



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 07:00 AM UTC

PDB ID : 5TR7 / pdb_00005tr7
Title : Crystal structure of a putative D-alanyl-D-alanine carboxypeptidase from *Vibrio cholerae* O1 biovar eltor str. N16961
Authors : Filippova, E.V.; Minasov, G.; Shuvalova, L.; Kiryukhina, O.; Dubrovskaya, I.; Shatsman, S.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2016-10-25
Resolution : 2.05 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

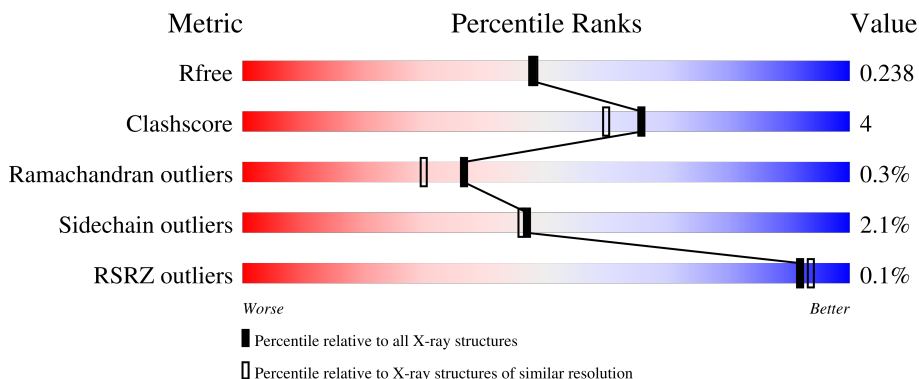
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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}	180053	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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 $\leq 5\%$. 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	341	
1	B	341	
1	C	341	

2 Entry composition [i](#)

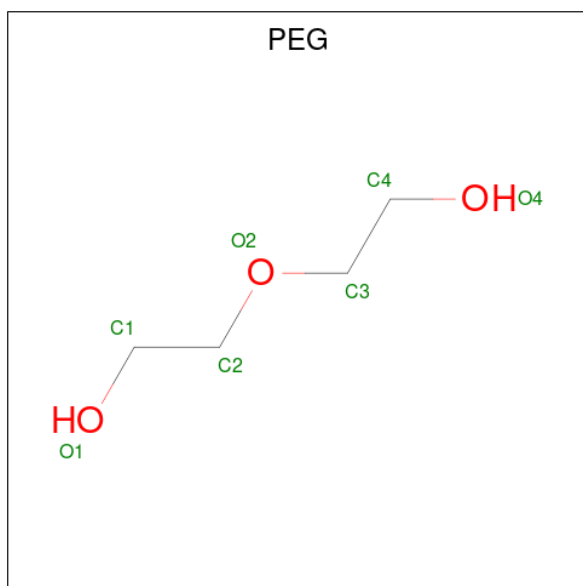
There are 5 unique types of molecules in this entry. The entry contains 5902 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-alanyl-D-alanine carboxypeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	0	0
			1930	1210	326	382	12			
1	B	249	Total	C	N	O	S	0	0	0
			1912	1200	324	376	12			
1	C	251	Total	C	N	O	S	0	1	0
			1936	1213	327	384	12			

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



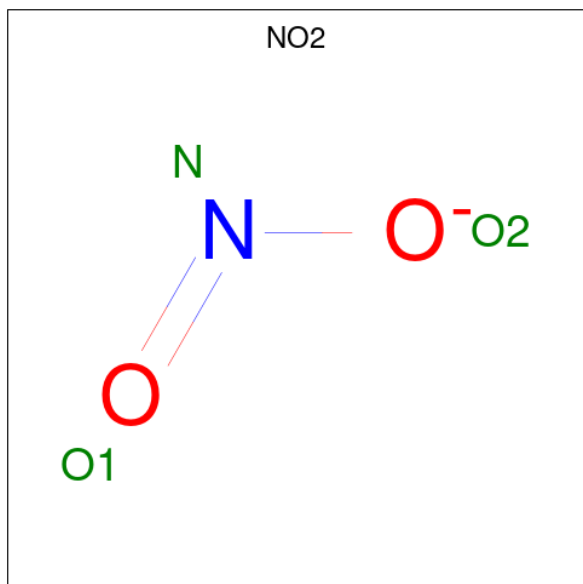
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		

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

- Molecule 3 is NITRITE ION (CCD ID: NO2) (formula: NO₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			3	1	2		
3	A	1	Total	N	O	0	0
			3	1	2		
3	C	1	Total	N	O	0	0
			3	1	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



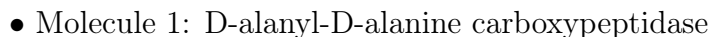
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	26	Total	O	0	0
			26	26		
5	B	38	Total	O	0	0
			38	38		
5	C	17	Total	O	0	0
			17	17		

i

- Molecule 1: D-alanyl-D-alanine carboxypeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	88.75Å 88.75Å 85.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.05 30.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-2.05) 99.9 (30.00-2.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 2.04Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.191 , 0.238 0.193 , 0.238	Depositor DCC
R_{free} test set	2246 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	48.6	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l 0.043 for h,-h-k,-l 0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5902	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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, 0.333 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: PEG, NO2, GOL

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	1.23	1/1965 (0.1%)	1.27	17/2656 (0.6%)
1	B	1.16	2/1947 (0.1%)	1.20	8/2631 (0.3%)
1	C	1.20	2/1971 (0.1%)	1.20	7/2664 (0.3%)
All	All	1.20	5/5883 (0.1%)	1.22	32/7951 (0.4%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	245	SER	C-O	-6.52	1.15	1.23
1	C	193	ARG	CD-NE	-5.64	1.38	1.46
1	B	195	VAL	CA-C	5.47	1.58	1.52
1	A	267	ARG	CA-C	5.20	1.59	1.52
1	B	126	ILE	C-O	-5.11	1.18	1.24

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ASP	N-CA-C	11.65	126.67	110.35
1	A	173	ASP	CB-CA-C	-11.36	91.67	109.61
1	A	237	SER	N-CA-C	8.49	121.61	111.33
1	C	237	SER	N-CA-C	8.41	121.51	111.33
1	B	173	ASP	N-CA-C	7.83	120.42	110.24
1	A	198	GLU	CG-CD-OE2	-7.17	101.92	118.40
1	B	60	THR	CB-CA-C	6.57	120.58	109.75
1	B	268	LYS	CB-CG-CD	5.83	124.70	111.30
1	A	63	SER	CB-CA-C	5.77	116.42	109.85
1	C	82	ARG	CB-CG-CD	5.68	124.36	111.30
1	A	111	GLU	N-CA-C	5.64	118.67	109.59
1	B	262	ASP	N-CA-C	5.57	118.28	111.82
1	C	193	ARG	NE-CZ-NH1	5.57	127.07	121.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	63	SER	N-CA-C	-5.56	100.59	109.04
1	A	262	ASP	N-CA-C	5.54	119.30	112.54
1	A	193	ARG	NE-CZ-NH1	5.50	127.00	121.50
1	A	98	TRP	CA-CB-CG	5.50	124.06	113.60
1	B	81	LYS	CB-CA-C	-5.50	101.67	110.79
1	B	63	SER	CB-CA-C	5.42	116.03	109.85
1	C	193	ARG	NE-CZ-NH2	-5.42	114.33	119.20
1	A	81	LYS	CB-CA-C	-5.34	101.92	110.79
1	A	59	ASP	CA-C-N	-5.33	115.37	122.72
1	A	59	ASP	C-N-CA	-5.33	115.37	122.72
1	A	233	THR	N-CA-CB	-5.29	102.71	110.85
1	A	217	ARG	NE-CZ-NH2	5.26	123.93	119.20
1	A	110	VAL	N-CA-CB	-5.25	103.31	110.56
1	C	81	LYS	CB-CA-C	-5.25	102.08	110.79
1	B	233	THR	N-CA-CB	-5.17	102.89	110.85
1	C	63	SER	N-CA-C	-5.15	102.08	109.24
1	A	63	SER	N-CA-C	-5.14	101.23	109.04
1	C	104	ASP	N-CA-C	5.10	118.48	111.54
1	A	110	VAL	CB-CA-C	5.04	118.12	111.21

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1930	0	1885	11	0
1	B	1912	0	1871	21	0
1	C	1936	0	1889	15	0
2	A	7	0	10	0	0
2	B	14	0	20	0	0
2	C	7	0	10	0	0
3	A	6	0	0	0	0
3	C	3	0	0	0	0
4	B	6	0	8	0	0
5	A	26	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	38	0	0	1	0
5	C	17	0	0	0	0
All	All	5902	0	5693	44	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 (44) 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:A:193:ARG:HD2	1:A:194:ASP:OD1	1.73	0.89
1:B:252:MET:HE3	1:B:282:PHE:CE1	2.08	0.89
1:C:193:ARG:HD2	1:C:194:ASP:OD1	1.79	0.82
1:B:252:MET:HE3	1:B:282:PHE:CD1	2.17	0.79
1:B:114:THR:HG21	5:B:512:HOH:O	1.84	0.78
1:B:167:THR:HB	1:C:224:LYS:HE2	1.67	0.76
1:C:217:ARG:HB2	1:C:233:THR:HG21	1.71	0.72
1:B:63:SER:O	1:B:236:THR:HG21	1.90	0.72
1:A:217:ARG:HB2	1:A:233:THR:HG21	1.74	0.69
1:B:236:THR:HG23	1:B:242:ASN:HD21	1.60	0.67
1:C:236:THR:HG23	1:C:242:ASN:HD21	1.61	0.66
1:B:217:ARG:HB2	1:B:233:THR:HG21	1.78	0.65
1:C:98:TRP:CG	1:C:112:VAL:HG23	2.33	0.63
1:B:62:LEU:HD13	1:B:259:MET:HE2	1.80	0.62
1:B:90:ASP:OD1	1:B:117:LYS:CE	2.47	0.62
1:C:58:MET:HG2	1:C:182:TYR:HB2	1.87	0.55
1:A:58:MET:HG2	1:A:182:TYR:HB2	1.89	0.53
1:C:78:GLN:O	1:C:82:ARG:HG3	2.09	0.53
1:B:58:MET:HG2	1:B:182:TYR:HB2	1.91	0.53
1:B:111:GLU:O	1:B:114:THR:HG22	2.10	0.52
1:C:236:THR:HG23	1:C:242:ASN:ND2	2.26	0.51
1:B:78:GLN:O	1:B:82:ARG:HG3	2.10	0.50
1:B:63:SER:O	1:B:236:THR:CG2	2.60	0.49
1:B:236:THR:HG23	1:B:242:ASN:ND2	2.27	0.48
1:A:227:ASN:HB3	5:A:524:HOH:O	2.14	0.48
1:C:98:TRP:CD2	1:C:112:VAL:HG23	2.49	0.47
1:C:236:THR:OG1	1:C:239:ALA:HB3	2.15	0.47
1:A:263:ASN:HB2	5:A:508:HOH:O	2.16	0.46
1:C:58:MET:CG	1:C:182:TYR:HB2	2.46	0.46
1:A:93:ILE:HG21	1:A:110:VAL:HG13	1.98	0.45
1:B:236:THR:OG1	1:B:239:ALA:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:SER:O	1:C:236:THR:HG21	2.17	0.45
1:B:270:GLU:HA	1:B:270:GLU:OE1	2.18	0.44
1:B:90:ASP:OD1	1:B:117:LYS:HE2	2.15	0.44
1:B:168:ASN:HD22	1:C:224:LYS:NZ	2.16	0.44
1:A:58:MET:CG	1:A:182:TYR:HB2	2.48	0.42
1:C:166:PHE:HA	1:C:177:LEU:HD13	2.02	0.42
1:B:167:THR:CB	1:C:224:LYS:HE2	2.45	0.42
1:B:58:MET:CG	1:B:182:TYR:HB2	2.49	0.42
1:A:219:GLY:HA3	1:A:268:LYS:HD3	2.02	0.42
1:A:171:GLY:HA2	1:A:177:LEU:HD21	2.03	0.41
1:A:63:SER:HB3	1:A:236:THR:HG23	2.03	0.40
1:B:166:PHE:HA	1:B:177:LEU:HD13	2.03	0.40
1:A:166:PHE:HA	1:A:177:LEU:HD13	2.03	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	249/341 (73%)	244 (98%)	4 (2%)	1 (0%)	30	22
1	B	247/341 (72%)	241 (98%)	5 (2%)	1 (0%)	30	22
1	C	250/341 (73%)	244 (98%)	6 (2%)	0	100	100
All	All	746/1023 (73%)	729 (98%)	15 (2%)	2 (0%)	36	30

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	PHE
1	B	109	PHE

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	209/283 (74%)	203 (97%)	6 (3%)	37	33
1	B	207/283 (73%)	203 (98%)	4 (2%)	50	50
1	C	210/283 (74%)	207 (99%)	3 (1%)	59	60
All	All	626/849 (74%)	613 (98%)	13 (2%)	47	46

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	TRP
1	A	104	ASP
1	A	110	VAL
1	A	122	ASN
1	A	198	GLU
1	A	217	ARG
1	B	60	THR
1	B	114	THR
1	B	173	ASP
1	B	268	LYS
1	C	95	LYS
1	C	110	VAL
1	C	173	ASP

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

Mol	Chain	Res	Type
1	B	140	HIS

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	PEG	A	401	-	6,6,6	0.56	0	5,5,5	0.31	0
2	PEG	B	401	-	6,6,6	0.49	0	5,5,5	0.40	0
2	PEG	B	402	-	6,6,6	0.47	0	5,5,5	0.32	0
3	NO2	C	402	-	1,2,2	0.00	0	0,1,1	-	-
2	PEG	C	401	-	6,6,6	0.47	0	5,5,5	0.28	0
3	NO2	A	403	-	1,2,2	0.17	0	0,1,1	-	-
4	GOL	B	403	-	5,5,5	0.32	0	5,5,5	0.52	0
3	NO2	A	402	-	1,2,2	0.04	0	0,1,1	-	-

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	PEG	A	401	-	-	3/4/4/4	-
2	PEG	B	401	-	-	1/4/4/4	-
2	PEG	B	402	-	-	1/4/4/4	-
2	PEG	C	401	-	-	2/4/4/4	-
4	GOL	B	403	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	401	PEG	O1-C1-C2-O2
2	B	402	PEG	O2-C3-C4-O4
2	A	401	PEG	O2-C3-C4-O4
2	A	401	PEG	C1-C2-O2-C3
2	B	401	PEG	C1-C2-O2-C3
2	C	401	PEG	O2-C3-C4-O4
2	A	401	PEG	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	251/341 (73%)	-0.17	1 (0%) 88 90	46, 61, 87, 125	0
1	B	249/341 (73%)	-0.19	0 100 100	42, 61, 93, 125	0
1	C	251/341 (73%)	-0.09	0 100 100	27, 64, 93, 109	1 (0%)
All	All	751/1023 (73%)	-0.15	1 (0%) 92 93	27, 62, 92, 125	1 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	62	LEU	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 oligosaccharides 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	NO2	C	402	3/3	0.76	0.12	87,87,89,92	0
3	NO2	A	402	3/3	0.79	0.10	80,80,87,100	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PEG	B	402	7/7	0.80	0.09	89,95,100,100	0
4	GOL	B	403	6/6	0.88	0.12	73,82,86,87	0
3	NO2	A	403	3/3	0.90	0.09	68,68,72,72	0
2	PEG	C	401	7/7	0.90	0.09	67,70,72,72	0
2	PEG	A	401	7/7	0.90	0.10	78,90,103,104	0
2	PEG	B	401	7/7	0.95	0.07	52,64,69,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.