Full wwPDB X-ray Structure Validation Report (i)

Sep 18, 2023 – 04:03 PM EDT

PDB ID	:	5CZO
Title	:	Structure of S. cerevisiae Hrr25:Mam1 complex, form 2
Authors	:	Ye, Q.; Corbett, K.D.
Deposited on	:	2015-07-31
Resolution	:	2.89 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

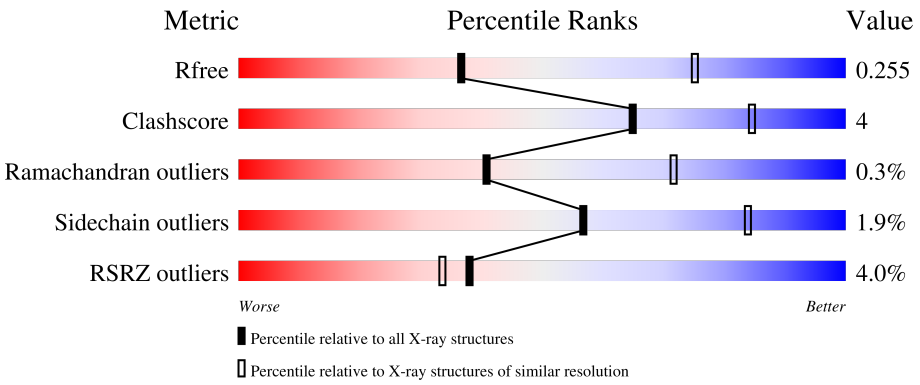
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.89 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	395	3% 81% 14% 5%
1	B	395	6% 83% 10% • 5%
2	C	105	2% 79% 16% • •
2	D	105	2% 92% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7958 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 Casein kinase I homolog HRR25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	374	3098	1984	548	551	15	0	0	0
1	B	374	3096	1981	546	554	15	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	expression tag	UNP P29295
A	38	ARG	LYS	engineered mutation	UNP P29295
B	0	ALA	-	expression tag	UNP P29295
B	38	ARG	LYS	engineered mutation	UNP P29295

- Molecule 2 is a protein called Monopolin complex subunit MAM1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	101	868	554	145	162	7	0	0	0
2	D	101	855	545	142	161	7	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
3	C	1	1	1	0	0
3	D	1	1	1	0	0

- Molecule 4 is water.

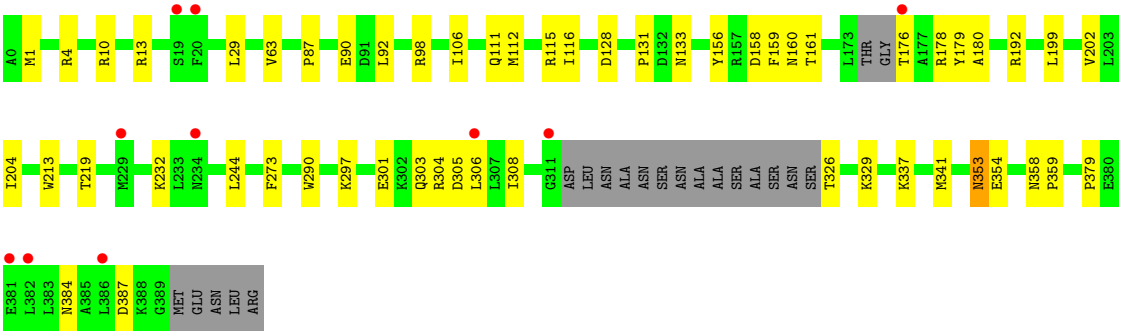
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	25	25	25	0	0
			Total	O		
4	C	4	4	4	0	0
			Total	O		
4	B	5	5	5	0	0
			Total	O		
4	D	5	5	5	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

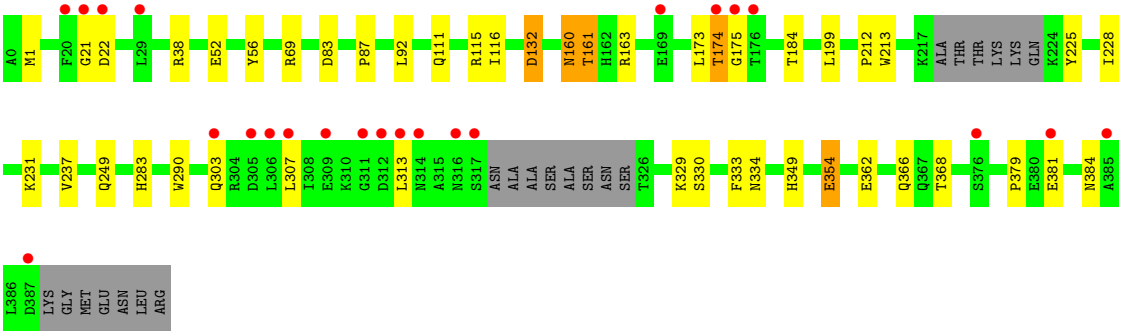
• Molecule 1: Casein kinase I homolog HRR25

Chain A:



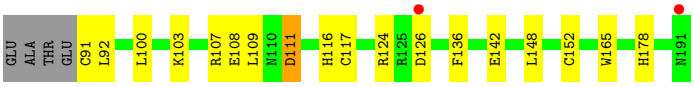
• Molecule 1: Casein kinase I homolog HRR25

Chain B:



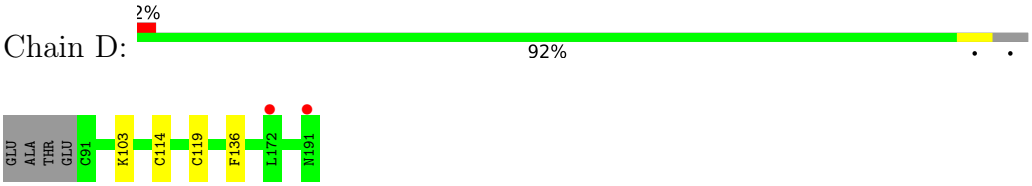
• Molecule 2: Monopolin complex subunit MAM1

Chain C:



• Molecule 2: Monopolin complex subunit MAM1

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	52.70Å 83.84Å 132.51Å 90.00° 92.57° 90.00°	Depositor
Resolution (Å)	49.69 – 2.89 49.69 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.69-2.89) 99.5 (49.69-2.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) >$ [1]	3.23 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.218 , 0.253 0.218 , 0.255	Depositor DCC
R_{free} test set	1320 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å^2)	72.2	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 45.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7958	wwPDB-VP
Average B, all atoms (Å^2)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: MLY, ZN

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	0.21	0/3068	0.36	0/4128
1	B	0.22	0/3066	0.35	0/4129
2	C	0.21	0/865	0.33	0/1168
2	D	0.21	0/851	0.32	0/1150
All	All	0.21	0/7850	0.35	0/10575

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	3098	0	3098	29	0
1	B	3096	0	3085	29	0
2	C	868	0	835	11	0
2	D	855	0	817	3	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	25	0	0	0	0
4	B	5	0	0	0	0
4	C	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	5	0	0	0	0
All	All	7958	0	7835	65	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 4.

All (65) 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)
2:C:108:GLU:O	2:C:124:ARG:NH1	2.20	0.74
1:A:63:VAL:HG22	2:C:178:HIS:HB3	1.74	0.68
1:B:303:GLN:HE21	1:B:384:ASN:HB3	1.59	0.66
1:A:180:ALA:O	1:A:232:LYS:NZ	2.32	0.62
1:B:174:THR:OG1	1:B:175:GLY:N	2.38	0.57
1:B:330:SER:O	1:B:334:ASN:ND2	2.29	0.56
1:B:307:LEU:HD22	1:B:313:LEU:HD23	1.87	0.56
1:B:160:ASN:OD1	1:B:160:ASN:N	2.30	0.55
2:C:91:CYS:SG	2:C:92:LEU:N	2.81	0.54
1:A:116:ILE:HD12	1:A:199:LEU:HD22	1.89	0.54
1:B:213:TRP:HB2	1:B:228:ILE:HG23	1.91	0.53
1:B:213:TRP:HB3	1:B:231:MLY:HG3	1.93	0.51
1:A:4:ARG:HD3	1:A:29:LEU:HD11	1.91	0.51
1:A:326:THR:HA	1:A:329:LYS:HE3	1.95	0.49
1:A:353:ASN:N	1:A:353:ASN:OD1	2.44	0.49
1:B:237:VAL:HG12	1:B:249:GLN:HG3	1.94	0.49
1:A:384:ASN:HA	1:A:387:ASP:HB3	1.94	0.48
1:B:329:LYS:O	1:B:333:PHE:HB3	2.14	0.48
1:A:92:LEU:HD13	1:A:290:TRP:CE2	2.49	0.48
1:B:21:GLY:HA3	1:B:38:ARG:NH1	2.29	0.47
1:B:174:THR:HG22	1:B:184:THR:HG21	1.96	0.47
1:B:1:MET:HE2	2:D:103:LYS:HD3	1.95	0.47
1:B:69:ARG:NH2	1:B:83:ASP:HB3	2.30	0.47
1:B:116:ILE:HD12	1:B:199:LEU:HD22	1.96	0.47
1:A:158:ASP:OD1	1:A:159:PHE:N	2.41	0.47
1:B:69:ARG:HH21	1:B:83:ASP:HB3	1.80	0.46
1:A:329:LYS:HZ3	2:C:126:ASP:HA	1.80	0.46
1:B:111:GLN:O	1:B:115:ARG:HG2	2.16	0.46
2:D:114:CYS:SG	2:D:119:CYS:HB2	2.55	0.46
2:C:116:HIS:CE1	2:C:152:CYS:SG	3.09	0.46
1:B:87:PRO:HG3	2:D:136:PHE:HB2	1.98	0.45
1:A:178:ARG:HD2	1:A:179:TYR:CZ	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:PRO:HG3	1:A:202:VAL:HG13	1.99	0.45
1:A:128:ASP:O	1:A:133:ASN:ND2	2.47	0.45
1:B:354:GLU:H	1:B:354:GLU:HG3	1.49	0.45
1:A:304:ARG:O	1:A:308:ILE:HG12	2.17	0.44
1:B:173:LEU:HD22	1:B:225:TYR:HE2	1.83	0.44
1:A:111:GLN:O	1:A:115:ARG:HG2	2.17	0.44
1:B:303:GLN:NE2	1:B:381:GLU:O	2.48	0.44
1:A:1:MET:HE3	2:C:100:LEU:HD12	1.99	0.44
1:B:349:HIS:NE2	1:B:368:THR:OG1	2.43	0.43
1:A:160:ASN:OD1	1:A:161:THR:N	2.51	0.43
1:B:212:PRO:HG2	1:B:213:TRP:CE3	2.53	0.43
1:A:87:PRO:HG3	2:C:136:PHE:HB2	2.00	0.43
1:B:52:GLU:HG2	1:B:56:TYR:CE2	2.54	0.43
1:A:10:ARG:HD2	2:C:165:TRP:CE2	2.53	0.43
1:B:132:ASP:N	1:B:132:ASP:OD1	2.51	0.43
1:A:204:ILE:HG23	1:A:244:LEU:HD11	2.01	0.43
1:B:362:GLU:O	1:B:366:GLN:HG3	2.19	0.43
2:C:103:LYS:O	2:C:107:ARG:HG3	2.19	0.42
1:A:112:MET:SD	1:A:199:LEU:HD11	2.59	0.42
1:A:358:ASN:HA	1:A:359:PRO:HD3	1.90	0.42
2:C:107:ARG:HA	2:C:111:ASP:HB2	2.01	0.42
1:A:156:TYR:CZ	1:A:192:ARG:HD3	2.54	0.42
1:B:92:LEU:HD13	1:B:290:TRP:CE2	2.55	0.42
1:A:297:LYS:O	1:A:301:GLU:HG2	2.20	0.41
1:A:303:GLN:HG3	1:A:306:LEU:HD12	2.03	0.41
1:B:161:THR:O	1:B:163:ARG:N	2.49	0.41
1:B:313:LEU:HD12	1:B:313:LEU:HA	1.81	0.41
1:A:90:GLU:OE2	1:A:98:ARG:NH2	2.54	0.41
1:B:303:GLN:HE22	1:B:381:GLU:HA	1.86	0.41
2:C:142:GLU:HB2	2:C:148:LEU:HD21	2.03	0.40
1:A:10:ARG:HB2	1:A:29:LEU:HD21	2.03	0.40
1:A:106:ILE:HG23	1:A:273:PHE:HB3	2.02	0.40
1:A:337:MLY:O	1:A:341:MET:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	359/395 (91%)	345 (96%)	13 (4%)	1 (0%)	41	71
1	B	359/395 (91%)	348 (97%)	9 (2%)	2 (1%)	25	58
2	C	97/105 (92%)	95 (98%)	2 (2%)	0	100	100
2	D	97/105 (92%)	91 (94%)	6 (6%)	0	100	100
All	All	912/1000 (91%)	879 (96%)	30 (3%)	3 (0%)	41	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	379	PRO
1	B	379	PRO
1	B	174	THR

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	324/342 (95%)	317 (98%)	7 (2%)	52	81
1	B	325/342 (95%)	319 (98%)	6 (2%)	59	85
2	C	95/99 (96%)	92 (97%)	3 (3%)	39	73
2	D	93/99 (94%)	93 (100%)	0	100	100
All	All	837/882 (95%)	821 (98%)	16 (2%)	57	84

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	176	THR
1	A	213	TRP
1	A	219	THR
1	A	305	ASP
1	A	353	ASN
1	A	354	GLU
2	C	109	LEU
2	C	111	ASP
2	C	117	CYS
1	B	22	ASP
1	B	132	ASP
1	B	160	ASN
1	B	161	THR
1	B	283	HIS
1	B	354	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	GLN
1	A	358	ASN
1	A	366	GLN
1	B	214	GLN
1	B	303	GLN
1	B	316	ASN
2	D	145	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

22 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	B	365	1	9,10,11	0.50	0	6,11,13	0.78	0
1	MLY	A	337	1	9,10,11	0.49	0	6,11,13	0.88	0
2	MLY	D	166	2	9,10,11	0.48	0	6,11,13	0.79	0
2	MLY	C	166	2	9,10,11	0.50	0	6,11,13	0.83	0
1	MLY	B	155	1	9,10,11	0.49	0	6,11,13	0.85	0
1	MLY	A	155	1	9,10,11	0.49	0	6,11,13	0.83	0
1	MLY	B	256	1	9,10,11	0.51	0	6,11,13	0.80	0
1	MLY	B	154	1	9,10,11	0.52	0	6,11,13	0.80	0
1	MLY	A	365	1	9,10,11	0.50	0	6,11,13	0.81	0
1	MLY	B	8	1	9,10,11	0.51	0	6,11,13	0.88	0
1	MLY	A	263	1	9,10,11	0.52	0	6,11,13	0.82	0
1	MLY	A	154	1	9,10,11	0.49	0	6,11,13	0.79	0
1	MLY	A	8	1	9,10,11	0.57	0	6,11,13	0.77	0
1	MLY	B	352	1	9,10,11	0.51	0	6,11,13	0.81	0
1	MLY	A	352	1	9,10,11	0.52	0	6,11,13	0.83	0
1	MLY	B	231	1	9,10,11	0.53	0	6,11,13	0.66	0
2	MLY	D	155	2	9,10,11	0.53	0	6,11,13	0.65	0
1	MLY	B	263	1	9,10,11	0.57	0	6,11,13	0.62	0
1	MLY	A	231	1	9,10,11	0.51	0	6,11,13	0.83	0
1	MLY	A	256	1	9,10,11	0.53	0	6,11,13	0.77	0
2	MLY	C	155	2	9,10,11	0.51	0	6,11,13	0.77	0
1	MLY	B	337	1	9,10,11	0.49	0	6,11,13	0.83	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
1	MLY	B	365	1	-	0/8/9/11	-
1	MLY	A	337	1	-	0/8/9/11	-
2	MLY	D	166	2	-	2/8/9/11	-
2	MLY	C	166	2	-	0/8/9/11	-
1	MLY	B	155	1	-	0/8/9/11	-
1	MLY	A	155	1	-	0/8/9/11	-
1	MLY	B	256	1	-	1/8/9/11	-
1	MLY	B	154	1	-	0/8/9/11	-
1	MLY	A	365	1	-	0/8/9/11	-
1	MLY	B	8	1	-	0/8/9/11	-
1	MLY	A	263	1	-	0/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	A	154	1	-	1/8/9/11	-
1	MLY	A	8	1	-	0/8/9/11	-
1	MLY	B	352	1	-	5/8/9/11	-
1	MLY	A	352	1	-	3/8/9/11	-
1	MLY	B	231	1	-	1/8/9/11	-
2	MLY	D	155	2	-	1/8/9/11	-
1	MLY	B	263	1	-	0/8/9/11	-
1	MLY	A	231	1	-	1/8/9/11	-
1	MLY	A	256	1	-	4/8/9/11	-
2	MLY	C	155	2	-	0/8/9/11	-
1	MLY	B	337	1	-	0/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	154	MLY	O-C-CA-CB
1	A	352	MLY	N-CA-CB-CG
1	A	352	MLY	C-CA-CB-CG
1	B	352	MLY	C-CA-CB-CG
2	D	166	MLY	O-C-CA-CB
1	B	352	MLY	CD-CE-NZ-CH2
1	B	231	MLY	CA-CB-CG-CD
1	B	352	MLY	CA-CB-CG-CD
1	A	256	MLY	CD-CE-NZ-CH2
1	A	352	MLY	CA-CB-CG-CD
2	D	155	MLY	CD-CE-NZ-CH2
1	B	256	MLY	CA-CB-CG-CD
1	A	256	MLY	CD-CE-NZ-CH1
1	B	352	MLY	CD-CE-NZ-CH1
1	A	256	MLY	CE-CD-CG-CB
1	B	352	MLY	CE-CD-CG-CB
1	A	231	MLY	CD-CE-NZ-CH2
1	A	256	MLY	CA-CB-CG-CD
2	D	166	MLY	CA-CB-CG-CD

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	337	MLY	1	0
1	B	231	MLY	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	365/395 (92%)	-0.04	10 (2%)	54	50	39, 72, 162, 268	0
1	B	365/395 (92%)	0.14	23 (6%)	20	16	47, 78, 184, 241	0
2	C	99/105 (94%)	-0.09	2 (2%)	65	63	52, 79, 145, 175	0
2	D	99/105 (94%)	0.07	2 (2%)	65	63	61, 101, 143, 165	0
All	All	928/1000 (92%)	0.04	37 (3%)	38	33	39, 79, 170, 268	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	176	THR	6.3
1	A	20	PHE	6.1
1	A	381	GLU	5.9
1	B	316	ASN	5.9
1	B	175	GLY	5.6
1	B	21	GLY	5.1
1	B	314	ASN	4.8
1	B	311	GLY	4.2
2	C	191	ASN	4.1
1	B	22	ASP	4.0
2	D	191	ASN	4.0
1	B	307	LEU	3.9
1	B	305	ASP	3.9
1	B	174	THR	3.9
1	A	386	LEU	3.8
1	B	312	ASP	3.5
1	B	381	GLU	3.4
1	A	176	THR	3.4
1	A	234	ASN	3.0
1	A	311	GLY	3.0
1	B	20	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	387	ASP	2.9
1	B	306	LEU	2.9
2	C	126	ASP	2.9
1	B	317	SER	2.8
1	B	376	SER	2.6
1	B	303	GLN	2.5
1	A	382	LEU	2.4
1	A	306	LEU	2.4
1	A	19	SER	2.3
1	B	309	GLU	2.3
1	B	313	LEU	2.3
2	D	172	LEU	2.2
1	B	29	LEU	2.2
1	B	385	ALA	2.1
1	B	169	GLU	2.1
1	A	229	MET	2.1

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MLY	D	166	11/12	0.86	0.27	88,103,115,116	0
1	MLY	A	352	11/12	0.88	0.20	98,111,120,121	0
1	MLY	B	365	11/12	0.90	0.21	74,88,95,96	0
1	MLY	A	154	11/12	0.91	0.23	49,56,92,95	0
2	MLY	C	166	11/12	0.91	0.23	63,73,104,107	0
1	MLY	A	337	11/12	0.92	0.17	74,81,87,91	0
2	MLY	D	155	11/12	0.92	0.16	99,104,114,115	0
1	MLY	B	8	11/12	0.92	0.15	95,101,109,112	0
1	MLY	B	231	11/12	0.93	0.20	88,93,97,102	0
1	MLY	B	256	11/12	0.93	0.16	61,69,89,90	0
1	MLY	B	352	11/12	0.93	0.19	62,72,89,90	0
1	MLY	A	365	11/12	0.93	0.30	91,103,114,118	0
1	MLY	A	231	11/12	0.93	0.20	92,97,106,107	0
1	MLY	A	256	11/12	0.93	0.17	65,75,90,91	0
1	MLY	B	154	11/12	0.94	0.23	57,67,99,100	0
1	MLY	B	337	11/12	0.94	0.24	78,88,100,101	0
2	MLY	C	155	11/12	0.94	0.25	72,86,98,101	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
1	MLY	A	8	11/12	0.95	0.27	71,75,104,106	0
1	MLY	A	263	11/12	0.96	0.13	58,69,87,92	0
1	MLY	A	155	11/12	0.96	0.23	43,45,61,64	0
1	MLY	B	263	11/12	0.96	0.15	79,86,117,118	0
1	MLY	B	155	11/12	0.96	0.34	61,67,79,83	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ZN	C	201	1/1	0.89	0.06	83,83,83,83	0
3	ZN	D	201	1/1	0.96	0.03	88,88,88,88	0

6.5 Other polymers (i)

There are no such residues in this entry.