Full wwPDB X-ray Structure Validation Report (i)

Jun 22, 2024 – 01:54 PM EDT

PDB ID	:	4Y6T
Title	:	Structure of Tobacco streak virus coat protein dimer at 2.4 Angstroms resolution
Authors	:	Gulati, A.; Murthy, M.R.N.
Deposited on	:	2015-02-13
Resolution	:	2.40 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:

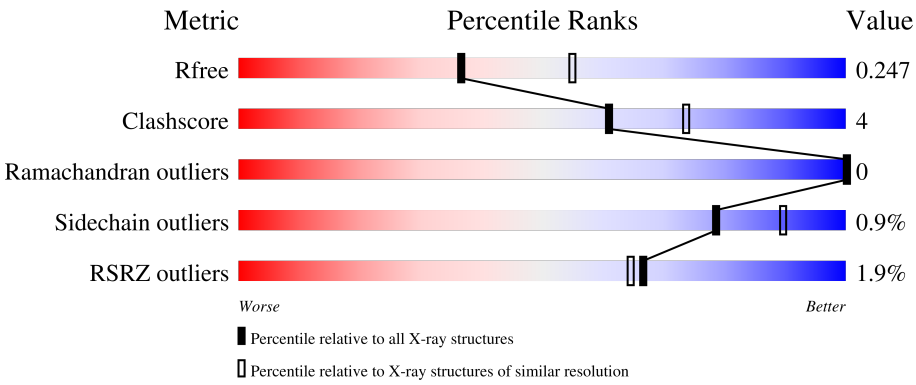
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	166	2%, 81%, 5%, 12%
1	B	166	2%, 83%, 13%
1	C	166	2%, 78%, 8%, 12%
1	D	166	78%, 9%, 13%
1	E	166	%, 84%, 11%

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Mol	Chain	Length	Quality of chain
1	F	166	2% 81% • 16%

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	EDO	A	301	-	-	X	-

2 Entry composition (i)

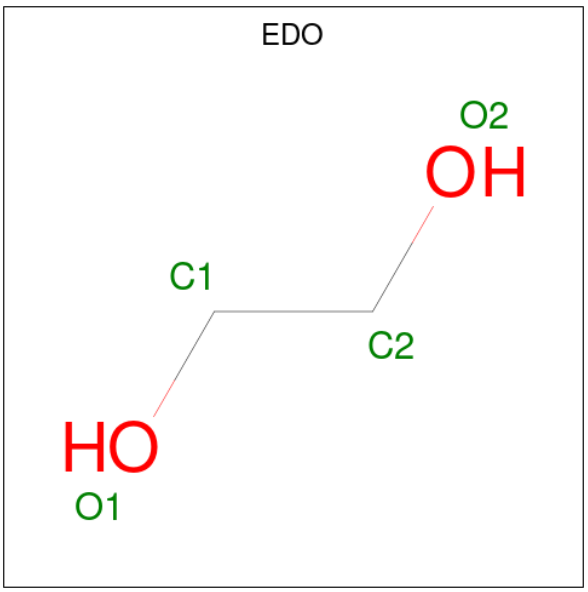
There are 3 unique types of molecules in this entry. The entry contains 7153 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	146	1152	738	195	215	4	0	1	0
			Total	C	N	O	S			
1	B	144	1098	704	191	200	3	0	0	0
			Total	C	N	O	S			
1	C	146	1131	726	196	205	4	0	1	0
			Total	C	N	O	S			
1	D	144	1125	723	193	205	4	0	1	0
			Total	C	N	O	S			
1	E	148	1153	739	197	213	4	0	0	0
			Total	C	N	O	S			
1	F	140	1063	688	175	197	3	0	0	0

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 3 is water.

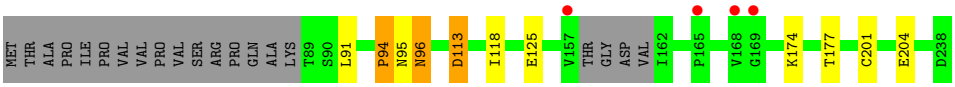
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	130	130	130	0	0
			Total	O		
3	B	36	36	36	0	0
			Total	O		
3	C	53	53	53	0	0
			Total	O		
3	D	86	86	86	0	0
			Total	O		
3	E	72	72	72	0	0
			Total	O		
3	F	50	50	50	0	0

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.

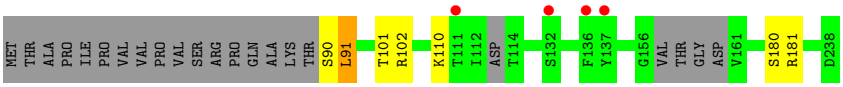
• Molecule 1: Coat protein

Chain A:



• Molecule 1: Coat protein

Chain B:



• Molecule 1: Coat protein

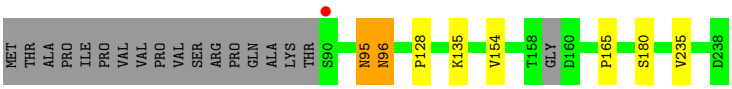
Chain C:

• Molecule 1: Coat protein

Chain D:

• Molecule 1: Coat protein

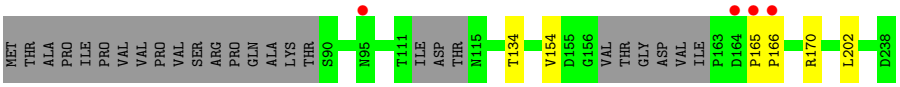
Chain E:



• Molecule 1: Coat protein

Chain F:

2%
81%
16%


MET
THR
ALA
PRO
ILE
PRO
VAL
VAL
PRO
VAL
SER
ARG
PRO
GLN
ALA
LYS
THR
S90
N95
T111
ILE
ASP
THR
N115
T134
V154
D155
G156
VAL
THR
GLY
ASP
VAL
ILE
P163
D164
P165
P166
R170
L202
D238

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	103.04Å 69.98Å 103.44Å 90.00° 105.51° 90.00°	Depositor
Resolution (Å)	99.67 – 2.40 40.34 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.5 (99.67-2.40) 97.5 (40.34-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$< I/\sigma(I) >$ [1]	2.78 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.207 , 0.244 0.211 , 0.247	Depositor DCC
R_{free} test set	2765 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	42.6	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 42.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.015 for l,-k,h	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7153	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% 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: EDO, CME

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	0.97	2/1167 (0.2%)	0.98	3/1586 (0.2%)
1	B	0.68	0/1116	0.82	1/1521 (0.1%)
1	C	0.71	0/1146	0.84	2/1562 (0.1%)
1	D	0.83	0/1139	0.88	0/1547
1	E	0.77	0/1168	0.84	0/1590
1	F	0.76	0/1081	0.87	2/1471 (0.1%)
All	All	0.79	2/6817 (0.0%)	0.88	8/9277 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	ASN	C-O	5.55	1.33	1.23
1	A	204	GLU	CD-OE1	5.37	1.31	1.25

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	102	ARG	NE-CZ-NH1	8.08	124.34	120.30
1	A	118	ILE	CA-CB-CG1	-7.22	97.27	111.00
1	C	102	ARG	NE-CZ-NH2	-6.84	116.88	120.30
1	F	202	LEU	CB-CG-CD1	6.28	121.67	111.00
1	B	91	LEU	N-CA-CB	5.50	121.39	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	113	ASP	CB-CG-OD1	-5.21	113.61	118.30
1	F	202	LEU	CB-CG-CD2	-5.09	102.34	111.00
1	A	113	ASP	CB-CG-OD2	5.08	122.87	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	94	PRO	Peptide

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	1152	0	1119	14	0
1	B	1098	0	1035	5	0
1	C	1131	0	1087	17	0
1	D	1125	0	1080	13	0
1	E	1153	0	1116	8	0
1	F	1063	0	994	4	0
2	A	4	0	6	8	0
3	A	130	0	0	5	0
3	B	36	0	0	1	0
3	C	53	0	0	1	0
3	D	86	0	0	6	0
3	E	72	0	0	0	0
3	F	50	0	0	1	0
All	All	7153	0	6437	59	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 (59) 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:213:GLN:HG2	3:D:371:HOH:O	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ASN:O	1:C:96:ASN:ND2	2.26	0.67
1:B:181:ARG:HD2	3:C:319:HOH:O	1.94	0.67
1:C:110:LYS:O	1:C:207[A]:ARG:HG2	1.95	0.66
1:D:134:THR:O	3:D:384:HOH:O	2.13	0.65
1:E:95:ASN:O	1:E:96:ASN:ND2	2.28	0.65
1:D:213:GLN:NE2	3:D:301:HOH:O	2.29	0.65
1:C:111:THR:O	1:C:207[A]:ARG:NH1	2.30	0.64
1:C:100:VAL:HG11	1:C:127:ILE:CG2	2.28	0.64
1:B:110:LYS:HB2	3:B:333:HOH:O	1.96	0.63
1:C:102:ARG:HG2	1:C:102:ARG:HH11	1.63	0.63
1:C:95:ASN:O	1:C:96:ASN:CB	2.47	0.62
1:A:94:PRO:CB	1:A:95:ASN:HB3	2.30	0.62
1:A:94:PRO:HB3	1:A:95:ASN:HB3	1.82	0.61
1:E:95:ASN:O	1:E:96:ASN:CB	2.47	0.61
1:D:205:ARG:HD3	3:D:379:HOH:O	2.02	0.60
2:A:301:EDO:H22	3:A:498:HOH:O	2.02	0.59
1:C:106:GLU:OE1	1:E:135:LYS:HD3	2.03	0.57
1:D:232[A]:ASP:HB2	3:D:306:HOH:O	2.05	0.55
1:D:154:VAL:HG12	1:D:165:PRO:HA	1.88	0.55
1:E:154:VAL:HG12	1:E:165:PRO:HA	1.88	0.55
1:E:95:ASN:O	1:E:96:ASN:HB3	2.07	0.55
1:B:101:THR:HG22	1:B:102:ARG:N	2.23	0.53
1:C:100:VAL:HG12	1:C:101:THR:N	2.23	0.53
1:C:100:VAL:HG11	1:C:127:ILE:HG23	1.90	0.53
1:C:102:ARG:NH1	1:C:103:LYS:O	2.41	0.53
1:F:154:VAL:HG12	1:F:165:PRO:HA	1.91	0.52
1:C:95:ASN:O	1:C:96:ASN:HB3	2.08	0.52
1:C:100:VAL:HG11	1:C:127:ILE:HG21	1.93	0.51
1:C:154:VAL:HG12	1:C:165:PRO:HA	1.92	0.51
1:D:101:THR:HG22	1:D:102:ARG:N	2.26	0.50
1:F:166:PRO:O	1:F:170:ARG:HG2	2.10	0.50
1:F:134:THR:O	3:F:331:HOH:O	2.20	0.50
1:E:154:VAL:CG1	1:E:165:PRO:HA	2.41	0.50
1:A:177:THR:CG2	2:A:301:EDO:H12	2.41	0.49
1:D:115:ASN:HA	3:D:324:HOH:O	2.12	0.48
1:A:113:ASP:OD1	1:A:113:ASP:C	2.51	0.48
1:A:125:GLU:CG	3:A:501:HOH:O	2.61	0.48
1:F:154:VAL:CG1	1:F:165:PRO:HA	2.42	0.48
1:A:91:LEU:HA	3:A:409:HOH:O	2.13	0.48
1:D:154:VAL:CG1	1:D:165:PRO:HA	2.43	0.47
1:C:154:VAL:CG1	1:C:165:PRO:HA	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:301:EDO:C2	3:A:498:HOH:O	2.62	0.47
1:A:95:ASN:O	1:A:96:ASN:HB2	2.16	0.46
1:B:90:SER:OG	1:B:91:LEU:N	2.48	0.46
1:A:174:LYS:HB2	2:A:301:EDO:H11	1.98	0.45
1:A:94:PRO:HB3	1:A:95:ASN:CB	2.48	0.43
1:A:177:THR:HG22	2:A:301:EDO:H12	2.00	0.43
1:A:177:THR:HB	2:A:301:EDO:H11	2.01	0.43
1:A:125:GLU:HG3	3:A:501:HOH:O	2.19	0.43
1:D:155:ASP:OD1	1:D:196:ARG:HG2	2.20	0.42
1:D:202:LEU:N	1:D:202:LEU:HD12	2.35	0.42
1:A:177:THR:HB	2:A:301:EDO:C1	2.48	0.42
1:B:180:SER:O	1:C:237:GLN:HA	2.20	0.41
1:D:237:GLN:HA	1:E:180:SER:O	2.21	0.41
1:D:178:TYR:HD2	1:E:235:VAL:HG23	1.85	0.41
1:A:174:LYS:H	2:A:301:EDO:C1	2.33	0.41
1:C:202:LEU:HD12	1:C:202:LEU:N	2.36	0.40
1:C:102:ARG:HD3	1:C:123:ILE:O	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	142/166 (86%)	140 (99%)	2 (1%)	0	100	100
1	B	137/166 (82%)	131 (96%)	6 (4%)	0	100	100
1	C	142/166 (86%)	140 (99%)	2 (1%)	0	100	100
1	D	138/166 (83%)	134 (97%)	4 (3%)	0	100	100
1	E	143/166 (86%)	140 (98%)	3 (2%)	0	100	100
1	F	133/166 (80%)	130 (98%)	3 (2%)	0	100	100
All	All	835/996 (84%)	815 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	119/143 (83%)	119 (100%)	0	100	100
1	B	108/143 (76%)	108 (100%)	0	100	100
1	C	112/143 (78%)	108 (96%)	4 (4%)	35	54
1	D	111/143 (78%)	111 (100%)	0	100	100
1	E	118/143 (82%)	115 (98%)	3 (2%)	47	67
1	F	102/143 (71%)	102 (100%)	0	100	100
All	All	670/858 (78%)	663 (99%)	7 (1%)	78	88

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	96	ASN
1	C	102	ARG
1	C	207[A]	ARG
1	C	207[B]	ARG
1	E	95	ASN
1	E	96	ASN
1	E	128	PRO

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

Mol	Chain	Res	Type
1	D	192	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

6 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	CME	B	201	1	4,5,10	1.15	0	1,5,11	0.17	0
1	CME	D	201	1	8,9,10	0.67	0	5,9,11	1.48	1 (20%)
1	CME	A	201	1	8,9,10	0.76	0	5,9,11	1.39	1 (20%)
1	CME	E	201	1	8,9,10	0.46	0	5,9,11	0.96	0
1	CME	C	201	1	8,9,10	0.72	0	5,9,11	1.04	0
1	CME	F	201	1	4,5,10	0.97	0	1,5,11	0.16	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	CME	B	201	1	-	0/1/4/10	-
1	CME	D	201	1	-	1/5/8/10	-
1	CME	A	201	1	-	1/5/8/10	-
1	CME	E	201	1	-	2/5/8/10	-
1	CME	C	201	1	-	2/5/8/10	-
1	CME	F	201	1	-	0/1/4/10	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	201	CME	CB-SG-SD	2.94	111.43	103.82
1	A	201	CME	CB-SG-SD	2.38	109.98	103.82

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	201	CME	SD-CE-CZ-OH
1	D	201	CME	SD-CE-CZ-OH
1	A	201	CME	SD-CE-CZ-OH
1	E	201	CME	SD-CE-CZ-OH
1	C	201	CME	CE-SD-SG-CB
1	E	201	CME	CE-SD-SG-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	EDO	A	301	-	3,3,3	0.47	0	2,2,2	0.89	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	301	-	-	0/1/1/1	-

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.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	EDO	8	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	145/166 (87%)	-0.09	4 (2%)	53	51	17, 30, 74, 113	0
1	B	143/166 (86%)	0.05	4 (2%)	53	51	36, 58, 87, 97	0
1	C	145/166 (87%)	-0.05	3 (2%)	63	61	30, 54, 86, 110	2 (1%)
1	D	143/166 (86%)	-0.23	0	100	100	23, 44, 71, 91	1 (0%)
1	E	147/166 (88%)	-0.35	1 (0%)	87	86	28, 42, 73, 86	0
1	F	139/166 (83%)	-0.10	4 (2%)	51	50	21, 50, 91, 141	0
All	All	862/996 (86%)	-0.13	16 (1%)	66	64	17, 47, 83, 141	3 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	111	THR	4.3
1	C	113	ASP	3.7
1	A	157	VAL	3.5
1	F	164	ASP	3.4
1	F	166	PRO	3.4
1	A	169	GLY	3.3
1	B	132	SER	3.1
1	F	165	PRO	2.9
1	F	95	ASN	2.9
1	A	168	VAL	2.7
1	A	165	PRO	2.6
1	B	111	THR	2.4
1	B	136	PHE	2.3
1	E	90	SER	2.3
1	B	137	TYR	2.3
1	C	114	THR	2.2

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
1	CME	A	201	10/11	0.92	0.20	20,32,82,87	0
1	CME	F	201	6/11	0.92	0.17	40,56,69,87	0
1	CME	B	201	6/11	0.93	0.14	47,51,56,57	0
1	CME	E	201	10/11	0.94	0.15	30,42,79,82	0
1	CME	C	201	10/11	0.95	0.17	35,48,92,97	0
1	CME	D	201	10/11	0.95	0.18	29,34,87,95	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
2	EDO	A	301	4/4	0.95	0.23	24,25,29,30	0

6.5 Other polymers (i)

There are no such residues in this entry.