



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:18 AM UTC

PDB ID : 7E6A / pdb_00007e6a
Title : Crystal structure of cysteine desulfurase SufS C361A from *Bacillus subtilis*
Authors : Nakamura, R.; Takahashi, Y.; Fujishiro, T.
Deposited on : 2021-02-22
Resolution : 1.96 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

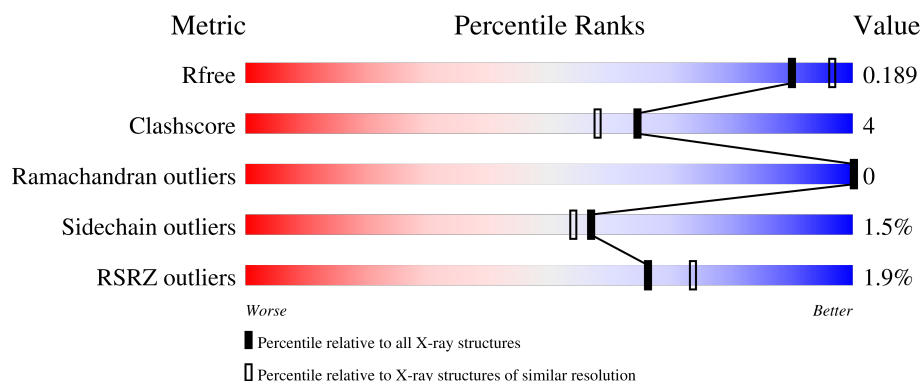
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

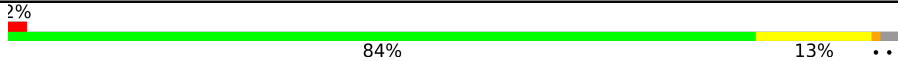
The reported resolution of this entry is 1.96 Å.

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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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 $\leq 5\%$. 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	419	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3553 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 Cysteine desulfurase SufS.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	412	3269	2073	557	626	1	12	0	6	0

There are 14 discrepancies between the modelled and reference sequences:

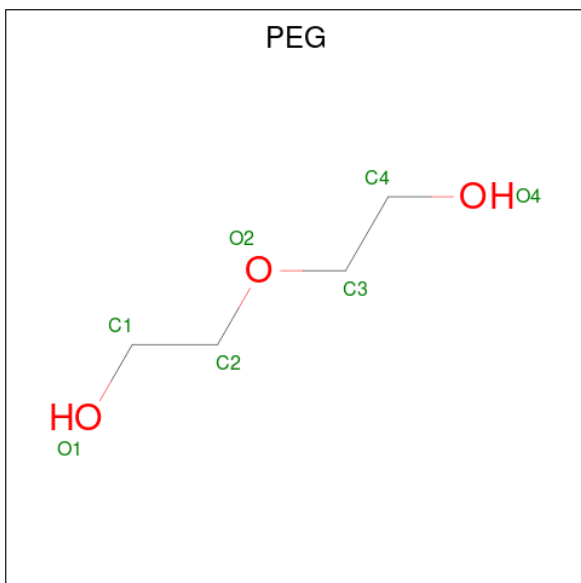
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP O32164
A	-1	GLY	-	expression tag	UNP O32164
A	0	HIS	-	expression tag	UNP O32164
A	361	ALA	CYS	engineered mutation	UNP O32164
A	407	VAL	-	expression tag	UNP O32164
A	408	ASP	-	expression tag	UNP O32164
A	409	LEU	-	expression tag	UNP O32164
A	410	GLU	-	expression tag	UNP O32164
A	411	HIS	-	expression tag	UNP O32164
A	412	HIS	-	expression tag	UNP O32164
A	413	HIS	-	expression tag	UNP O32164
A	414	HIS	-	expression tag	UNP O32164
A	415	HIS	-	expression tag	UNP O32164
A	416	HIS	-	expression tag	UNP O32164

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



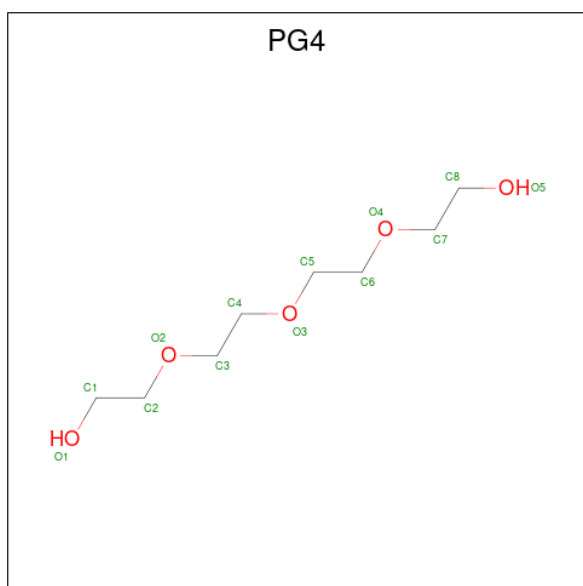
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 13 8 5	0	0

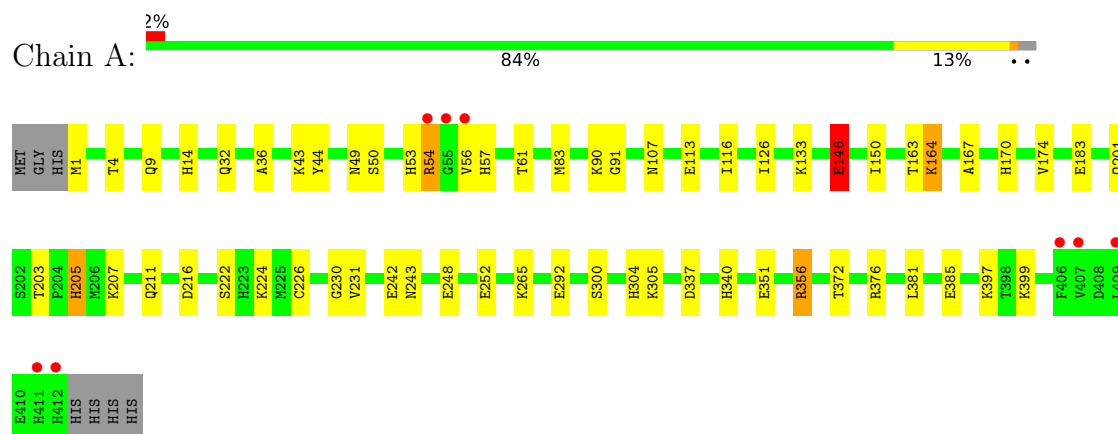
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	206	Total O 206 206	0	0

3 Residue-property plots

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: Cysteine desulfurase SufS



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	92.60Å 92.60Å 129.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.34 – 1.96 46.34 – 1.96	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.34-1.96) 99.8 (46.34-1.96)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.155 , 0.179 (Not available) , 0.189	Depositor DCC
R_{free} test set	2323 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3553	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, PEG, LLP, EDO

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	1.24	13/3330 (0.4%)	1.47	31/4519 (0.7%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	53	HIS	CE1-NE2	10.76	1.43	1.32
1	A	340	HIS	CE1-NE2	7.38	1.40	1.32
1	A	205	HIS	CA-CB	7.05	1.63	1.54
1	A	57	HIS	CE1-NE2	6.99	1.39	1.32
1	A	183	GLU	CD-OE1	6.57	1.37	1.25
1	A	170	HIS	CE1-NE2	5.90	1.38	1.32
1	A	340	HIS	CG-ND1	5.90	1.44	1.38
1	A	50	SER	CA-CB	5.70	1.63	1.53
1	A	304	HIS	CG-ND1	5.53	1.44	1.38
1	A	53	HIS	CG-ND1	5.45	1.44	1.38
1	A	91	GLY	N-CA	5.34	1.50	1.45
1	A	356	ARG	NE-CZ	5.14	1.38	1.33
1	A	304	HIS	CE1-NE2	5.05	1.37	1.32

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ILE	N-CA-CB	8.87	116.08	110.50
1	A	356	ARG	CG-CD-NE	8.72	131.19	112.00
1	A	337	ASP	CA-C-N	-8.67	108.88	122.67
1	A	337	ASP	C-N-CA	-8.67	108.88	122.67
1	A	14	HIS	CA-CB-CG	-8.16	105.64	113.80
1	A	107	ASN	CA-C-N	-7.25	112.42	122.93
1	A	107	ASN	C-N-CA	-7.25	112.42	122.93
1	A	146	GLU	CB-CG-CD	6.51	123.67	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	LYS	N-CA-CB	-6.33	99.78	110.41
1	A	61	THR	CA-CB-OG1	-6.25	100.22	109.60
1	A	150	ILE	CA-C-N	-6.24	111.39	122.07
1	A	150	ILE	C-N-CA	-6.24	111.39	122.07
1	A	300	SER	CB-CA-C	-6.08	100.51	110.85
1	A	265	LYS	CB-CA-C	-6.00	97.98	110.40
1	A	203	THR	CA-CB-OG1	-5.90	100.75	109.60
1	A	399	LYS	CB-CA-C	5.80	119.99	110.88
1	A	56	VAL	CA-C-N	-5.78	113.51	122.59
1	A	56	VAL	C-N-CA	-5.78	113.51	122.59
1	A	226	CYS	CA-C-N	-5.69	112.94	121.87
1	A	226	CYS	C-N-CA	-5.69	112.94	121.87
1	A	216	ASP	CA-CB-CG	-5.65	106.95	112.60
1	A	356	ARG	CB-CG-CD	5.64	124.27	111.30
1	A	351	GLU	CB-CA-C	-5.58	101.13	110.56
1	A	248	GLU	CB-CA-C	-5.52	97.93	110.07
1	A	230	GLY	CA-C-N	-5.21	114.86	122.58
1	A	230	GLY	C-N-CA	-5.21	114.86	122.58
1	A	243	ASN	CA-CB-CG	-5.21	107.39	112.60
1	A	242	GLU	CB-CG-CD	5.18	121.41	112.60
1	A	203	THR	O-C-N	-5.15	115.60	120.70
1	A	372	THR	CA-CB-OG1	-5.11	101.93	109.60
1	A	174	VAL	CB-CA-C	-5.07	106.54	111.05

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	3269	0	3228	26	0
2	A	16	0	24	1	0
3	A	49	0	70	8	0
4	A	13	0	18	0	0
5	A	206	0	0	2	0
All	All	3553	0	3340	26	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 (26) 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:146:GLU:H	1:A:146:GLU:CD	1.81	0.86
1:A:164:LYS:HE2	5:A:643:HOH:O	1.91	0.71
1:A:43:LYS:NZ	3:A:509:PEG:H42	2.06	0.70
1:A:397:LYS:HB3	3:A:510:PEG:H41	1.74	0.68
1:A:83:MET:HE1	3:A:506:PEG:H32	1.84	0.59
1:A:54[B]:ARG:NH2	1:A:252:GLU:OE2	2.36	0.56
1:A:397:LYS:CB	3:A:510:PEG:H41	2.36	0.56
1:A:146:GLU:CD	1:A:146:GLU:N	2.58	0.56
1:A:43:LYS:HZ1	3:A:509:PEG:H42	1.70	0.55
1:A:381:LEU:H	1:A:381:LEU:HD23	1.73	0.53
1:A:43:LYS:HZ2	3:A:509:PEG:H42	1.72	0.52
1:A:113:GLU:HB2	1:A:163:THR:HA	1.92	0.51
1:A:211[B]:GLN:HA	1:A:211[B]:GLN:OE1	2.09	0.50
1:A:356:ARG:HH11	1:A:376:ARG:HH22	1.58	0.50
1:A:36:ALA:HB1	3:A:507:PEG:H41	1.95	0.49
1:A:44:TYR:CZ	1:A:49:ASN:HA	2.47	0.49
1:A:356:ARG:NH1	1:A:376:ARG:HH22	2.12	0.46
1:A:9[B]:GLN:HE22	2:A:504:EDO:C1	2.29	0.46
1:A:222:SER:CB	1:A:231:VAL:HG13	2.46	0.46
1:A:32:GLN:HA	1:A:32:GLN:OE1	2.17	0.44
1:A:207:LYS:NZ	1:A:292:GLU:OE2	2.51	0.43
1:A:116:ILE:HG22	1:A:167:ALA:HB3	2.00	0.42
1:A:305[A]:LYS:HE3	1:A:385:GLU:OE2	2.19	0.41
1:A:201:GLN:O	1:A:205:HIS:HD2	2.03	0.41
1:A:164:LYS:CE	5:A:643:HOH:O	2.62	0.40
1:A:4:THR:OG1	3:A:505:PEG:H12	2.21	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	415/419 (99%)	406 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	347/347 (100%)	341 (98%)	6 (2%)	53	49

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	54[A]	ARG
1	A	54[B]	ARG
1	A	133	LYS
1	A	146	GLU
1	A	164	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	HIS
1	A	18	ASN
1	A	20	HIS
1	A	335	ASN
1	A	359	HIS
1	A	411	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	$\# Z > 2$
1	LLP	A	224	1	23,24,25	1.03	2 (8%)	25,32,34	1.15	3 (12%)

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	LLP	A	224	1	-	1/16/17/19	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	224	LLP	P-OP2	-2.46	1.45	1.54
1	A	224	LLP	C4-C5	-2.41	1.38	1.42

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	LLP	C2'-C2-C3	2.39	123.59	120.80
1	A	224	LLP	C4-C4'-NZ	-2.01	114.74	124.04
1	A	224	LLP	C3-C2-N1	-2.01	118.43	120.96

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	224	LLP	C4-C4'-NZ-CE

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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	PEG	A	508	-	6,6,6	0.58	0	5,5,5	0.44	0
3	PEG	A	511	-	6,6,6	0.23	0	5,5,5	0.14	0
2	EDO	A	501	-	3,3,3	0.19	0	2,2,2	0.57	0
2	EDO	A	502	-	3,3,3	0.41	0	2,2,2	0.49	0
3	PEG	A	506	-	6,6,6	0.21	0	5,5,5	0.17	0
3	PEG	A	505	-	6,6,6	0.42	0	5,5,5	0.23	0
3	PEG	A	510	-	6,6,6	0.16	0	5,5,5	0.14	0
3	PEG	A	509	-	6,6,6	0.09	0	5,5,5	0.24	0
2	EDO	A	504	-	3,3,3	0.13	0	2,2,2	0.28	0
3	PEG	A	507	-	6,6,6	0.19	0	5,5,5	0.30	0
4	PG4	A	512	-	12,12,12	0.29	0	11,11,11	0.51	0
2	EDO	A	503	-	3,3,3	0.31	0	2,2,2	0.30	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	PEG	A	508	-	-	3/4/4/4	-
3	PEG	A	511	-	-	1/4/4/4	-
2	EDO	A	501	-	-	1/1/1/1	-
2	EDO	A	502	-	-	1/1/1/1	-
3	PEG	A	506	-	-	2/4/4/4	-
3	PEG	A	505	-	-	1/4/4/4	-
3	PEG	A	510	-	-	0/4/4/4	-
3	PEG	A	509	-	-	4/4/4/4	-
2	EDO	A	504	-	-	0/1/1/1	-
3	PEG	A	507	-	-	2/4/4/4	-
4	PG4	A	512	-	-	1/10/10/10	-
2	EDO	A	503	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	507	PEG	O2-C3-C4-O4
3	A	507	PEG	O1-C1-C2-O2
3	A	506	PEG	O2-C3-C4-O4
2	A	501	EDO	O1-C1-C2-O2
2	A	503	EDO	O1-C1-C2-O2
3	A	508	PEG	O1-C1-C2-O2
3	A	506	PEG	C1-C2-O2-C3
3	A	508	PEG	O2-C3-C4-O4
3	A	509	PEG	O1-C1-C2-O2
3	A	508	PEG	C1-C2-O2-C3
3	A	509	PEG	O2-C3-C4-O4
3	A	509	PEG	C1-C2-O2-C3
3	A	509	PEG	C4-C3-O2-C2
3	A	505	PEG	O1-C1-C2-O2
2	A	502	EDO	O1-C1-C2-O2
3	A	511	PEG	O2-C3-C4-O4
4	A	512	PG4	C1-C2-O2-C3

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	506	PEG	1	0
3	A	505	PEG	1	0
3	A	510	PEG	2	0
3	A	509	PEG	3	0
2	A	504	EDO	1	0
3	A	507	PEG	1	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	411/419 (98%)	-0.17	8 (1%) 66 74	21, 38, 63, 129	6 (1%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	409	LEU	5.8
1	A	407	VAL	5.2
1	A	406	PHE	3.6
1	A	55	GLY	3.0
1	A	54[A]	ARG	2.7
1	A	56	VAL	2.7
1	A	412	HIS	2.7
1	A	411	HIS	2.1

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	LLP	A	224	24/25	0.99	0.05	26,33,38,42	0

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($\text{\AA}^2$)	Q<0.9
3	PEG	A	508	7/7	0.82	0.22	45,56,74,74	0
3	PEG	A	511	7/7	0.85	0.23	85,92,99,108	0
3	PEG	A	510	7/7	0.87	0.16	75,80,86,87	0
2	EDO	A	504	4/4	0.88	0.15	64,67,68,70	0
3	PEG	A	509	7/7	0.89	0.19	70,76,91,92	0
3	PEG	A	506	7/7	0.92	0.13	60,66,75,85	0
3	PEG	A	507	7/7	0.92	0.13	59,63,76,77	0
2	EDO	A	502	4/4	0.92	0.13	54,60,65,68	0
4	PG4	A	512	13/13	0.92	0.13	48,58,83,86	0
2	EDO	A	501	4/4	0.93	0.13	57,58,60,63	0
2	EDO	A	503	4/4	0.93	0.13	56,62,65,66	0
3	PEG	A	505	7/7	0.94	0.10	45,52,63,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.