Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 10:10 PM EDT

PDB ID	:	3DG7
Title	:	Crystal structure of muconate lactonizing enzyme from Mucobacterium Smegmatis complexed with muconolactone
Authors	:	Fedorov, A.A.; Fedorov, E.V.; Sakai, A.; Gerlt, J.A.; Almo, S.C.
Deposited on	:	2008-06-12
Resolution	:	2.00 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:

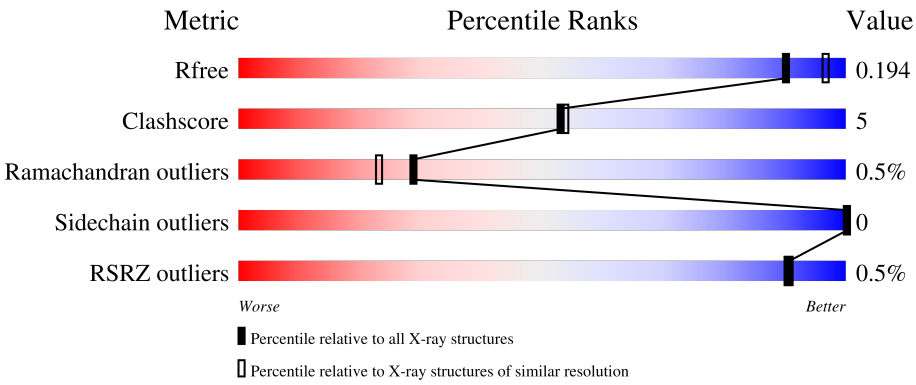
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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	367	% 90% 10%
1	B	367	89% 10%
1	C	367	% 89% 10%
1	D	367	89% 11%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12017 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.

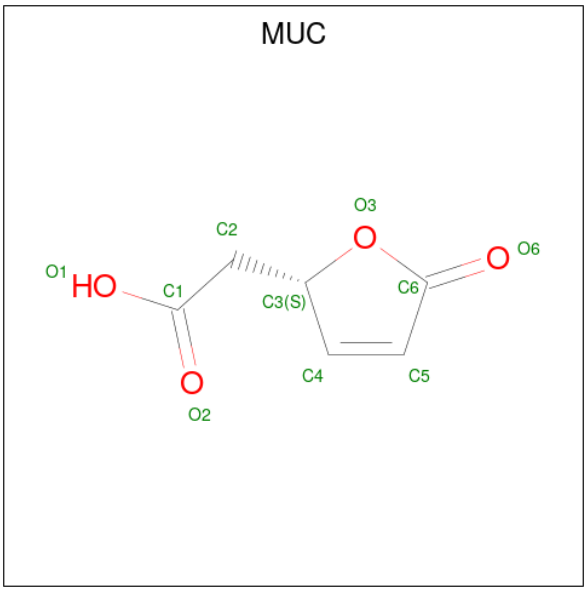
- Molecule 1 is a protein called Muconate cycloisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	366	2796	1750	500	530	16	0	0	0
1	B	366	2796	1750	500	530	16	0	0	0
1	C	366	2796	1750	500	530	16	0	0	0
1	D	366	2796	1750	500	530	16	0	0	0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0
2	D	1	1	1	0	0

- Molecule 3 is [(2S)-5-oxo-2,5-dihydrofuran-2-yl]acetic acid (three-letter code: MUC) (formula: $C_6H_6O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	10	6	4	0	0
			Total	C	O		
3	B	1	10	6	4	0	0
			Total	C	O		
3	C	1	10	6	4	0	0
			Total	C	O		
3	D	1	10	6	4	0	0

- Molecule 4 is water.

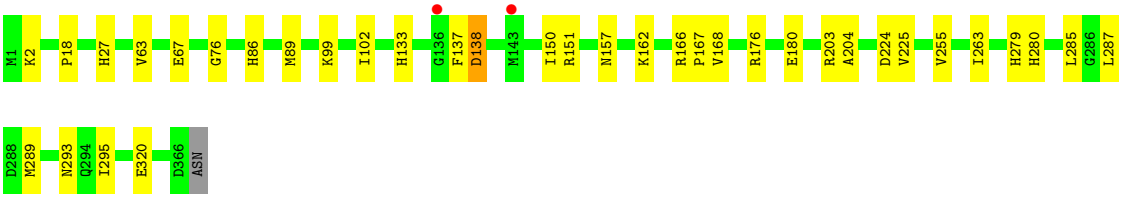
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	201	201	201	0	0
			Total	O		
4	B	198	198	198	0	0
			Total	O		
4	C	202	202	202	0	0
			Total	O		
4	D	188	188	188	0	0

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.

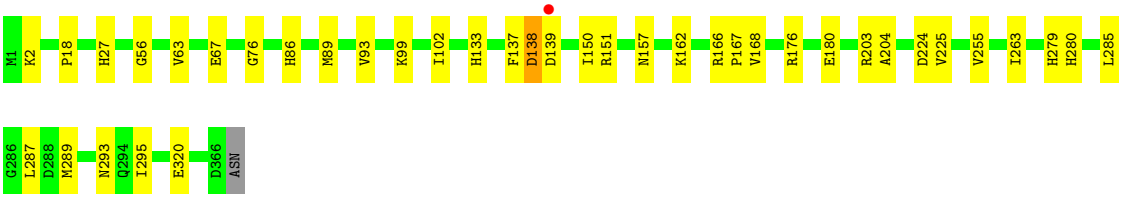
• Molecule 1: Muconate cycloisomerase

Chain A:



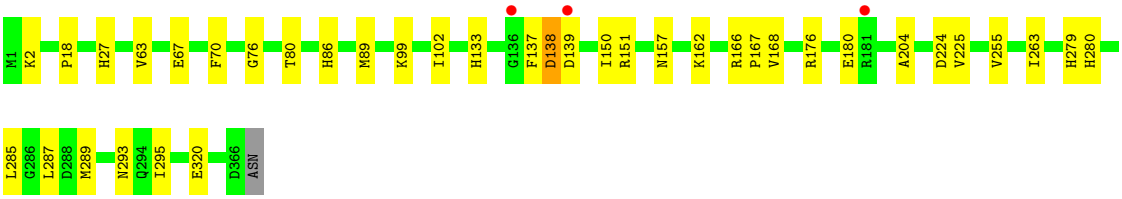
• Molecule 1: Muconate cycloisomerase

Chain B:



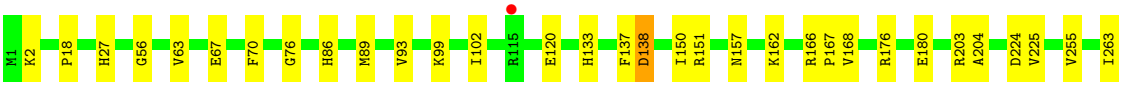
• Molecule 1: Muconate cycloisomerase

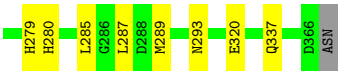
Chain C:



• Molecule 1: Muconate cycloisomerase

Chain D:




H279
H280
L285
G286
L287
D288
M289
N293
E320
Q337
D366
ASN

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	171.05Å 124.27Å 117.48Å 90.00° 133.39° 90.00°	Depositor
Resolution (Å)	24.42 – 2.00 39.31 – 1.99	Depositor EDS
% Data completeness (in resolution range)	96.5 (24.42-2.00) 95.9 (39.31-1.99)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.60 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.188 , 0.203 0.178 , 0.194	Depositor DCC
R_{free} test set	5835 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	19.3	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 28.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.460 for k+l,h+l,-l 0.457 for -k+l,-h-l,-l 0.459 for -h-2*l,-k,l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	12017	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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: MUC, MG

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	0.31	0/2847	0.60	1/3860 (0.0%)
1	B	0.31	0/2847	0.60	1/3860 (0.0%)
1	C	0.31	0/2847	0.60	1/3860 (0.0%)
1	D	0.31	0/2847	0.60	1/3860 (0.0%)
All	All	0.31	0/11388	0.60	4/15440 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	224	ASP	N-CA-C	-5.43	96.34	111.00
1	B	224	ASP	N-CA-C	-5.42	96.35	111.00
1	D	224	ASP	N-CA-C	-5.41	96.40	111.00
1	A	224	ASP	N-CA-C	-5.37	96.49	111.00

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	2796	0	2793	30	0
1	B	2796	0	2793	33	0
1	C	2796	0	2793	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2796	0	2793	33	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	10	0	5	2	0
3	B	10	0	5	2	0
3	C	10	0	5	2	0
3	D	10	0	5	2	0
4	A	201	0	0	4	0
4	B	198	0	0	1	0
4	C	202	0	0	1	0
4	D	188	0	0	3	0
All	All	12017	0	11192	122	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 5.

All (122) 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:B:89:MET:HE1	1:B:99:LYS:HA	1.55	0.87
1:A:89:MET:HE1	1:A:99:LYS:HA	1.55	0.86
1:D:89:MET:HE1	1:D:99:LYS:HA	1.59	0.83
1:C:89:MET:HE1	1:C:99:LYS:HA	1.61	0.82
1:A:151:ARG:HE	1:A:157:ASN:HD22	1.37	0.73
1:B:151:ARG:HE	1:B:157:ASN:HD22	1.36	0.72
1:D:151:ARG:HE	1:D:157:ASN:HD22	1.37	0.72
1:C:133:HIS:CD2	1:C:150:ILE:HD13	2.24	0.72
1:D:133:HIS:CD2	1:D:150:ILE:HD13	2.24	0.72
1:B:133:HIS:CD2	1:B:150:ILE:HD13	2.24	0.72
1:C:168:VAL:HG11	1:C:204:ALA:HB2	1.72	0.72
1:A:168:VAL:HG11	1:A:204:ALA:HB2	1.72	0.71
1:B:168:VAL:HG11	1:B:204:ALA:HB2	1.72	0.71
1:A:133:HIS:CD2	1:A:150:ILE:HD13	2.25	0.71
1:A:151:ARG:HE	1:A:157:ASN:ND2	1.88	0.71
1:B:151:ARG:HE	1:B:157:ASN:ND2	1.88	0.71
1:C:151:ARG:HE	1:C:157:ASN:ND2	1.89	0.71
1:D:151:ARG:HE	1:D:157:ASN:ND2	1.89	0.70
1:D:168:VAL:HG11	1:D:204:ALA:HB2	1.73	0.70
1:C:151:ARG:HE	1:C:157:ASN:HD22	1.37	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:MET:CE	1:C:99:LYS:HA	2.26	0.66
1:D:89:MET:CE	1:D:99:LYS:HA	2.27	0.65
1:A:166:ARG:HH21	1:A:166:ARG:HG3	1.62	0.65
1:B:166:ARG:HG3	1:B:166:ARG:HH21	1.62	0.64
1:A:89:MET:CE	1:A:99:LYS:HA	2.26	0.64
1:D:166:ARG:HH21	1:D:166:ARG:HG3	1.62	0.63
1:C:166:ARG:HH21	1:C:166:ARG:HG3	1.62	0.63
1:B:89:MET:CE	1:B:99:LYS:HA	2.26	0.62
1:A:279:HIS:ND1	1:A:280:HIS:HD2	1.99	0.61
1:C:279:HIS:ND1	1:C:280:HIS:HD2	2.00	0.60
1:B:176:ARG:O	1:B:180:GLU:HG3	2.01	0.60
1:B:279:HIS:ND1	1:B:280:HIS:HD2	2.00	0.60
1:A:176:ARG:O	1:A:180:GLU:HG3	2.01	0.60
1:D:176:ARG:O	1:D:180:GLU:HG3	2.01	0.60
1:C:176:ARG:O	1:C:180:GLU:HG3	2.01	0.60
1:D:279:HIS:ND1	1:D:280:HIS:HD2	1.99	0.60
1:D:255:VAL:HG11	1:D:287:LEU:HD11	1.85	0.58
1:B:255:VAL:HG11	1:B:287:LEU:HD11	1.85	0.58
1:C:255:VAL:HG11	1:C:287:LEU:HD11	1.86	0.58
1:A:162:LYS:NZ	3:A:1001:MUC:H2	2.18	0.57
1:A:255:VAL:HG11	1:A:287:LEU:HD11	1.86	0.57
1:C:2:LYS:HE2	1:C:76:GLY:HA2	1.86	0.57
1:D:2:LYS:HE2	1:D:76:GLY:HA2	1.87	0.57
1:A:2:LYS:HE2	1:A:76:GLY:HA2	1.86	0.57
1:A:168:VAL:CG1	1:A:204:ALA:HB2	2.35	0.56
1:B:168:VAL:CG1	1:B:204:ALA:HB2	2.36	0.56
1:B:2:LYS:HE2	1:B:76:GLY:HA2	1.86	0.56
1:C:168:VAL:CG1	1:C:204:ALA:HB2	2.36	0.55
1:D:86:HIS:HD2	4:D:2037:HOH:O	1.90	0.55
1:D:168:VAL:CG1	1:D:204:ALA:HB2	2.36	0.54
1:A:162:LYS:HZ1	3:A:1001:MUC:H2	1.72	0.54
1:D:162:LYS:HZ1	3:D:1004:MUC:H2	1.73	0.54
1:D:337:GLN:NE2	4:D:2186:HOH:O	2.38	0.53
1:A:151:ARG:NE	1:A:157:ASN:ND2	2.57	0.53
1:B:151:ARG:NE	1:B:157:ASN:ND2	2.57	0.53
1:B:137:PHE:O	1:B:138:ASP:HB2	2.09	0.52
1:D:89:MET:HE2	1:D:99:LYS:HG2	1.91	0.52
1:C:89:MET:HE2	1:C:99:LYS:HG2	1.91	0.52
1:A:137:PHE:O	1:A:138:ASP:HB2	2.09	0.52
4:A:2196:HOH:O	1:B:86:HIS:HE1	1.92	0.52
1:D:137:PHE:O	1:D:138:ASP:HB2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:HIS:HD2	4:A:2050:HOH:O	1.92	0.51
1:B:203:ARG:HD2	4:C:2122:HOH:O	2.10	0.51
1:B:162:LYS:NZ	3:B:1002:MUC:H2	2.26	0.51
4:A:2137:HOH:O	1:D:203:ARG:HD2	2.10	0.51
1:A:166:ARG:HG3	1:A:166:ARG:NH2	2.26	0.50
1:D:151:ARG:NE	1:D:157:ASN:ND2	2.57	0.50
1:A:89:MET:HE2	1:A:99:LYS:HG2	1.94	0.50
1:C:137:PHE:O	1:C:138:ASP:HB2	2.10	0.50
1:C:151:ARG:NE	1:C:157:ASN:ND2	2.57	0.50
1:B:89:MET:CE	1:B:102:ILE:HD12	2.42	0.50
1:B:86:HIS:HD2	4:B:2056:HOH:O	1.95	0.49
1:B:162:LYS:HZ1	3:B:1002:MUC:H2	1.76	0.49
1:C:89:MET:CE	1:C:102:ILE:HD12	2.43	0.49
1:B:166:ARG:HG3	1:B:166:ARG:NH2	2.26	0.49
1:D:162:LYS:NZ	3:D:1004:MUC:H2	2.28	0.49
1:B:89:MET:HE3	1:B:102:ILE:HD12	1.94	0.48
1:B:166:ARG:HA	1:B:167:PRO:C	2.34	0.48
1:D:89:MET:CE	1:D:102:ILE:HD12	2.43	0.48
1:A:89:MET:CE	1:A:102:ILE:HD12	2.43	0.48
1:C:166:ARG:HG3	1:C:166:ARG:NH2	2.26	0.48
1:A:89:MET:HE3	1:A:102:ILE:HD12	1.96	0.48
1:A:166:ARG:HA	1:A:167:PRO:C	2.34	0.48
1:C:285:LEU:HD21	1:D:285:LEU:HD21	1.96	0.47
1:D:166:ARG:HA	1:D:167:PRO:C	2.34	0.47
1:C:166:ARG:HA	1:C:167:PRO:C	2.34	0.47
1:D:293:ASN:HB2	1:D:320:GLU:HB3	1.97	0.47
1:A:293:ASN:HB2	1:A:320:GLU:HB3	1.97	0.47
1:B:293:ASN:HB2	1:B:320:GLU:HB3	1.96	0.47
1:B:89:MET:HE2	1:B:99:LYS:HG2	1.95	0.46
1:C:86:HIS:HE1	4:D:2136:HOH:O	1.98	0.46
1:C:162:LYS:NZ	3:C:1003:MUC:H2	2.31	0.46
1:A:285:LEU:HD21	1:B:285:LEU:HD21	1.98	0.45
1:C:293:ASN:HB2	1:C:320:GLU:HB3	1.97	0.45
1:A:63:VAL:O	1:A:67:GLU:HG3	2.16	0.45
1:D:166:ARG:HG3	1:D:166:ARG:NH2	2.27	0.45
1:D:63:VAL:O	1:D:67:GLU:HG3	2.17	0.44
1:C:63:VAL:O	1:C:67:GLU:HG3	2.18	0.44
1:B:63:VAL:O	1:B:67:GLU:HG3	2.17	0.43
1:A:203:ARG:HD2	4:A:2131:HOH:O	2.17	0.43
1:C:263:ILE:O	1:C:289:MET:HA	2.19	0.43
1:A:263:ILE:O	1:A:289:MET:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ARG:NE	1:A:157:ASN:HD22	2.12	0.43
1:C:18:PRO:HB3	1:C:27:HIS:ND1	2.33	0.43
1:D:263:ILE:O	1:D:289:MET:HA	2.19	0.43
1:D:18:PRO:HB3	1:D:27:HIS:ND1	2.33	0.42
1:A:18:PRO:HB3	1:A:27:HIS:ND1	2.34	0.42
1:B:18:PRO:HB3	1:B:27:HIS:ND1	2.34	0.42
1:B:56:GLY:HA3	1:D:93:VAL:O	2.20	0.42
1:B:263:ILE:O	1:B:289:MET:HA	2.20	0.41
1:C:162:LYS:HZ1	3:C:1003:MUC:H2	1.84	0.41
1:D:70:PHE:CG	1:D:89:MET:HE3	2.56	0.41
1:B:293:ASN:OD1	1:B:295:ILE:N	2.54	0.41
1:D:89:MET:HE3	1:D:102:ILE:HD12	2.01	0.41
1:A:293:ASN:OD1	1:A:295:ILE:N	2.54	0.41
1:B:138:ASP:OD1	1:B:139:ASP:N	2.51	0.41
1:C:293:ASN:OD1	1:C:295:ILE:N	2.53	0.41
1:C:138:ASP:OD1	1:C:139:ASP:N	2.51	0.41
1:C:89:MET:HE3	1:C:102:ILE:HD12	2.02	0.41
1:B:93:VAL:O	1:D:56:GLY:HA3	2.20	0.41
1:C:70:PHE:CG	1:C:89:MET:HE3	2.55	0.40
1:C:80:THR:HB	1:D:120:GLU:HB2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	364/367 (99%)	354 (97%)	8 (2%)	2 (0%)	29	23
1	B	364/367 (99%)	354 (97%)	8 (2%)	2 (0%)	29	23
1	C	364/367 (99%)	354 (97%)	8 (2%)	2 (0%)	29	23
1	D	364/367 (99%)	354 (97%)	8 (2%)	2 (0%)	29	23
All	All	1456/1468 (99%)	1416 (97%)	32 (2%)	8 (0%)	29	23

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	ASP
1	A	225	VAL
1	B	138	ASP
1	B	225	VAL
1	C	138	ASP
1	D	138	ASP
1	D	225	VAL
1	C	225	VAL

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	295/296 (100%)	295 (100%)	0	100	100
1	B	295/296 (100%)	295 (100%)	0	100	100
1	C	295/296 (100%)	295 (100%)	0	100	100
1	D	295/296 (100%)	295 (100%)	0	100	100
All	All	1180/1184 (100%)	1180 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	HIS
1	A	86	HIS
1	A	157	ASN
1	A	280	HIS
1	B	37	HIS
1	B	86	HIS
1	B	157	ASN
1	B	280	HIS
1	C	37	HIS
1	C	86	HIS

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Mol	Chain	Res	Type
1	C	157	ASN
1	C	280	HIS
1	D	37	HIS
1	D	86	HIS
1	D	157	ASN
1	D	280	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
3	MUC	B	1002	2	10,10,10	2.63	3 (30%)	13,13,13	3.65	5 (38%)
3	MUC	D	1004	2	10,10,10	2.65	4 (40%)	13,13,13	3.64	5 (38%)
3	MUC	C	1003	2	10,10,10	2.66	3 (30%)	13,13,13	3.63	5 (38%)
3	MUC	A	1001	2	10,10,10	2.62	3 (30%)	13,13,13	3.77	5 (38%)

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
3	MUC	B	1002	2	-	0/4/13/13	0/1/1/1
3	MUC	D	1004	2	-	0/4/13/13	0/1/1/1
3	MUC	C	1003	2	-	0/4/13/13	0/1/1/1
3	MUC	A	1001	2	-	0/4/13/13	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1003	MUC	C5-C6	-7.05	1.33	1.46
3	D	1004	MUC	C5-C6	-7.01	1.33	1.46
3	B	1002	MUC	C5-C6	-6.89	1.34	1.46
3	A	1001	MUC	C5-C6	-6.86	1.34	1.46
3	B	1002	MUC	C2-C3	2.33	1.57	1.54
3	C	1003	MUC	C2-C3	2.23	1.57	1.54
3	A	1001	MUC	C2-C3	2.23	1.57	1.54
3	D	1004	MUC	C3-C4	2.16	1.54	1.50
3	D	1004	MUC	C2-C3	2.10	1.57	1.54
3	C	1003	MUC	C2-C1	2.09	1.56	1.51
3	B	1002	MUC	C2-C1	2.06	1.56	1.51
3	A	1001	MUC	C2-C1	2.06	1.56	1.51
3	D	1004	MUC	C2-C1	2.01	1.56	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1001	MUC	C3-O3-C6	-8.65	102.20	109.16
3	D	1004	MUC	C3-O3-C6	-8.59	102.24	109.16
3	B	1002	MUC	C3-O3-C6	-8.52	102.31	109.16
3	C	1003	MUC	C3-O3-C6	-8.13	102.61	109.16
3	C	1003	MUC	O3-C6-C5	7.00	112.84	108.17
3	A	1001	MUC	O3-C6-C5	6.95	112.81	108.17
3	B	1002	MUC	O3-C6-C5	6.68	112.63	108.17
3	D	1004	MUC	O3-C6-C5	6.53	112.53	108.17
3	A	1001	MUC	C3-C4-C5	-5.47	104.19	110.45
3	B	1002	MUC	C3-C4-C5	-5.06	104.65	110.45
3	D	1004	MUC	C3-C4-C5	-5.03	104.69	110.45
3	C	1003	MUC	C3-C4-C5	-4.81	104.94	110.45
3	C	1003	MUC	O6-C6-C5	-4.20	123.90	130.85
3	A	1001	MUC	O6-C6-C5	-4.08	124.08	130.85
3	B	1002	MUC	O6-C6-C5	-4.02	124.18	130.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	1004	MUC	O6-C6-C5	-4.01	124.21	130.85
3	C	1003	MUC	O3-C3-C2	2.27	111.74	108.76
3	A	1001	MUC	O3-C3-C2	2.24	111.69	108.76
3	B	1002	MUC	O3-C3-C2	2.17	111.60	108.76
3	D	1004	MUC	O3-C3-C2	2.08	111.48	108.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1002	MUC	2	0
3	D	1004	MUC	2	0
3	C	1003	MUC	2	0
3	A	1001	MUC	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (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	366/367 (99%)	-0.38	2 (0%)	91	90	12, 19, 33, 44	0
1	B	366/367 (99%)	-0.42	1 (0%)	94	93	11, 19, 33, 44	0
1	C	366/367 (99%)	-0.42	3 (0%)	86	85	11, 19, 33, 44	0
1	D	366/367 (99%)	-0.38	1 (0%)	94	93	12, 19, 33, 44	0
All	All	1464/1468 (99%)	-0.40	7 (0%)	91	90	11, 19, 33, 44	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	ASP	3.1
1	A	136	GLY	2.7
1	A	143	MET	2.6
1	C	139	ASP	2.2
1	C	136	GLY	2.2
1	C	181	ARG	2.1
1	D	115	ARG	2.0

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

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

6.3 Carbohydrates (i)

There are no 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
2	MG	B	2002	1/1	0.85	0.08	20,20,20,20	0
2	MG	C	2003	1/1	0.88	0.09	20,20,20,20	0
3	MUC	A	1001	10/10	0.88	0.18	21,27,29,29	0
2	MG	D	2004	1/1	0.89	0.09	19,19,19,19	0
3	MUC	C	1003	10/10	0.89	0.21	19,27,29,29	0
3	MUC	B	1002	10/10	0.91	0.19	19,27,29,30	0
3	MUC	D	1004	10/10	0.93	0.14	20,26,29,29	0
2	MG	A	2001	1/1	0.96	0.06	20,20,20,20	0

6.5 Other polymers (i)

There are no such residues in this entry.