	11/12	0.14	1.19	20,23,38,38	0
1	MLY	A	107	11/12	0.20	1.08	22,26,46,47	0
1	MLY	C	62	11/12	0.20	0.94	24,28,37,37	0
1	MLY	A	38	11/12	0.12	0.94	14,18,34,34	0
1	MLY	D	18	11/12	0.13	0.93	17,19,36,36	0
1	MLY	A	18	11/12	0.11	0.67	10,15,24,24	0
1	MLY	B	72	11/12	0.14	0.66	14,20,34,34	0
1	MLY	B	107	11/12	0.12	0.64	18,23,33,33	0
1	MLY	E	117	11/12	0.14	0.48	16,21,27,27	0
1	MLY	A	92	11/12	0.12	0.45	13,14,28,29	0
1	MLY	C	117	11/12	0.14	0.44	14,17,30,33	0
1	MLY	E	38	11/12	0.20	0.43	26,32,37,37	0
1	MLY	A	56	11/12	0.16	0.37	17,21,37,38	0
1	MLY	A	63	11/12	0.13	0.35	16,18,29,30	0
1	MLY	C	56	11/12	0.13	0.21	20,25,31,33	0
1	MLY	D	92	11/12	0.13	0.15	15,16,32,33	0
1	MLY	D	62	9/12	0.13	0.06	30,32,39,40	0
1	MLY	D	104	11/12	0.11	0.02	14,16,17,17	0
1	MLY	B	117	9/12	0.12	0.01	14,16,24,28	0
1	MLY	D	63	9/12	0.14	-0.02	25,29,41,41	0
1	MLY	E	92	11/12	0.11	-0.04	15,16,24,26	0
1	MLY	A	49	11/12	0.12	-0.09	10,14,29,32	0
1	MLY	A	104	11/12	0.10	-0.20	13,15,21,22	0
1	MLY	C	63	11/12	0.12	-0.20	19,22,25,27	0
1	MLY	D	107	11/12	0.11	-0.29	24,27,37,38	0
1	MLY	D	38	11/12	0.11	-0.31	16,19,22,25	0
1	MLY	A	117	11/12	0.11	-0.33	10,15,32,34	0
1	MLY	D	56	11/12	0.10	-0.35	18,21,29,31	0
1	MLY	B	49	11/12	0.11	-0.37	15,17,29,30	0
1	MLY	E	49	9/12	0.09	-0.49	16,18,27,28	0
1	MLY	D	117	11/12	0.09	-0.56	11,16,29,29	0
1	MLY	E	63	5/12	0.19	-0.63	37,37,39,42	0
1	MLY	C	92	11/12	0.08	-0.68	14,15,24,24	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	MLY	E	104	11/12	0.10	-0.76	20,21,23,26	0
1	MLY	A	72	11/12	0.09	-0.80	11,17,31,32	0
1	MLY	E	18	5/12	0.14	-1.13	35,35,36,36	0
1	MLY	C	18	5/12	0.07	-1.46	14,16,18,19	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, 95<sup>th</sup> 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GOL	D	3126	6/6	0.28	7.89	61,62,63,64	0
2	GOL	B	1126	6/6	0.23	5.25	57,59,60,60	0
2	GOL	A	128	6/6	0.17	4.84	48,52,52,54	0
3	PEG	D	3127	7/7	0.16	4.79	39,40,41,42	0
2	GOL	C	2126	6/6	0.24	2.61	54,55,56,57	0
3	PEG	E	4126	7/7	0.17	1.02	45,46,48,49	0
2	GOL	A	127	6/6	0.10	-0.09	22,25,31,36	0
2	GOL	A	126	6/6	0.09	-0.38	20,22,23,23	0

### 6.5 Other polymers

There are no such residues in this entry.