



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 08:30 PM GMT

PDB ID : 4Y68
Title : Structure of a lipoprotein from Streptococcus agalactiae
Authors : Khosa, S.; Hoepfner, A.; Smits, S.H.
Deposited on : 2015-02-12
Resolution : 2.21 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

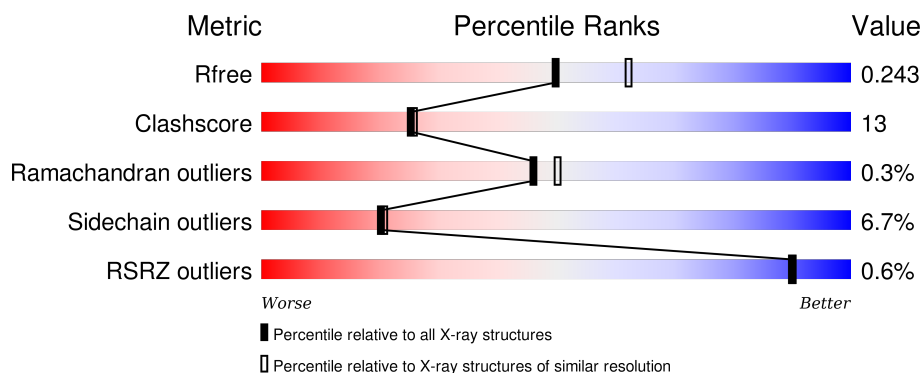
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (2.24-2.20)

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	312	<div> <div>69%</div> <div>21%</div> <div>8%</div> </div>
1	B	312	<div> <div>2%</div> <div>65%</div> <div>22%</div> <div>8%</div> </div>
1	C	312	<div> <div>79%</div> <div>11%</div> <div>8%</div> </div>
1	D	312	<div> <div>74%</div> <div>16%</div> <div>8%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MES	C	401	-	-	X	X
2	MES	C	402	-	-	-	X
2	MES	D	401	-	-	X	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9601 atoms, of which 13 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 Putative nisin-resistance protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	287	Total	C	N	O	S	0	0	0
			2295	1450	408	431	6			
1	A	287	Total	C	N	O	S	0	0	0
			2251	1424	391	430	6			
1	B	286	Total	C	N	O	S	0	0	0
			2186	1378	379	423	6			
1	D	287	Total	C	N	O	S	0	0	0
			2285	1445	405	429	6			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	9	MET	-	expression tag	UNP Q3DLX7
C	10	GLY	-	expression tag	UNP Q3DLX7
C	11	SER	-	expression tag	UNP Q3DLX7
C	12	SER	-	expression tag	UNP Q3DLX7
C	13	HIS	-	expression tag	UNP Q3DLX7
C	14	HIS	-	expression tag	UNP Q3DLX7
C	15	HIS	-	expression tag	UNP Q3DLX7
C	16	HIS	-	expression tag	UNP Q3DLX7
C	17	HIS	-	expression tag	UNP Q3DLX7
C	18	HIS	-	expression tag	UNP Q3DLX7
C	19	HIS	-	expression tag	UNP Q3DLX7
C	20	HIS	-	expression tag	UNP Q3DLX7
C	21	SER	-	expression tag	UNP Q3DLX7
C	22	SER	-	expression tag	UNP Q3DLX7
C	23	GLY	-	expression tag	UNP Q3DLX7
C	24	LEU	-	expression tag	UNP Q3DLX7
C	25	VAL	-	expression tag	UNP Q3DLX7
C	26	PRO	-	expression tag	UNP Q3DLX7
C	27	ARG	-	expression tag	UNP Q3DLX7
C	28	GLY	-	expression tag	UNP Q3DLX7
C	29	SER	-	expression tag	UNP Q3DLX7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	30	HIS	-	expression tag	UNP Q3DLX7
A	9	MET	-	expression tag	UNP Q3DLX7
A	10	GLY	-	expression tag	UNP Q3DLX7
A	11	SER	-	expression tag	UNP Q3DLX7
A	12	SER	-	expression tag	UNP Q3DLX7
A	13	HIS	-	expression tag	UNP Q3DLX7
A	14	HIS	-	expression tag	UNP Q3DLX7
A	15	HIS	-	expression tag	UNP Q3DLX7
A	16	HIS	-	expression tag	UNP Q3DLX7
A	17	HIS	-	expression tag	UNP Q3DLX7
A	18	HIS	-	expression tag	UNP Q3DLX7
A	19	HIS	-	expression tag	UNP Q3DLX7
A	20	HIS	-	expression tag	UNP Q3DLX7
A	21	SER	-	expression tag	UNP Q3DLX7
A	22	SER	-	expression tag	UNP Q3DLX7
A	23	GLY	-	expression tag	UNP Q3DLX7
A	24	LEU	-	expression tag	UNP Q3DLX7
A	25	VAL	-	expression tag	UNP Q3DLX7
A	26	PRO	-	expression tag	UNP Q3DLX7
A	27	ARG	-	expression tag	UNP Q3DLX7
A	28	GLY	-	expression tag	UNP Q3DLX7
A	29	SER	-	expression tag	UNP Q3DLX7
A	30	HIS	-	expression tag	UNP Q3DLX7
B	9	MET	-	expression tag	UNP Q3DLX7
B	10	GLY	-	expression tag	UNP Q3DLX7
B	11	SER	-	expression tag	UNP Q3DLX7
B	12	SER	-	expression tag	UNP Q3DLX7
B	13	HIS	-	expression tag	UNP Q3DLX7
B	14	HIS	-	expression tag	UNP Q3DLX7
B	15	HIS	-	expression tag	UNP Q3DLX7
B	16	HIS	-	expression tag	UNP Q3DLX7
B	17	HIS	-	expression tag	UNP Q3DLX7
B	18	HIS	-	expression tag	UNP Q3DLX7
B	19	HIS	-	expression tag	UNP Q3DLX7
B	20	HIS	-	expression tag	UNP Q3DLX7
B	21	SER	-	expression tag	UNP Q3DLX7
B	22	SER	-	expression tag	UNP Q3DLX7
B	23	GLY	-	expression tag	UNP Q3DLX7
B	24	LEU	-	expression tag	UNP Q3DLX7
B	25	VAL	-	expression tag	UNP Q3DLX7
B	26	PRO	-	expression tag	UNP Q3DLX7
B	27	ARG	-	expression tag	UNP Q3DLX7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	28	GLY	-	expression tag	UNP Q3DLX7
B	29	SER	-	expression tag	UNP Q3DLX7
B	30	HIS	-	expression tag	UNP Q3DLX7
D	9	MET	-	expression tag	UNP Q3DLX7
D	10	GLY	-	expression tag	UNP Q3DLX7
D	11	SER	-	expression tag	UNP Q3DLX7
D	12	SER	-	expression tag	UNP Q3DLX7
D	13	HIS	-	expression tag	UNP Q3DLX7
D	14	HIS	-	expression tag	UNP Q3DLX7
D	15	HIS	-	expression tag	UNP Q3DLX7
D	16	HIS	-	expression tag	UNP Q3DLX7
D	17	HIS	-	expression tag	UNP Q3DLX7
D	18	HIS	-	expression tag	UNP Q3DLX7
D	19	HIS	-	expression tag	UNP Q3DLX7
D	20	HIS	-	expression tag	UNP Q3DLX7
D	21	SER	-	expression tag	UNP Q3DLX7
D	22	SER	-	expression tag	UNP Q3DLX7
D	23	GLY	-	expression tag	UNP Q3DLX7
D	24	LEU	-	expression tag	UNP Q3DLX7
D	25	VAL	-	expression tag	UNP Q3DLX7
D	26	PRO	-	expression tag	UNP Q3DLX7
D	27	ARG	-	expression tag	UNP Q3DLX7
D	28	GLY	-	expression tag	UNP Q3DLX7
D	29	SER	-	expression tag	UNP Q3DLX7
D	30	HIS	-	expression tag	UNP Q3DLX7

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	C	1	Total	C	H	N	O	S	0
			25	6	13	1	4	1	
2	A	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		

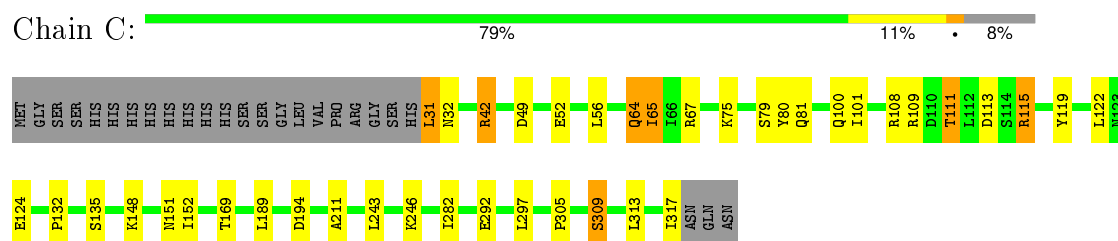
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	193	Total	O	0	0
			193	193		
3	A	110	Total	O	0	0
			110	110		
3	B	73	Total	O	0	0
			73	73		
3	D	147	Total	O	0	0
			147	147		

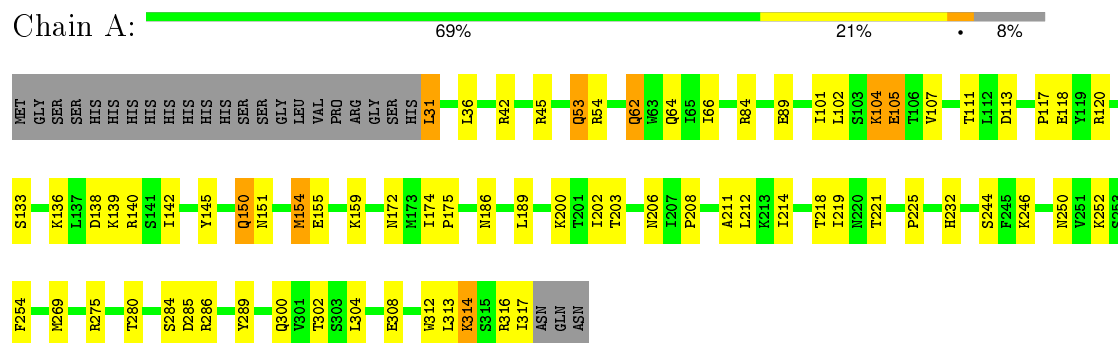
3 Residue-property plots

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

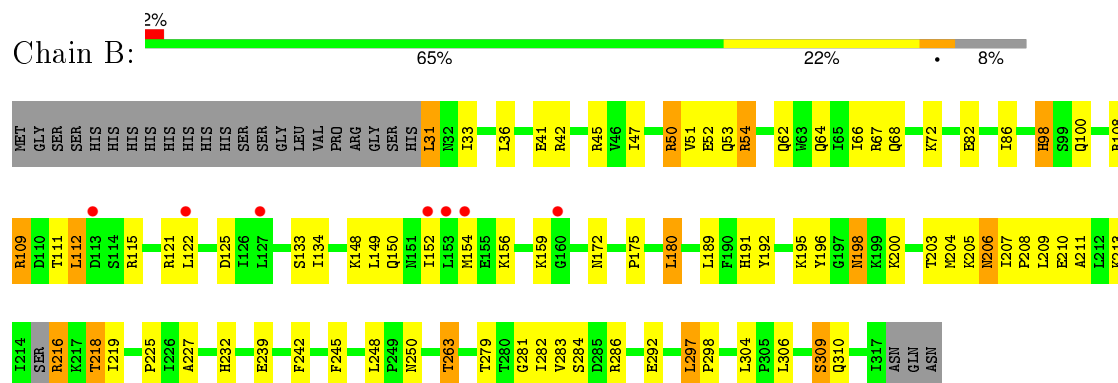
- Molecule 1: Putative nisin-resistance protein



- Molecule 1: Putative nisin-resistance protein

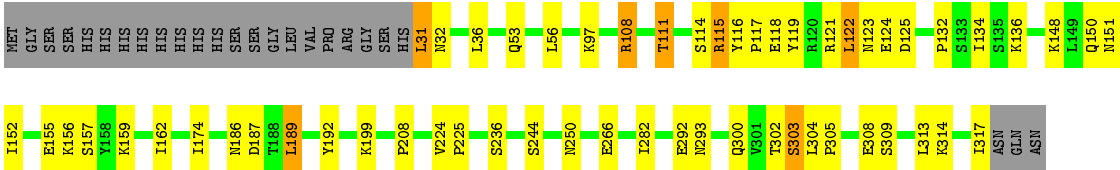


- Molecule 1: Putative nisin-resistance protein



- Molecule 1: Putative nisin-resistance protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.81Å 137.24Å 164.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	55.36 – 2.21 55.36 – 2.21	Depositor EDS
% Data completeness (in resolution range)	99.6 (55.36-2.21) 99.6 (55.36-2.21)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.194 , 0.243 0.194 , 0.243	Depositor DCC
R_{free} test set	2001 reflections (2.98%)	DCC
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 67172 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9601	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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: MES

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	0.38	0/2291	0.54	0/3100
1	B	0.34	0/2224	0.52	0/3018
1	C	0.47	0/2335	0.59	0/3149
1	D	0.44	0/2325	0.59	0/3137
All	All	0.41	0/9175	0.56	0/12404

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	2251	0	2272	64	0
1	B	2186	0	2138	77	0
1	C	2295	0	2366	40	0
1	D	2285	0	2351	57	0
2	A	12	0	12	0	0
2	C	24	13	24	8	0
2	D	12	0	12	9	0
3	A	110	0	0	9	0
3	B	73	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	193	0	0	13	1
3	D	147	0	0	7	1
All	All	9588	13	9175	229	1

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 13.

All (229) 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:D:121:ARG:HE	2:D:401:MES:H52	1.09	1.16
1:C:42:ARG:NH2	3:C:607:HOH:O	1.89	1.03
1:A:53:GLN:HG2	1:A:54:ARG:HG2	1.47	0.96
1:C:49:ASP:OD2	3:C:607:HOH:O	1.83	0.95
1:C:80:TYR:H	2:C:401:MES:H51	1.27	0.95
1:C:124:GLU:OE1	3:C:676:HOH:O	1.87	0.91
1:B:82:GLU:OE1	3:B:467:HOH:O	1.89	0.90
1:D:121:ARG:NE	2:D:401:MES:H52	1.88	0.89
1:D:300:GLN:NE2	1:D:308:GLU:OE1	2.11	0.83
1:B:50:ARG:HG2	1:B:50:ARG:HH11	1.44	0.82
1:B:111:THR:HG23	1:B:112:LEU:HD13	1.60	0.82
1:B:115:ARG:CB	1:D:115:ARG:HH12	1.93	0.81
1:C:64:GLN:NE2	3:C:501:HOH:O	1.91	0.81
1:A:139:LYS:HE2	1:A:211:ALA:O	1.80	0.80
1:A:202:ILE:HA	1:A:206:ASN:HD22	1.45	0.79
1:B:232:HIS:O	3:B:452:HOH:O	2.02	0.77
1:D:282:ILE:HD12	1:D:292:GLU:HG3	1.68	0.74
1:D:121:ARG:NH2	1:D:157:SER:O	2.18	0.74
1:A:316:ARG:NE	3:A:581:HOH:O	2.19	0.74
2:C:401:MES:O1S	2:C:401:MES:H32	1.87	0.74
1:A:31:LEU:HD11	1:B:100:GLN:HE22	1.53	0.73
1:C:108:ARG:O	1:C:111:THR:HG22	1.87	0.73
1:A:202:ILE:HA	1:A:206:ASN:ND2	2.03	0.73
1:C:81:GLN:NE2	3:C:655:HOH:O	1.97	0.73
1:A:118:GLU:OE1	3:A:594:HOH:O	2.05	0.73
1:B:108:ARG:O	1:B:111:THR:HG22	1.92	0.70
1:A:250:ASN:OD1	3:A:598:HOH:O	2.08	0.70
1:A:154:MET:CE	1:A:154:MET:HA	2.21	0.70
1:B:45:ARG:HG3	1:B:45:ARG:HH11	1.57	0.69
1:D:150:GLN:HG3	3:D:627:HOH:O	1.92	0.69
1:A:104:LYS:N	1:A:104:LYS:HD3	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:MET:SD	1:A:275:ARG:HG2	2.34	0.68
1:D:114:SER:HB3	1:D:134:ILE:HG13	1.77	0.67
1:A:31:LEU:CD1	1:B:100:GLN:HE22	2.08	0.67
1:A:45:ARG:NH1	3:A:595:HOH:O	2.27	0.67
1:C:313:LEU:O	1:C:317:ILE:HG12	1.95	0.66
1:A:232:HIS:HB3	3:A:562:HOH:O	1.96	0.66
1:D:53:GLN:OE1	3:D:637:HOH:O	2.13	0.66
1:A:154:MET:HA	1:A:154:MET:HE2	1.76	0.65
1:B:227:ALA:HB1	1:B:309:SER:HB3	1.79	0.64
1:B:225:PRO:HA	1:B:250:ASN:HB2	1.79	0.64
1:C:56:LEU:HD13	1:C:282:ILE:HD11	1.79	0.63
1:B:150:GLN:O	1:B:154:MET:HG2	1.98	0.63
1:A:138:ASP:O	1:A:142:ILE:HG12	1.97	0.63
1:B:209:LEU:HD13	1:B:219:ILE:HD11	1.80	0.63
1:C:246:LYS:NZ	3:C:649:HOH:O	2.31	0.63
1:C:115:ARG:HG3	1:C:132:PRO:HB2	1.82	0.62
1:B:41:GLU:O	1:B:45:ARG:HG2	1.99	0.62
1:B:200:LYS:HD3	3:B:444:HOH:O	1.99	0.62
1:A:64:GLN:NE2	3:A:585:HOH:O	2.14	0.62
1:D:121:ARG:NH2	2:D:401:MES:O2S	2.30	0.62
1:C:64:GLN:HB3	3:C:668:HOH:O	1.99	0.62
1:B:245:PHE:HA	1:B:248:LEU:HD12	1.80	0.62
1:A:53:GLN:NE2	3:A:550:HOH:O	2.33	0.62
1:C:100:GLN:HE22	1:D:31:LEU:HD13	1.65	0.61
1:B:156:LYS:O	3:B:445:HOH:O	2.16	0.61
1:C:56:LEU:HD13	1:C:282:ILE:CD1	2.31	0.61
1:A:174:ILE:HB	1:A:175:PRO:HD3	1.83	0.60
1:C:80:TYR:CE2	2:C:401:MES:H31	2.37	0.60
1:D:150:GLN:NE2	3:D:627:HOH:O	2.33	0.60
1:B:31:LEU:O	3:B:453:HOH:O	2.16	0.60
1:C:64:GLN:O	1:C:65:ILE:HB	2.01	0.60
1:A:136:LYS:O	1:A:142:ILE:HD11	2.02	0.60
1:D:115:ARG:HG3	1:D:132:PRO:HG3	1.84	0.60
1:B:200:LYS:NZ	3:B:444:HOH:O	2.10	0.59
1:B:148:LYS:O	1:B:152:ILE:HG13	2.02	0.59
1:D:313:LEU:O	1:D:317:ILE:HG13	2.03	0.59
1:D:156:LYS:HD3	2:D:401:MES:C7	2.32	0.59
1:A:203:THR:H	1:A:206:ASN:ND2	2.01	0.59
1:A:285:ASP:OD2	1:A:289:TYR:HB2	2.04	0.58
1:B:82:GLU:O	1:B:86:ILE:HG13	2.03	0.58
1:C:151:ASN:HB2	3:C:659:HOH:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:LYS:HA	1:B:216:ARG:HH12	1.69	0.57
1:B:47:ILE:O	1:B:51:VAL:HG23	2.04	0.57
1:C:79:SER:HB2	2:C:401:MES:H52	1.86	0.57
1:D:156:LYS:HD3	2:D:401:MES:H71	1.87	0.57
1:B:286:ARG:NE	3:B:458:HOH:O	2.31	0.56
1:B:50:ARG:HG2	1:B:50:ARG:NH1	2.16	0.56
1:B:192:TYR:HA	1:B:282:ILE:O	2.05	0.56
1:B:195:LYS:HG3	1:B:196:TYR:CE1	2.40	0.56
1:D:156:LYS:HD3	2:D:401:MES:H82	1.88	0.56
1:D:115:ARG:HD3	3:D:571:HOH:O	2.06	0.56
1:B:209:LEU:CD1	1:B:219:ILE:HD11	2.35	0.55
1:B:306:LEU:O	1:B:310:GLN:HG3	2.07	0.55
1:C:80:TYR:H	2:C:401:MES:C5	2.11	0.55
1:D:282:ILE:HD12	1:D:292:GLU:CG	2.37	0.55
1:D:305:PRO:O	1:D:309:SER:OG	2.19	0.55
1:C:52:GLU:OE1	1:C:67:ARG:NH1	2.39	0.54
1:A:31:LEU:HD12	1:A:31:LEU:N	2.22	0.54
1:A:107:VAL:O	1:A:111:THR:HG23	2.07	0.54
1:B:263:THR:HG22	1:B:281:GLY:O	2.08	0.54
1:A:107:VAL:HG21	1:A:275:ARG:HD2	1.90	0.54
1:C:64:GLN:O	1:C:65:ILE:CB	2.54	0.53
1:D:266:GLU:OE2	3:D:576:HOH:O	2.19	0.53
1:A:118:GLU:HG2	1:A:120:ARG:HH12	1.74	0.53
1:C:100:GLN:NE2	1:D:31:LEU:HD13	2.24	0.53
1:D:121:ARG:NH1	1:D:124:GLU:O	2.42	0.53
1:D:56:LEU:HD13	1:D:282:ILE:HD11	1.91	0.53
1:B:245:PHE:HA	1:B:248:LEU:CD1	2.38	0.53
1:A:139:LYS:O	1:A:139:LYS:HD3	2.09	0.52
1:A:159:LYS:HE3	1:A:159:LYS:HA	1.90	0.52
1:D:174:ILE:HG12	1:D:208:PRO:HD2	1.92	0.52
1:A:142:ILE:HG22	1:A:212:LEU:HD21	1.92	0.52
1:D:282:ILE:CD1	1:D:292:GLU:HG3	2.39	0.52
1:D:250:ASN:OD1	3:D:595:HOH:O	2.19	0.52
1:D:187:ASP:OD2	1:D:187:ASP:N	2.41	0.51
1:B:121:ARG:NH2	3:B:416:HOH:O	2.42	0.51
1:C:80:TYR:CD2	2:C:401:MES:H31	2.46	0.51
1:D:159:LYS:O	1:D:225:PRO:HD2	2.10	0.51
1:A:252:LYS:HG2	1:A:254:PHE:CZ	2.46	0.51
1:A:66:ILE:HD12	1:A:89:GLU:OE2	2.11	0.51
1:B:216:ARG:C	1:B:218:THR:H	2.13	0.51
1:A:218:THR:O	1:A:221:THR:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:LEU:CD2	1:B:180:LEU:HD13	2.41	0.51
1:A:107:VAL:HG21	1:A:275:ARG:CD	2.42	0.50
1:A:244:SER:C	1:A:286:ARG:HH22	2.14	0.50
1:A:62:GLN:O	1:A:66:ILE:HG12	2.11	0.50
1:B:62:GLN:O	1:B:66:ILE:HG13	2.11	0.50
1:C:109:ARG:NH2	3:C:636:HOH:O	2.45	0.50
1:D:123:ASN:OD1	1:D:125:ASP:N	2.30	0.50
1:C:31:LEU:HD23	1:C:32:ASN:H	1.77	0.49
1:B:149:LEU:HD21	1:B:180:LEU:HD13	1.93	0.49
1:C:56:LEU:CD1	1:C:282:ILE:HD12	2.43	0.49
1:B:204:MET:HG2	1:B:219:ILE:HD13	1.93	0.49
1:A:104:LYS:HE2	1:A:105:GLU:CB	2.43	0.49
1:A:186:ASN:HB2	1:A:219:ILE:HG21	1.95	0.49
1:C:64:GLN:NE2	3:C:503:HOH:O	2.45	0.49
1:A:302:THR:HG21	1:A:308:GLU:HG3	1.94	0.48
1:D:151:ASN:O	1:D:155:GLU:HB2	2.12	0.48
1:A:104:LYS:HD3	1:A:105:GLU:H	1.78	0.48
1:B:210:GLU:HG2	1:B:211:ALA:N	2.28	0.48
1:A:54:ARG:O	1:A:280:THR:HG22	2.14	0.47
1:A:174:ILE:HD12	1:A:208:PRO:HD2	1.96	0.47
1:B:210:GLU:O	1:B:213:LYS:N	2.47	0.47
1:B:239:GLU:O	1:B:242:PHE:HB3	2.14	0.47
1:B:198:ASN:OD1	1:B:198:ASN:N	2.47	0.47
1:B:203:THR:O	1:B:206:ASN:HB2	2.13	0.47
1:D:111:THR:HG23	1:D:136:LYS:HG2	1.97	0.47
1:A:159:LYS:CE	1:A:159:LYS:HA	2.44	0.47
1:C:108:ARG:NE	3:C:664:HOH:O	2.13	0.47
1:B:195:LYS:HG3	1:B:196:TYR:CD1	2.50	0.47
1:B:125:ASP:OD1	1:B:159:LYS:HE3	2.15	0.47
1:C:56:LEU:CD1	1:C:282:ILE:CD1	2.93	0.47
1:D:186:ASN:ND2	1:D:186:ASN:O	2.47	0.47
1:C:246:LYS:HB3	1:C:246:LYS:HE3	1.70	0.46
1:C:119:TYR:HB3	1:C:152:ILE:HD13	1.96	0.46
1:D:115:ARG:CG	1:D:132:PRO:HG3	2.43	0.46
1:A:104:LYS:HE2	1:A:105:GLU:HB3	1.98	0.46
1:D:119:TYR:OH	2:D:401:MES:H71	2.16	0.46
1:A:174:ILE:HB	1:A:175:PRO:CD	2.45	0.46
1:D:156:LYS:HD3	2:D:401:MES:C8	2.45	0.46
1:C:124:GLU:CD	1:C:124:GLU:H	2.18	0.46
1:B:108:ARG:O	1:B:112:LEU:HD22	2.16	0.46
1:B:45:ARG:NH1	1:B:45:ARG:HG3	2.26	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:ILE:CD1	1:B:292:GLU:HG3	2.45	0.46
1:B:53:GLN:HG2	1:B:54:ARG:HG2	1.97	0.46
1:A:155:GLU:HB2	3:A:503:HOH:O	2.16	0.46
1:A:174:ILE:HD11	1:B:36:LEU:HD12	1.98	0.46
1:D:116:TYR:HB3	1:D:117:PRO:HD2	1.98	0.46
1:B:98:HIS:HB3	1:B:279:THR:OG1	2.16	0.45
1:B:112:LEU:HA	1:B:112:LEU:HD12	1.69	0.45
1:D:115:ARG:HG3	1:D:116:TYR:N	2.28	0.45
1:A:104:LYS:H	1:A:104:LYS:HD3	1.78	0.45
1:D:192:TYR:HA	1:D:282:ILE:O	2.16	0.45
1:A:174:ILE:HD12	1:A:208:PRO:CD	2.46	0.45
1:B:297:LEU:HD22	1:B:298:PRO:HD2	1.99	0.45
1:B:50:ARG:HA	1:B:50:ARG:HD3	1.83	0.45
1:B:204:MET:HG3	1:B:209:LEU:HD11	1.98	0.45
1:D:199:LYS:HE3	3:D:607:HOH:O	2.16	0.45
1:B:227:ALA:CB	1:B:309:SER:HB3	2.46	0.45
1:A:142:ILE:HG22	1:A:212:LEU:CD2	2.47	0.45
1:D:97:LYS:HE2	1:D:293:ASN:HB3	1.97	0.44
1:A:314:LYS:HB2	1:A:314:LYS:NZ	2.32	0.44
1:B:204:MET:CG	1:B:219:ILE:HD13	2.46	0.44
1:A:84:ARG:HD2	1:A:101:ILE:HG21	1.99	0.44
1:A:31:LEU:HD11	1:B:100:GLN:NE2	2.28	0.44
1:D:189:LEU:HG	1:D:244:SER:HB3	1.99	0.44
1:B:191:HIS:HB2	1:B:284:SER:OG	2.18	0.44
1:A:312:TRP:CZ2	1:A:316:ARG:HG3	2.52	0.44
1:D:31:LEU:HD22	1:D:31:LEU:O	2.17	0.44
1:A:172:ASN:O	1:A:175:PRO:HD2	2.17	0.44
1:D:119:TYR:HB3	1:D:152:ILE:HD13	2.00	0.44
1:C:211:ALA:CB	2:C:402:MES:H82	2.48	0.43
1:A:316:ARG:CD	3:A:581:HOH:O	2.64	0.43
1:A:150:GLN:O	1:A:154:MET:HB2	2.17	0.43
1:A:174:ILE:HD11	1:B:36:LEU:CD1	2.48	0.43
1:A:118:GLU:HG2	1:A:120:ARG:NH1	2.33	0.43
2:C:402:MES:H31	2:C:402:MES:O2S	2.18	0.43
1:C:194:ASP:HB2	3:C:559:HOH:O	2.18	0.43
1:B:109:ARG:HE	1:B:109:ARG:HB2	1.68	0.43
1:B:72:LYS:HD3	1:B:72:LYS:C	2.38	0.43
1:D:303:SER:HB3	1:D:304:LEU:HD13	2.01	0.43
1:D:115:ARG:HG3	1:D:132:PRO:CG	2.48	0.43
1:B:250:ASN:N	1:B:250:ASN:OD1	2.41	0.43
1:B:152:ILE:O	1:B:156:LYS:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:LEU:HD11	1:C:282:ILE:HD12	2.00	0.43
1:B:125:ASP:HA	1:B:159:LYS:HE2	2.00	0.43
1:D:162:ILE:HG12	1:D:313:LEU:HD11	2.01	0.42
1:B:31:LEU:HA	1:B:31:LEU:HD13	1.86	0.42
1:B:31:LEU:HB3	1:B:33:ILE:HD11	1.99	0.42
1:B:282:ILE:HD12	1:B:292:GLU:HG3	2.01	0.42
1:B:72:LYS:HD3	1:B:72:LYS:O	2.20	0.42
1:B:50:ARG:NH1	1:B:50:ARG:CG	2.81	0.42
1:A:225:PRO:HB2	1:A:313:LEU:HD23	2.01	0.42
1:C:305:PRO:O	1:C:309:SER:HB2	2.19	0.42
1:C:81:GLN:CD	3:C:655:HOH:O	2.48	0.42
1:D:117:PRO:O	1:D:148:LYS:HE2	2.19	0.42
1:B:52:GLU:OE1	1:B:67:ARG:NH1	2.53	0.41
1:B:115:ARG:CB	1:D:115:ARG:HH22	2.33	0.41
1:B:149:LEU:HD12	1:B:149:LEU:HA	1.77	0.41
1:D:108:ARG:O	1:D:111:THR:HB	2.19	0.41
1:A:84:ARG:HD2	1:A:101:ILE:CG2	2.49	0.41
1:D:303:SER:O	1:D:304:LEU:HD12	2.21	0.41
1:C:169:THR:O	1:D:32:ASN:HA	2.20	0.41
1:A:313:LEU:O	1:A:317:ILE:HG12	2.19	0.41
1:C:243:LEU:HD23	1:C:243:LEU:HA	1.82	0.41
1:A:203:THR:H	1:A:206:ASN:HD22	1.68	0.41
1:B:64:GLN:O	1:B:68:GLN:HG3	2.21	0.41
1:B:172:ASN:O	1:B:175:PRO:HD2	2.20	0.41
1:C:282:ILE:HG13	1:C:292:GLU:HG3	2.03	0.41
1:A:252:LYS:HG2	1:A:254:PHE:CE1	2.56	0.41
1:B:283:VAL:HG12	1:B:284:SER:N	2.36	0.41
1:D:115:ARG:HD2	1:D:116:TYR:O	2.21	0.40
1:D:121:ARG:HD2	2:D:401:MES:H61	2.03	0.40
1:B:115:ARG:CB	1:D:115:ARG:NH1	2.74	0.40
1:B:207:ILE:O	1:B:209:LEU:N	2.53	0.40
1:A:117:PRO:HD3	1:A:145:TYR:CD2	2.57	0.40
1:D:159:LYS:O	1:D:224:VAL:HB	2.22	0.40
1:D:122:LEU:HA	1:D:122:LEU:HD12	1.55	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:526:HOH:O	3:D:524:HOH:O[4_475]	2.04	0.16

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	285/312 (91%)	272 (95%)	13 (5%)	0	100	100
1	B	282/312 (90%)	264 (94%)	17 (6%)	1 (0%)	39	41
1	C	285/312 (91%)	278 (98%)	6 (2%)	1 (0%)	39	41
1	D	285/312 (91%)	277 (97%)	7 (2%)	1 (0%)	39	41
All	All	1137/1248 (91%)	1091 (96%)	43 (4%)	3 (0%)	46	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	65	ILE
1	D	236	SER
1	B	208	PRO

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	249/280 (89%)	227 (91%)	22 (9%)	12	11
1	B	234/280 (84%)	214 (92%)	20 (8%)	13	12
1	C	258/280 (92%)	244 (95%)	14 (5%)	27	31
1	D	256/280 (91%)	245 (96%)	11 (4%)	35	42
All	All	997/1120 (89%)	930 (93%)	67 (7%)	20	21

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	31	LEU
1	C	42	ARG
1	C	64	GLN
1	C	75	LYS
1	C	101	ILE
1	C	111	THR
1	C	113	ASP
1	C	115	ARG
1	C	122	LEU
1	C	135	SER
1	C	148	LYS
1	C	189	LEU
1	C	297	LEU
1	C	309	SER
1	A	31	LEU
1	A	36	LEU
1	A	42	ARG
1	A	53	GLN
1	A	62	GLN
1	A	102	LEU
1	A	104	LYS
1	A	105	GLU
1	A	113	ASP
1	A	133	SER
1	A	140	ARG
1	A	150	GLN
1	A	151	ASN
1	A	154	MET
1	A	189	LEU
1	A	200	LYS
1	A	214	ILE
1	A	246	LYS
1	A	284	SER
1	A	300	GLN
1	A	304	LEU
1	A	314	LYS
1	B	31	LEU
1	B	42	ARG
1	B	50	ARG
1	B	54	ARG
1	B	98	HIS
1	B	109	ARG
1	B	112	LEU

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Mol	Chain	Res	Type
1	B	122	LEU
1	B	133	SER
1	B	134	ILE
1	B	180	LEU
1	B	189	LEU
1	B	198	ASN
1	B	206	ASN
1	B	216	ARG
1	B	218	THR
1	B	263	THR
1	B	297	LEU
1	B	304	LEU
1	B	309	SER
1	D	31	LEU
1	D	36	LEU
1	D	108	ARG
1	D	111	THR
1	D	115	ARG
1	D	118	GLU
1	D	122	LEU
1	D	189	LEU
1	D	302	THR
1	D	303	SER
1	D	314	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	206	ASN
1	B	100	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 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	MES	A	401	-	12,12,12	2.07	1 (8%)	15,16,16	2.91	8 (53%)
2	MES	C	401	-	12,12,12	2.06	1 (8%)	15,16,16	2.97	9 (60%)
2	MES	C	402	-	12,12,12	2.26	2 (16%)	16,16,16	2.52	8 (50%)
2	MES	D	401	-	12,12,12	2.20	1 (8%)	15,16,16	2.98	7 (46%)

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
2	MES	C	401	-	-	0/6/14/14	0/1/1/1
2	MES	C	402	-	-	0/6/14/14	0/1/1/1
2	MES	D	401	-	-	0/6/14/14	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	MES	C8-S	-7.18	1.66	1.77
2	A	401	MES	C8-S	-6.82	1.67	1.77
2	C	401	MES	C8-S	-6.72	1.67	1.77
2	C	402	MES	C8-S	-6.41	1.66	1.78
2	C	402	MES	C5-N4	-2.25	1.44	1.49

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	402	MES	O1-C6-C5	-5.97	105.34	111.42
2	D	401	MES	C2-C3-N4	-5.36	101.91	110.11
2	C	401	MES	C6-C5-N4	-4.52	103.19	110.11
2	D	401	MES	C6-C5-N4	-3.80	104.30	110.11
2	A	401	MES	C2-C3-N4	-3.34	105.01	110.11
2	C	402	MES	C2-C3-N4	-3.26	105.95	109.97
2	C	402	MES	C6-C5-N4	-3.22	105.99	109.97
2	C	401	MES	C2-C3-N4	-2.97	105.57	110.11
2	A	401	MES	C6-C5-N4	-2.72	105.95	110.11
2	C	402	MES	O3S-S-O1S	-2.70	105.75	112.39
2	C	402	MES	O1-C2-C3	-2.50	108.87	111.42
2	C	401	MES	O2S-S-O1S	-2.37	107.25	113.96
2	C	401	MES	O3S-S-C8	2.16	109.47	104.99
2	C	401	MES	C7-N4-C5	2.19	116.02	111.25
2	A	401	MES	O3S-S-C8	2.37	109.92	104.99
2	C	402	MES	O3S-S-C8	2.57	112.89	105.91
2	D	401	MES	O1S-S-C8	2.71	108.78	106.87
2	C	401	MES	C7-N4-C3	2.71	117.14	111.25
2	A	401	MES	O1S-S-C8	2.87	108.90	106.87
2	C	402	MES	C5-N4-C3	2.88	116.02	109.75
2	C	402	MES	C7-N4-C3	3.05	119.46	111.62
2	D	401	MES	C7-N4-C3	3.38	118.60	111.25
2	C	401	MES	O1S-S-C8	3.44	109.30	106.87
2	D	401	MES	C5-N4-C3	3.48	116.65	108.87
2	A	401	MES	C7-N4-C3	3.94	119.83	111.25
2	A	401	MES	C7-N4-C5	4.12	120.22	111.25
2	D	401	MES	C7-N4-C5	4.31	120.62	111.25
2	C	401	MES	C5-N4-C3	4.35	118.60	108.87
2	A	401	MES	C5-N4-C3	4.80	119.62	108.87
2	D	401	MES	O2S-S-C8	5.50	110.76	106.87
2	A	401	MES	O2S-S-C8	5.69	110.89	106.87
2	C	401	MES	O2S-S-C8	6.33	111.34	106.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	MES	6	0
2	C	402	MES	2	0
2	D	401	MES	9	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	287/312 (91%)	-0.36	0 100 100	28, 41, 58, 66	0
1	B	286/312 (91%)	0.02	7 (2%) 62 61	35, 52, 71, 81	0
1	C	287/312 (91%)	-0.30	0 100 100	19, 29, 44, 57	0
1	D	287/312 (91%)	-0.33	0 100 100	20, 34, 54, 61	0
All	All	1147/1248 (91%)	-0.24	7 (0%) 90 90	19, 39, 61, 81	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	160	GLY	4.0
1	B	122	LEU	3.3
1	B	153	LEU	3.0
1	B	154	MET	2.7
1	B	152	ILE	2.6
1	B	127	LEU	2.5
1	B	113	ASP	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MES	C	402	12/12	0.82	0.27	7.27	25,43,54,61	0
2	MES	C	401	12/12	0.88	0.22	5.94	34,44,60,64	0
2	MES	D	401	12/12	0.81	0.28	3.50	58,68,81,88	0
2	MES	A	401	12/12	0.73	0.33	-	105,113,119,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.