Full wwPDB X-ray Structure Validation Report (i)

Sep 17, 2023 – 09:01 AM EDT

PDB ID	:	4WMF
Title	:	Crystal structure of catalytically inactive MERS-CoV 3CL protease (C148A) in spacegroup P212121
Authors	:	Lountos, G.T.; Needle, D.; Waugh, D.S.
Deposited on	:	2014-10-08
Resolution	:	1.97 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:

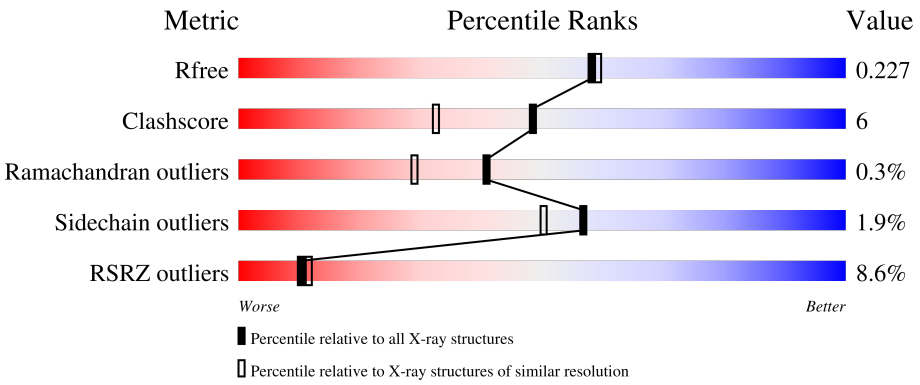
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 1.97 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	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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	306	5% 87% 11% ••
1	B	306	10% 89% 10% •
1	C	306	10% 85% 9% • 6%

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	PEG	A	403	-	-	X	-
3	PG4	B	406	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8040 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.

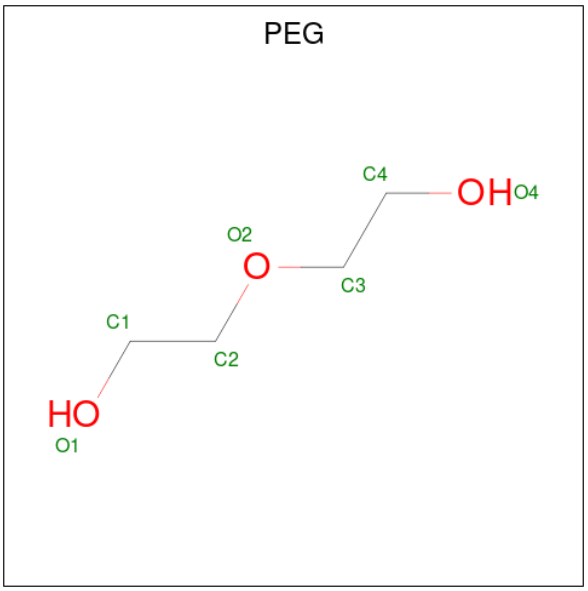
- Molecule 1 is a protein called MERS-CoV 3CL protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	304	2477	1566	414	467	30	0	21	0
1	B	303	2435	1543	407	458	27	7	16	0
1	C	289	2264	1437	370	431	26	0	11	0

There are 3 discrepancies between the modelled and reference sequences:

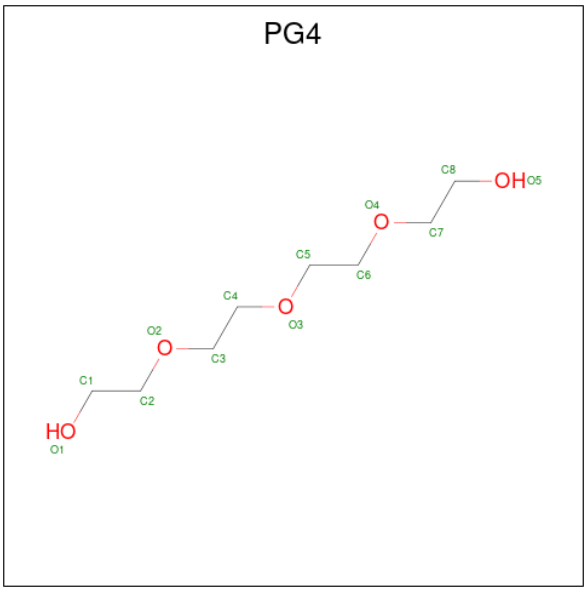
Chain	Residue	Modelled	Actual	Comment	Reference
A	148	ALA	CYS	engineered mutation	UNP W6A941
B	148	ALA	CYS	engineered mutation	UNP W6A941
C	148	ALA	CYS	engineered mutation	UNP W6A941

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O		0	0
			7	4	3			
2	A	1	Total	C	O		0	0
			7	4	3			
2	A	1	Total	C	O		0	0
			7	4	3			
2	B	1	Total	C	O		0	0
			7	4	3			
2	B	1	Total	C	O		0	0
			7	4	3			
2	C	1	Total	C	O		0	0
			7	4	3			

- Molecule 3 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	O		0	0
			13	8	5			
3	B	1	Total	C	O		0	0
			13	8	5			
3	B	1	Total	C	O		0	0
			13	8	5			
3	B	1	Total	C	O		0	0
			13	8	5			
3	C	1	Total	C	O		0	0
			12	8	4			

- Molecule 4 is water.

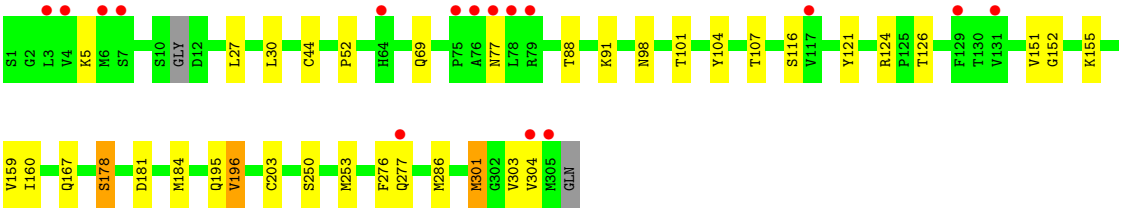
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	307	307	307	0	0
			Total	O		
4	B	253	253	253	0	0
			Total	O		
4	C	198	198	198	0	0

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.

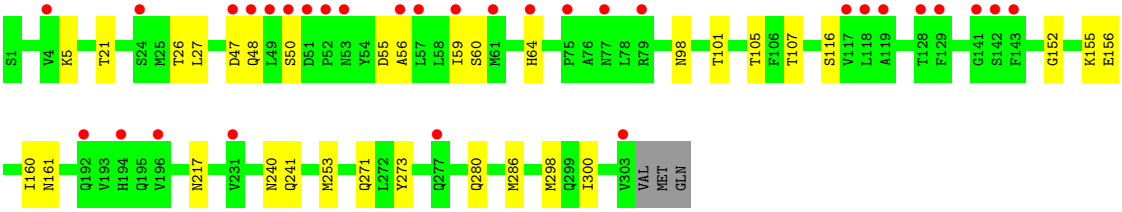
• Molecule 1: MERS-CoV 3CL protease

Chain A:



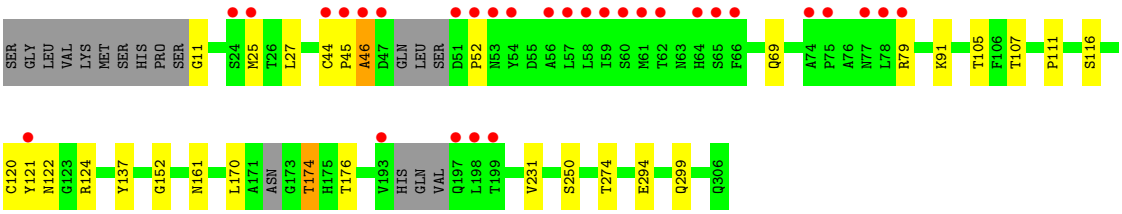
• Molecule 1: MERS-CoV 3CL protease

Chain B:



• Molecule 1: MERS-CoV 3CL protease

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	94.06Å 120.39Å 138.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.97 45.49 – 1.97	Depositor EDS
% Data completeness (in resolution range)	96.2 (50.00-1.97) 96.2 (45.49-1.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	2.33 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.5.0104	Depositor
R, R_{free}	0.193 , 0.226 0.194 , 0.227	Depositor DCC
R_{free} test set	5370 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	35.3	Xtriage
Anisotropy	0.123	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 49.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8040	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: PG4, PEG

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	0.88	1/2532 (0.0%)	0.80	2/3443 (0.1%)
1	B	0.84	0/2491	0.76	0/3391
1	C	0.76	0/2313	0.70	0/3147
All	All	0.83	1/7336 (0.0%)	0.75	2/9981 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	159	VAL	CB-CG2	6.27	1.66	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	124	ARG	NE-CZ-NH1	5.55	123.07	120.30
1	A	181	ASP	CB-CG-OD1	5.20	122.98	118.30

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	2477	0	2400	30	0
1	B	2435	0	2371	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2264	0	2175	24	0
2	A	21	0	30	5	0
2	B	14	0	20	1	0
2	C	7	0	10	0	0
3	B	52	0	72	11	0
3	C	12	0	15	4	0
4	A	307	0	0	13	0
4	B	253	0	0	10	0
4	C	198	0	0	8	0
All	All	8040	0	7093	86	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 6.

All (86) 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:240:ASN:O	1:B:241[B]:GLN:HG2	1.80	0.82
3:B:406:PG4:H62	4:B:550:HOH:O	1.82	0.79
4:A:548:HOH:O	3:B:406:PG4:H41	1.88	0.73
4:A:548:HOH:O	3:B:406:PG4:C4	2.36	0.73
1:B:105:THR:OG1	1:B:161[B]:ASN:ND2	2.23	0.71
1:C:121[B]:TYR:CE2	1:C:122:ASN:ND2	2.60	0.70
1:B:107:THR:OG1	1:B:161[A]:ASN:ND2	2.25	0.69
1:B:5[A]:LYS:HD2	4:B:624:HOH:O	1.94	0.68
1:A:91:LYS:HZ1	2:A:403:PEG:H32	1.59	0.67
3:B:406:PG4:C6	4:B:550:HOH:O	2.41	0.66
1:A:167[A]:GLN:OE1	1:A:178:SER:OG	2.14	0.64
4:A:539:HOH:O	3:B:402:PG4:H21	2.01	0.61
1:A:250[B]:SER:OG	4:A:501:HOH:O	2.13	0.60
1:B:155:LYS:HG2	1:B:160:ILE:HD13	1.82	0.60
1:C:11:GLY:N	4:C:504:HOH:O	2.34	0.60
1:B:298[B]:MET:SD	3:B:406:PG4:H62	2.41	0.60
1:C:124:ARG:NH1	4:C:503:HOH:O	2.32	0.59
1:C:105:THR:OG1	1:C:161[B]:ASN:ND2	2.33	0.58
1:A:126[B]:THR:HG23	4:A:548:HOH:O	2.02	0.58
1:B:271:GLN:NE2	4:B:503:HOH:O	2.36	0.58
1:C:274[A]:THR:HG22	1:C:274[A]:THR:O	2.04	0.58
4:A:548:HOH:O	3:B:406:PG4:H42	2.02	0.57
1:C:250[B]:SER:OG	4:C:501:HOH:O	2.17	0.57
1:A:304:VAL:HG12	4:A:666:HOH:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:298[B]:MET:HG2	3:B:406:PG4:O3	2.05	0.57
1:B:47:ASP:OD1	1:B:48:GLN:OE1	2.23	0.56
1:A:277[A]:GLN:O	1:A:277[A]:GLN:CG	2.53	0.56
1:C:170:LEU:HD11	1:C:176:THR:CG2	2.35	0.56
1:B:55:ASP:O	4:B:501:HOH:O	2.17	0.56
1:C:45:PRO:O	1:C:46:ALA:HB3	2.06	0.55
1:A:253[B]:MET:SD	1:A:303:VAL:HG11	2.47	0.54
1:C:231[B]:VAL:HG23	4:C:613:HOH:O	2.08	0.54
1:A:88:THR:HG21	1:A:184[B]:MET:CG	2.38	0.54
2:A:403:PEG:H32	4:A:694:HOH:O	2.06	0.54
1:B:59:ILE:HG23	4:B:501:HOH:O	2.08	0.53
1:B:241[B]:GLN:NE2	1:B:273:TYR:OH	2.41	0.53
1:A:107[A]:THR:OG1	4:A:502:HOH:O	2.19	0.53
1:A:30:LEU:HD13	1:A:151[A]:VAL:HG21	1.91	0.53
1:B:59:ILE:HD12	1:B:60:SER:N	2.23	0.52
1:B:155:LYS:HE2	1:B:160:ILE:CD1	2.38	0.52
1:B:56:ALA:O	1:B:59:ILE:HG13	2.09	0.52
1:A:253[B]:MET:HG3	4:A:630:HOH:O	2.10	0.50
1:A:91:LYS:HZ1	2:A:403:PEG:H21	1.78	0.49
1:C:174:THR:HG21	4:C:604:HOH:O	2.12	0.48
1:A:5:LYS:HE3	4:A:543:HOH:O	2.14	0.48
1:C:120[A]:CYS:C	1:C:121[A]:TYR:CD2	2.87	0.48
1:A:277[A]:GLN:O	1:A:277[A]:GLN:HG3	2.14	0.47
1:A:91:LYS:HZ1	2:A:403:PEG:C3	2.25	0.47
1:B:5[B]:LYS:HB3	1:B:5[B]:LYS:HE2	1.71	0.47
1:B:21:THR:HG23	1:B:26[B]:THR:HG22	1.96	0.47
1:A:301:MET:HB2	1:A:301:MET:HE3	1.66	0.47
1:C:294:GLU:OE2	4:C:501:HOH:O	2.20	0.47
1:B:155:LYS:HE2	1:B:160:ILE:HD11	1.96	0.46
1:C:69:GLN:NE2	4:C:515:HOH:O	2.48	0.46
1:C:111:PRO:HD3	1:C:137:TYR:CE1	2.51	0.46
1:A:155[A]:LYS:HE3	4:A:506:HOH:O	2.16	0.46
1:C:91:LYS:HZ1	3:C:402:PG4:C5	2.29	0.45
1:C:91:LYS:HZ3	3:C:402:PG4:H62	1.81	0.45
1:C:107:THR:HG21	1:C:161[A]:ASN:HD21	1.82	0.45
1:C:231[A]:VAL:HG22	4:C:613:HOH:O	2.16	0.45
1:A:126[B]:THR:HG23	3:B:406:PG4:H41	2.00	0.44
1:C:91:LYS:NZ	3:C:402:PG4:H72	2.33	0.44
1:A:116:SER:O	1:A:152:GLY:HA2	2.18	0.44
1:B:5[B]:LYS:HE2	4:B:668:HOH:O	2.17	0.43
1:A:276:PHE:O	1:A:277[A]:GLN:HG2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:25:MET:HG3	1:C:44:CYS:O	2.18	0.43
1:A:203[B]:CYS:SG	4:A:754:HOH:O	2.44	0.43
1:C:91:LYS:NZ	3:C:402:PG4:H62	2.33	0.43
1:A:69:GLN:HG2	1:A:77[A]:ASN:OD1	2.19	0.43
1:B:5[A]:LYS:CD	4:B:624:HOH:O	2.59	0.42
1:A:91:LYS:NZ	2:A:403:PEG:H21	2.34	0.42
1:A:155[B]:LYS:HG2	1:A:160:ILE:HD13	2.02	0.42
2:B:404:PEG:H11	4:B:705:HOH:O	2.20	0.42
1:A:44:CYS:SG	1:A:52:PRO:HB3	2.59	0.42
1:A:121:TYR:CD1	3:B:406:PG4:H21	2.55	0.42
1:C:116:SER:O	1:C:152:GLY:HA2	2.20	0.42
1:B:298[B]:MET:HE3	1:B:298[B]:MET:HB2	1.89	0.41
1:A:98:ASN:HB3	1:A:101:THR:OG1	2.20	0.41
1:B:98:ASN:HB3	1:B:101:THR:OG1	2.21	0.41
1:B:217:ASN:HD21	1:B:300:ILE:HD12	1.86	0.41
1:B:253:MET:HE2	4:B:745:HOH:O	2.21	0.41
1:C:45:PRO:O	1:C:46:ALA:CB	2.69	0.41
1:A:104:TYR:HA	1:A:160:ILE:O	2.21	0.40
1:A:196:VAL:HG21	1:C:299:GLN:CD	2.41	0.40
1:B:116:SER:O	1:B:152:GLY:HA2	2.21	0.40
1:A:121:TYR:CE1	3:B:406:PG4:H21	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	321/306 (105%)	316 (98%)	4 (1%)	1 (0%)	41 29
1	B	317/306 (104%)	315 (99%)	2 (1%)	0	100 100
1	C	292/306 (95%)	278 (95%)	12 (4%)	2 (1%)	22 11
All	All	930/918 (101%)	909 (98%)	18 (2%)	3 (0%)	41 29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	VAL
1	C	52	PRO
1	C	46	ALA

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	275/256 (107%)	270 (98%)	5 (2%)	59	51
1	B	269/256 (105%)	263 (98%)	6 (2%)	52	46
1	C	246/256 (96%)	243 (99%)	3 (1%)	71	67
All	All	790/768 (103%)	776 (98%)	14 (2%)	57	51

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	178	SER
1	A	195	GLN
1	A	286	MET
1	A	301	MET
1	B	27	LEU
1	B	50	SER
1	B	64	HIS
1	B	156	GLU
1	B	280	GLN
1	B	286	MET
1	C	27	LEU
1	C	79	ARG
1	C	174	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	122	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)

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)

11 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	$\#\|Z\| > 2$
3	PG4	B	402	-	12,12,12	0.60	0	11,11,11	0.53	0
2	PEG	A	401	-	6,6,6	0.32	0	5,5,5	0.24	0
3	PG4	B	401	-	12,12,12	0.68	0	11,11,11	0.26	0
2	PEG	A	402	-	6,6,6	0.39	0	5,5,5	0.29	0
2	PEG	B	404	-	6,6,6	0.62	0	5,5,5	0.44	0
2	PEG	A	403	-	6,6,6	0.60	0	5,5,5	0.74	0
3	PG4	B	405	-	12,12,12	0.41	0	11,11,11	0.37	0
2	PEG	B	403	-	6,6,6	0.34	0	5,5,5	0.41	0
2	PEG	C	401	-	6,6,6	0.43	0	5,5,5	0.50	0
3	PG4	B	406	-	12,12,12	0.56	0	11,11,11	0.81	0
3	PG4	C	402	-	11,11,12	0.80	0	10,10,11	0.95	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
3	PG4	B	402	-	-	3/10/10/10	-
2	PEG	A	401	-	-	3/4/4/4	-
3	PG4	B	401	-	-	1/10/10/10	-
2	PEG	A	402	-	-	3/4/4/4	-
2	PEG	B	404	-	-	3/4/4/4	-
2	PEG	A	403	-	-	3/4/4/4	-
3	PG4	B	405	-	-	3/10/10/10	-
2	PEG	B	403	-	-	2/4/4/4	-
2	PEG	C	401	-	-	3/4/4/4	-
3	PG4	B	406	-	-	3/10/10/10	-
3	PG4	C	402	-	-	3/9/9/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	PEG	O2-C3-C4-O4
2	B	404	PEG	O2-C3-C4-O4
2	A	403	PEG	C4-C3-O2-C2
2	A	401	PEG	O1-C1-C2-O2
3	B	402	PG4	O1-C1-C2-O2
2	A	401	PEG	O2-C3-C4-O4
2	B	403	PEG	O2-C3-C4-O4
3	B	405	PG4	O3-C5-C6-O4
2	A	402	PEG	O1-C1-C2-O2
3	B	402	PG4	O2-C3-C4-O3
2	C	401	PEG	O1-C1-C2-O2
2	B	404	PEG	O1-C1-C2-O2
2	A	403	PEG	O2-C3-C4-O4
3	B	406	PG4	O4-C7-C8-O5
3	B	406	PG4	O3-C5-C6-O4
3	C	402	PG4	O3-C5-C6-O4
3	B	401	PG4	O1-C1-C2-O2
3	C	402	PG4	C4-C3-O2-C2
2	C	401	PEG	C1-C2-O2-C3

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Mol	Chain	Res	Type	Atoms
2	C	401	PEG	C4-C3-O2-C2
2	A	402	PEG	C4-C3-O2-C2
3	C	402	PG4	C5-C6-O4-C7
3	B	406	PG4	C4-C3-O2-C2
2	B	403	PEG	O1-C1-C2-O2
2	A	403	PEG	O1-C1-C2-O2
3	B	405	PG4	C4-C3-O2-C2
2	B	404	PEG	C4-C3-O2-C2
3	B	402	PG4	C8-C7-O4-C6
2	A	401	PEG	C1-C2-O2-C3
3	B	405	PG4	C5-C6-O4-C7

There are no ring outliers.

5 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	PG4	1	0
2	B	404	PEG	1	0
2	A	403	PEG	5	0
3	B	406	PG4	10	0
3	C	402	PG4	4	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	304/306 (99%)	0.33	16 (5%)	26	28	20, 28, 47, 71	0
1	B	303/306 (99%)	0.55	31 (10%)	6	7	22, 33, 56, 67	0
1	C	289/306 (94%)	0.45	30 (10%)	6	7	25, 39, 69, 84	0
All	All	896/918 (97%)	0.44	77 (8%)	10	11	20, 34, 59, 84	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	198	LEU	8.1
1	C	59	ILE	7.6
1	A	304	VAL	6.4
1	B	303	VAL	6.4
1	C	45	PRO	5.5
1	C	197	GLN	5.5
1	C	57	LEU	5.1
1	A	77[A]	ASN	4.6
1	C	193	VAL	4.1
1	B	48	GLN	4.0
1	A	64[A]	HIS	3.8
1	B	51	ASP	3.7
1	B	50	SER	3.7
1	C	44	CYS	3.7
1	B	59	ILE	3.6
1	C	58	LEU	3.5
1	C	51	ASP	3.4
1	A	75	PRO	3.4
1	B	277	GLN	3.3
1	C	74	ALA	3.3
1	B	194	HIS	3.3
1	C	54	TYR	3.3
1	C	25	MET	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	199	THR	3.2
1	C	77	ASN	3.2
1	C	65	SER	3.2
1	B	118	LEU	3.2
1	C	56	ALA	3.2
1	C	64	HIS	3.1
1	C	60	SER	3.1
1	C	47	ASP	3.1
1	A	3	LEU	3.1
1	A	4	VAL	3.1
1	B	196	VAL	3.1
1	B	117	VAL	3.0
1	A	79[A]	ARG	2.9
1	B	129	PHE	2.9
1	C	24	SER	2.9
1	A	6	MET	2.9
1	B	53	ASN	2.9
1	C	61	MET	2.8
1	C	79	ARG	2.8
1	A	76	ALA	2.7
1	B	64	HIS	2.7
1	B	119	ALA	2.7
1	C	52	PRO	2.7
1	A	129	PHE	2.7
1	A	305	MET	2.7
1	B	4	VAL	2.6
1	B	128	THR	2.6
1	C	62	THR	2.6
1	C	46	ALA	2.6
1	B	192	GLN	2.6
1	A	277[A]	GLN	2.5
1	B	61	MET	2.5
1	C	121[A]	TYR	2.5
1	B	77	ASN	2.5
1	C	66	PHE	2.4
1	C	75	PRO	2.4
1	B	47	ASP	2.3
1	B	79	ARG	2.3
1	B	56	ALA	2.3
1	B	75	PRO	2.2
1	A	131	VAL	2.2
1	B	141	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	78	LEU	2.2
1	B	49	LEU	2.2
1	B	143	PHE	2.2
1	B	142	SER	2.1
1	A	117	VAL	2.1
1	B	231	VAL	2.1
1	B	24	SER	2.1
1	B	52	PRO	2.0
1	C	53	ASN	2.0
1	B	57	LEU	2.0
1	C	78	LEU	2.0
1	A	7	SER	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
3	PG4	B	405	13/13	0.61	0.30	81,83,85,85	0
3	PG4	C	402	12/13	0.64	0.25	61,64,66,66	0
3	PG4	B	406	13/13	0.76	0.21	54,64,66,66	0
2	PEG	A	403	7/7	0.78	0.25	48,55,62,65	0
2	PEG	C	401	7/7	0.82	0.23	58,61,65,65	0
2	PEG	B	404	7/7	0.84	0.15	47,48,50,53	0
2	PEG	A	402	7/7	0.86	0.15	40,47,51,53	0
2	PEG	B	403	7/7	0.87	0.18	54,55,58,59	0
3	PG4	B	402	13/13	0.89	0.24	44,50,57,57	0
3	PG4	B	401	13/13	0.92	0.14	29,40,44,48	0
2	PEG	A	401	7/7	0.93	0.17	43,47,55,56	0

6.5 Other polymers (i)

There are no such residues in this entry.