



# wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 16, 2017 – 10:39 AM EST

PDB ID : 6EQS  
Title : Human Sirt5 in complex with stalled peptidylimide intermediate of inhibitory compound 29  
Authors : Pannek, M.; Steegborn, C.  
Deposited on : unknown  
Resolution : 1.32 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

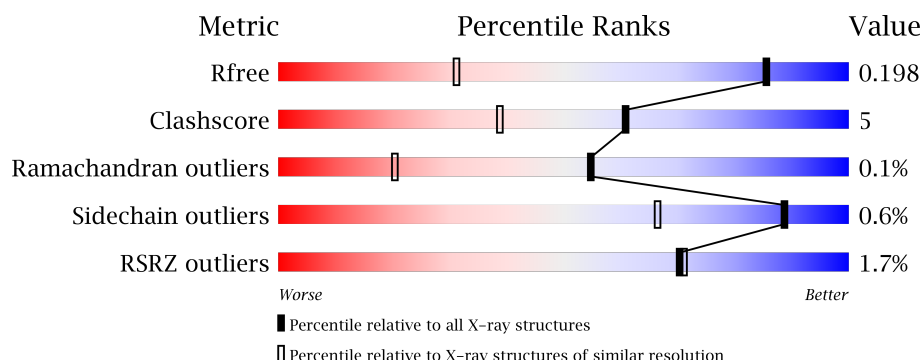
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.32 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))
$R_{free}$	100719	1164 (1.34-1.30)
Clashscore	112137	1223 (1.34-1.30)
Ramachandran outliers	110173	1179 (1.34-1.30)
Sidechain outliers	110143	1179 (1.34-1.30)
RSRZ outliers	101464	1167 (1.34-1.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	275	<div> <div>0.1%</div> <div> <div></div> <div>79%</div> <div>17%</div> <div>...</div> </div> </div>
1	B	275	<div> <div></div> <div> <div>82%</div> <div>13%</div> <div>...</div> </div> </div>
1	C	275	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>16%</div> <div>.</div> </div> </div>
1	D	275	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>15%</div> <div>.</div> </div> </div>

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.

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	EDO	A	403[A]	-	-	-	X
4	EDO	A	403[B]	-	-	-	X
4	EDO	A	404[A]	-	-	-	X
4	EDO	A	404[B]	-	-	-	X
4	EDO	A	405[A]	-	-	-	X
4	EDO	A	405[B]	-	-	-	X
4	EDO	A	406	-	-	-	X
4	EDO	B	403[A]	-	-	-	X
4	EDO	B	403[B]	-	-	-	X
4	EDO	B	404[A]	-	-	-	X
4	EDO	B	404[B]	-	-	-	X
4	EDO	B	405[A]	-	-	-	X
4	EDO	B	405[B]	-	-	-	X
4	EDO	B	406	-	-	X	-
4	EDO	C	403[A]	-	-	-	X
4	EDO	C	403[B]	-	-	-	X
4	EDO	C	404	-	-	X	-
4	EDO	C	405	-	-	-	X
4	EDO	D	403[A]	-	-	-	X
4	EDO	D	403[B]	-	-	-	X
4	EDO	D	404[A]	-	-	-	X
4	EDO	D	404[B]	-	-	-	X
5	BU2	A	408	-	-	X	X
5	BU2	A	409	-	-	-	X
5	BU2	B	408[A]	-	-	X	X
5	BU2	B	408[B]	-	-	-	X
5	BU2	B	409	-	-	-	X
5	BU2	B	410	-	-	X	-
5	BU2	C	406	-	-	-	X
5	BU2	D	405	-	-	X	X
5	BU2	D	406	-	-	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 10389 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 NAD-dependent protein deacetylase sirtuin-5, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	9	0
			2113	1334	387	377	15			
1	B	268	Total	C	N	O	S	0	10	0
			2107	1333	384	376	14			
1	C	274	Total	C	N	O	S	0	6	0
			2136	1351	388	384	13			
1	D	275	Total	C	N	O	S	0	7	0
			2145	1356	389	387	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLY	-	expression tag	UNP Q9NXA8
A	29	ILE	-	expression tag	UNP Q9NXA8
A	30	ASP	-	expression tag	UNP Q9NXA8
A	31	PRO	-	expression tag	UNP Q9NXA8
A	32	PHE	-	expression tag	UNP Q9NXA8
A	33	THR	-	expression tag	UNP Q9NXA8
B	28	GLY	-	expression tag	UNP Q9NXA8
B	29	ILE	-	expression tag	UNP Q9NXA8
B	30	ASP	-	expression tag	UNP Q9NXA8
B	31	PRO	-	expression tag	UNP Q9NXA8
B	32	PHE	-	expression tag	UNP Q9NXA8
B	33	THR	-	expression tag	UNP Q9NXA8
C	28	GLY	-	expression tag	UNP Q9NXA8
C	29	ILE	-	expression tag	UNP Q9NXA8
C	30	ASP	-	expression tag	UNP Q9NXA8
C	31	PRO	-	expression tag	UNP Q9NXA8
C	32	PHE	-	expression tag	UNP Q9NXA8
C	33	THR	-	expression tag	UNP Q9NXA8
D	28	GLY	-	expression tag	UNP Q9NXA8
D	29	ILE	-	expression tag	UNP Q9NXA8
D	30	ASP	-	expression tag	UNP Q9NXA8

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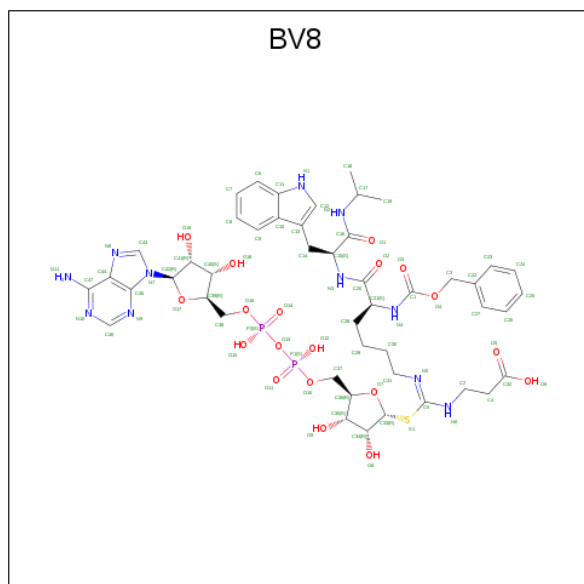
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Chain	Residue	Modelled	Actual	Comment	Reference
D	31	PRO	-	expression tag	UNP Q9NXA8
D	32	PHE	-	expression tag	UNP Q9NXA8
D	33	THR	-	expression tag	UNP Q9NXA8

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is 3-[[({Z})- {C}-(2 {R},3 {R},4 {S},5 {R})-5-[[[(2 {R},3 {S},4 {R},5 {R})-5-(6-aminopurin-9-yl)-3,4-bis(oxidanyl)oxolan-2-yl]methoxy-oxidanyl-phosphoryl]oxy-oxidanyl-phosphoryl]oxymethyl]-3,4-bis(oxidanyl)oxolan-2-yl]sulfanyl- {N}-(5 {S})-6-[(2 {S})-3-(1 {H}-indol-3-yl)-1-oxidanylidene-1-(propan-2-ylamino)propan-2-yl]amino]-6-oxidanylidene-5-(phenylmethoxycarbonylamino)hexyl]carbonimidoyl]amino]propanoic acid (three-letter code: BV8) (formula: C<sub>47</sub>H<sub>63</sub>N<sub>11</sub>O<sub>19</sub>P<sub>2</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P S 160 94 22 38 4 2	0	1

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total 160	C 94	N 22	O 38	P 4	S 2	0	1
3	C	1	Total 160	C 94	N 22	O 38	P 4	S 2	0	1
3	D	1	Total 160	C 94	N 22	O 38	P 4	S 2	0	1

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



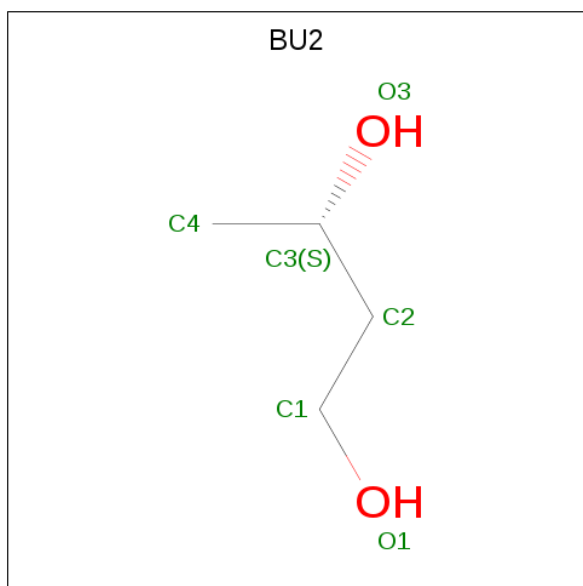
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 8	C 4	O 4	0	1
4	A	1	Total 8	C 4	O 4	0	1
4	A	1	Total 8	C 4	O 4	0	1
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	B	1	Total 8	C 4	O 4	0	1
4	B	1	Total 8	C 4	O 4	0	1
4	B	1	Total 8	C 4	O 4	0	1

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	1
			8	4	4		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	1
			8	4	4		
4	D	1	Total	C	O	0	1
			8	4	4		

- Molecule 5 is 1,3-BUTANEDIOL (three-letter code: BU2) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>2</sub>).



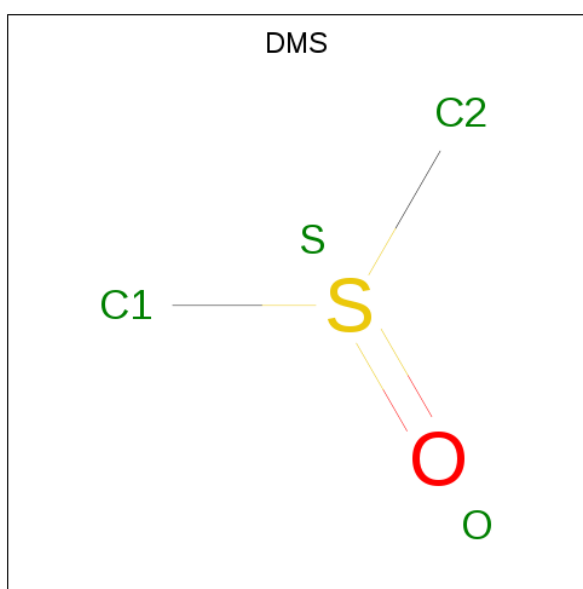
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	4	2		
5	A	1	Total	C	O	0	0
			6	4	2		
5	B	1	Total	C	O	0	1
			12	8	4		
5	B	1	Total	C	O	0	0
			6	4	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	4	2		
5	C	1	Total	C	O	0	0
			6	4	2		
5	D	1	Total	C	O	0	0
			6	4	2		
5	D	1	Total	C	O	0	0
			6	4	2		

- Molecule 6 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C<sub>2</sub>H<sub>6</sub>OS).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	294	Total	O	0	0
			294	294		
7	B	288	Total	O	0	0
			288	288		
7	C	255	Total	O	0	0
			255	255		

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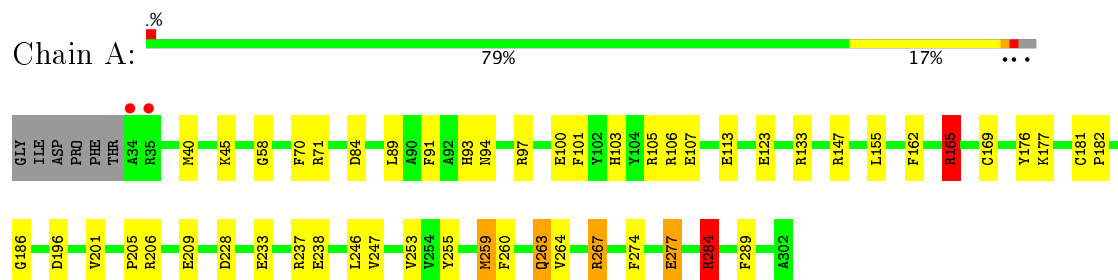
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	249	Total 249	O 249	0	0

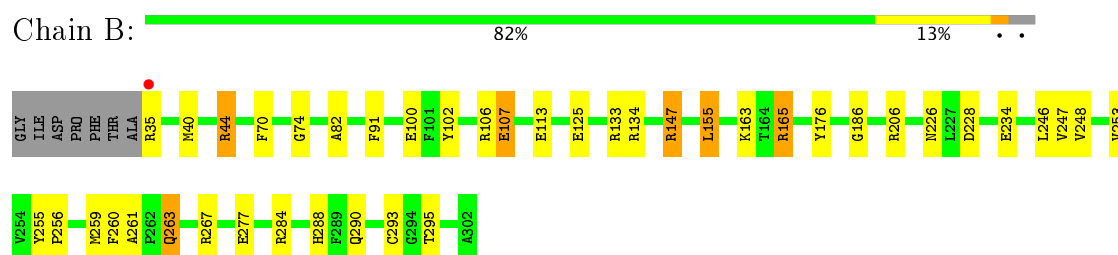
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

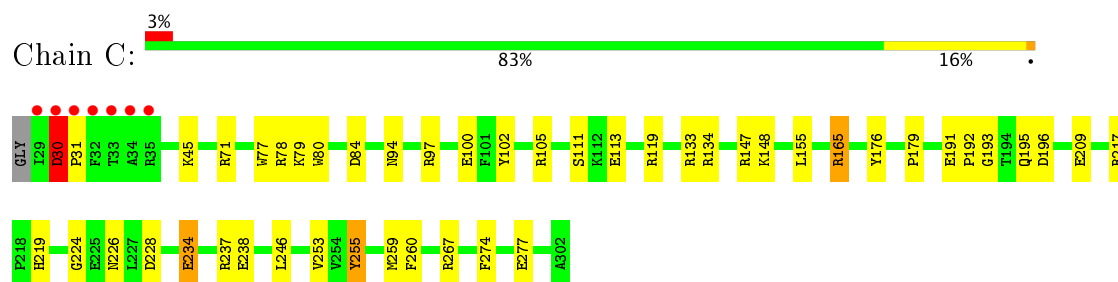
- Molecule 1: NAD-dependent protein deacylase sirtuin-5, mitochondrial



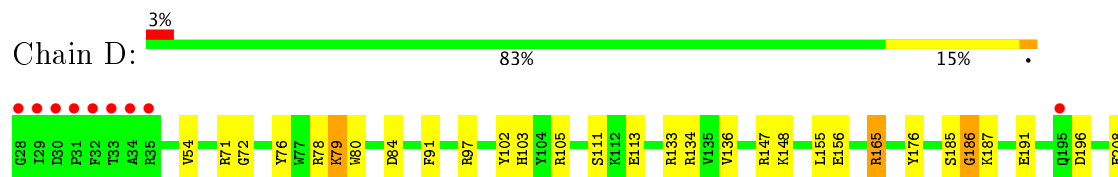
- Molecule 1: NAD-dependent protein deacylase sirtuin-5, mitochondrial



- Molecule 1: NAD-dependent protein deacylase sirtuin-5, mitochondrial



- Molecule 1: NAD-dependent protein deacylase sirtuin-5, mitochondrial





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	40.33Å 55.97Å 123.03Å 97.39° 99.29° 90.52°	Depositor
Resolution (Å)	43.77 – 1.32 43.77 – 1.32	Depositor EDS
% Data completeness (in resolution range)	91.9 (43.77-1.32) 88.1 (43.77-1.32)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.40 (at 1.32Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.160 , 0.209 0.159 , 0.198	Depositor DCC
$R_{free}$ test set	2388 reflections (1.06%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.2	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 61.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	10389	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 91.69 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3245e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: BU2, ZN, BV8, DMS, 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	1.58	17/2187 (0.8%)	1.59	37/2964 (1.2%)
1	B	1.52	19/2187 (0.9%)	1.52	27/2965 (0.9%)
1	C	1.60	21/2203 (1.0%)	1.44	29/2990 (1.0%)
1	D	1.69	23/2218 (1.0%)	1.43	25/3009 (0.8%)
All	All	1.60	80/8795 (0.9%)	1.50	118/11928 (1.0%)

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
1	B	0	1
1	C	0	2
1	D	0	2
All	All	0	6

The worst 5 of 80 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	234	GLU	CD-OE2	30.55	1.59	1.25
1	C	234	GLU	CD-OE2	18.00	1.45	1.25
1	D	234	GLU	CD-OE1	13.73	1.40	1.25
1	A	238	GLU	CD-OE1	-11.62	1.12	1.25
1	C	165	ARG	CZ-NH1	-9.30	1.21	1.33

The worst 5 of 118 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	284[A]	ARG	NE-CZ-NH2	17.53	129.06	120.30
1	A	284[B]	ARG	NE-CZ-NH2	17.53	129.06	120.30
1	B	267[A]	ARG	NE-CZ-NH1	15.83	128.22	120.30
1	B	267[B]	ARG	NE-CZ-NH1	15.83	128.22	120.30
1	D	267	ARG	NE-CZ-NH1	12.89	126.75	120.30

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	ARG	Sidechain
1	B	165	ARG	Sidechain
1	C	165	ARG	Sidechain
1	C	234	GLU	Sidechain
1	D	165	ARG	Sidechain

## 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	2113	0	2106	17	0
1	B	2107	0	2110	27	1
1	C	2136	0	2118	15	0
1	D	2145	0	2127	18	1
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	160	0	0	0	0
3	B	160	0	0	9	0
3	C	160	0	0	3	0
3	D	160	0	0	2	0
4	A	32	0	47	5	0
4	B	32	0	45	4	0
4	C	16	0	23	5	0
4	D	16	0	24	0	0
5	A	12	0	19	9	0
5	B	24	0	40	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	6	0	9	0	0
5	D	12	0	20	4	0
6	A	8	0	12	0	0
7	A	294	0	0	8	2
7	B	288	0	0	10	2
7	C	255	0	0	3	1
7	D	249	0	0	7	2
All	All	10389	0	8700	94	5

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.

The worst 5 of 94 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:103:HIS:NE2	5:D:405:BU2:HC11	1.73	1.01
5:B:408[B]:BU2:O1	5:B:408[B]:BU2:O3	1.63	0.98
1:B:295[B]:THR:HG23	7:B:509:HOH:O	1.70	0.92
1:B:234:GLU:OE2	7:B:501:HOH:O	1.91	0.88
3:B:402[A]:BV8:C8	7:B:728:HOH:O	2.28	0.81

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:547:HOH:O	7:D:729:HOH:O[1_455]	1.75	0.45
7:A:608:HOH:O	7:C:696:HOH:O[1_444]	1.76	0.44
7:A:580:HOH:O	7:A:768:HOH:O[1_455]	1.87	0.33
1:B:234:GLU:OE1	7:B:785:HOH:O[1_655]	2.11	0.09
1:D:147[A]:ARG:NH2	7:D:748:HOH:O[1_655]	2.12	0.08

## 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	276/275 (100%)	273 (99%)	3 (1%)	0	100	100
1	B	276/275 (100%)	273 (99%)	3 (1%)	0	100	100
1	C	278/275 (101%)	274 (99%)	3 (1%)	1 (0%)	38	11
1	D	280/275 (102%)	277 (99%)	3 (1%)	0	100	100
All	All	1110/1100 (101%)	1097 (99%)	12 (1%)	1 (0%)	55	20

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	30	ASP

### 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	223/219 (102%)	219 (98%)	4 (2%)	64	24
1	B	224/219 (102%)	222 (99%)	2 (1%)	82	53
1	C	225/219 (103%)	225 (100%)	0	100	100
1	D	226/219 (103%)	225 (100%)	1 (0%)	93	78
All	All	898/876 (102%)	891 (99%)	7 (1%)	89	59

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	284[B]	ARG
1	D	79	LYS
1	B	44	ARG
1	A	259[B]	MET
1	B	290	GLN

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



Mol	Chain	Res	Type
1	A	263	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 47 ligands modelled in this entry, 4 are monoatomic - leaving 43 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	BV8	A	402[A]	-	76,86,86	1.99	14 (18%)	81,122,122	2.05	14 (17%)
3	BV8	A	402[B]	-	76,86,86	2.14	15 (19%)	81,122,122	2.13	18 (22%)
4	EDO	A	403[A]	-	3,3,3	1.20	1 (33%)	2,2,2	0.69	0
4	EDO	A	403[B]	-	3,3,3	1.15	0	2,2,2	0.58	0
4	EDO	A	404[A]	-	3,3,3	0.41	0	2,2,2	1.02	0
4	EDO	A	404[B]	-	3,3,3	0.84	0	2,2,2	2.13	1 (50%)
4	EDO	A	405[A]	-	3,3,3	0.90	0	2,2,2	0.16	0
4	EDO	A	405[B]	-	3,3,3	0.37	0	2,2,2	0.71	0
4	EDO	A	406	-	3,3,3	1.26	0	2,2,2	1.22	0
4	EDO	A	407	-	3,3,3	0.78	0	2,2,2	0.75	0
5	BU2	A	408	-	5,5,5	0.96	0	5,5,5	3.47	1 (20%)
5	BU2	A	409	-	5,5,5	1.40	1 (20%)	5,5,5	1.55	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	DMS	A	410	-	3,3,3	0.49	0	3,3,3	1.64	1 (33%)
6	DMS	A	411	-	3,3,3	0.36	0	3,3,3	1.86	1 (33%)
3	BV8	B	402[A]	-	76,86,86	1.58	10 (13%)	81,122,122	2.32	16 (19%)
3	BV8	B	402[B]	-	76,86,86	1.50	14 (18%)	81,122,122	2.03	16 (19%)
4	EDO	B	403[A]	-	3,3,3	1.26	0	2,2,2	0.59	0
4	EDO	B	403[B]	-	3,3,3	0.87	0	2,2,2	0.35	0
4	EDO	B	404[A]	-	3,3,3	0.36	0	2,2,2	0.49	0
4	EDO	B	404[B]	-	3,3,3	0.40	0	2,2,2	0.59	0
4	EDO	B	405[A]	-	3,3,3	0.69	0	2,2,2	0.10	0
4	EDO	B	405[B]	-	3,3,3	0.84	0	2,2,2	1.00	0
4	EDO	B	406	-	3,3,3	1.34	0	2,2,2	0.73	0
4	EDO	B	407	-	3,3,3	0.50	0	2,2,2	0.79	0
5	BU2	B	408[A]	-	5,5,5	0.42	0	5,5,5	1.05	0
5	BU2	B	408[B]	-	5,5,5	0.48	0	5,5,5	2.72	2 (40%)
5	BU2	B	409	-	5,5,5	0.60	0	5,5,5	1.28	0
5	BU2	B	410	-	5,5,5	0.80	0	5,5,5	0.85	0
3	BV8	C	402[A]	-	76,86,86	1.43	11 (14%)	81,122,122	1.60	12 (14%)
3	BV8	C	402[D]	-	76,86,86	1.45	10 (13%)	81,122,122	1.79	14 (17%)
4	EDO	C	403[A]	-	3,3,3	1.58	1 (33%)	2,2,2	1.10	0
4	EDO	C	403[B]	-	3,3,3	0.84	0	2,2,2	0.49	0
4	EDO	C	404	-	3,3,3	0.95	0	2,2,2	0.38	0
4	EDO	C	405	-	3,3,3	0.79	0	2,2,2	0.40	0
5	BU2	C	406	-	5,5,5	0.81	0	5,5,5	1.55	1 (20%)
3	BV8	D	402[B]	-	76,86,86	1.79	15 (19%)	81,122,122	1.89	14 (17%)
3	BV8	D	402[C]	-	76,86,86	1.83	16 (21%)	81,122,122	2.05	18 (22%)
4	EDO	D	403[A]	-	3,3,3	0.75	0	2,2,2	0.58	0
4	EDO	D	403[B]	-	3,3,3	0.92	0	2,2,2	0.61	0
4	EDO	D	404[A]	-	3,3,3	0.85	0	2,2,2	0.27	0
4	EDO	D	404[B]	-	3,3,3	0.47	0	2,2,2	1.38	0
5	BU2	D	405	-	5,5,5	1.31	0	5,5,5	2.43	2 (40%)
5	BU2	D	406	-	5,5,5	0.61	0	5,5,5	1.96	2 (40%)

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	BV8	A	402[A]	-	-	0/56/101/101	0/7/7/7
3	BV8	A	402[B]	-	-	0/56/101/101	0/7/7/7

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	403[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	A	403[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	A	404[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	A	404[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	A	405[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	A	405[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	A	406	-	-	0/1/1/1	0/0/0/0
4	EDO	A	407	-	-	0/1/1/1	0/0/0/0
5	BU2	A	408	-	-	0/3/3/3	0/0/0/0
5	BU2	A	409	-	-	0/3/3/3	0/0/0/0
6	DMS	A	410	-	-	0/0/0/0	0/0/0/0
6	DMS	A	411	-	-	0/0/0/0	0/0/0/0
3	BV8	B	402[A]	-	-	0/56/101/101	0/7/7/7
3	BV8	B	402[B]	-	-	0/56/101/101	0/7/7/7
4	EDO	B	403[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	403[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	404[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	404[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	405[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	405[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	406	-	-	0/1/1/1	0/0/0/0
4	EDO	B	407	-	-	0/1/1/1	0/0/0/0
5	BU2	B	408[A]	-	-	0/3/3/3	0/0/0/0
5	BU2	B	408[B]	-	-	0/3/3/3	0/0/0/0
5	BU2	B	409	-	-	0/3/3/3	0/0/0/0
5	BU2	B	410	-	-	0/3/3/3	0/0/0/0
3	BV8	C	402[A]	-	-	0/56/101/101	0/7/7/7
3	BV8	C	402[D]	-	-	0/56/101/101	0/7/7/7
4	EDO	C	403[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	C	403[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	C	404	-	-	0/1/1/1	0/0/0/0
4	EDO	C	405	-	-	0/1/1/1	0/0/0/0
5	BU2	C	406	-	-	0/3/3/3	0/0/0/0
3	BV8	D	402[B]	-	-	0/56/101/101	0/7/7/7
3	BV8	D	402[C]	-	-	0/56/101/101	0/7/7/7
4	EDO	D	403[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	D	403[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	D	404[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	D	404[B]	-	-	0/1/1/1	0/0/0/0
5	BU2	D	405	-	-	0/3/3/3	0/0/0/0
5	BU2	D	406	-	-	0/3/3/3	0/0/0/0

The worst 5 of 108 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402[A]	BV8	C6-C11	-6.09	1.31	1.41
3	A	402[B]	BV8	C6-C11	-6.00	1.31	1.41
3	D	402[B]	BV8	C41-C42	-5.45	1.45	1.53
3	A	402[B]	BV8	C41-C42	-5.42	1.45	1.53
3	D	402[C]	BV8	C41-C42	-5.09	1.45	1.53

The worst 5 of 134 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402[B]	BV8	N9-C46-N10	-12.92	117.60	128.86
3	B	402[A]	BV8	N9-C46-N10	-12.35	118.11	128.86
3	B	402[B]	BV8	N9-C46-N10	-12.26	118.18	128.86
3	A	402[A]	BV8	N9-C46-N10	-12.14	118.29	128.86
3	D	402[B]	BV8	N9-C46-N10	-9.98	120.17	128.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

18 monomers are involved in 50 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	403[A]	EDO	2	0
4	A	403[B]	EDO	1	0
4	A	404[A]	EDO	2	0
4	A	406	EDO	2	0
5	A	408	BU2	9	0
3	B	402[A]	BV8	7	0
3	B	402[B]	BV8	2	0
4	B	404[B]	EDO	1	0
4	B	406	EDO	4	0
5	B	408[A]	BU2	4	0
5	B	408[B]	BU2	2	0
5	B	409	BU2	3	0
5	B	410	BU2	4	0
3	C	402[A]	BV8	2	0
3	C	402[D]	BV8	1	0
4	C	404	EDO	5	0
3	D	402[B]	BV8	2	0
5	D	405	BU2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	269/275 (97%)	-0.28	2 (0%) 87 86	8, 16, 30, 63	0
1	B	268/275 (97%)	-0.30	1 (0%) 92 91	10, 17, 31, 76	0
1	C	274/275 (99%)	-0.07	7 (2%) 56 57	10, 20, 42, 97	0
1	D	275/275 (100%)	-0.06	9 (3%) 47 47	8, 19, 41, 113	0
All	All	1086/1100 (98%)	-0.18	19 (1%) 70 71	8, 17, 37, 113	0

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	29	ILE	15.8
1	D	29	ILE	11.2
1	D	28	GLY	7.8
1	A	34	ALA	7.8
1	B	35	ARG	4.6

### 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	A	404[B]	4/4	0.93	0.27	18.25	16,18,20,25	4
4	EDO	B	405[B]	4/4	0.85	0.23	17.43	27,29,41,42	4
4	EDO	B	405[A]	4/4	0.85	0.23	16.47	33,43,46,49	4
4	EDO	A	404[A]	4/4	0.93	0.27	15.35	13,22,25,33	4
4	EDO	A	405[A]	4/4	0.83	0.16	13.38	30,36,42,46	4
4	EDO	B	404[A]	4/4	0.90	0.13	13.08	26,32,34,35	4
4	EDO	C	403[B]	4/4	0.96	0.17	12.63	10,11,15,15	4
4	EDO	D	403[B]	4/4	0.96	0.16	12.25	16,19,20,22	4
4	EDO	D	403[A]	4/4	0.96	0.16	12.03	10,10,12,13	4
4	EDO	A	405[B]	4/4	0.83	0.16	11.98	20,27,32,34	4
4	EDO	C	403[A]	4/4	0.96	0.17	9.61	13,14,15,17	4
4	EDO	B	404[B]	4/4	0.90	0.13	9.01	15,22,26,29	4
5	BU2	A	408	6/6	0.86	0.24	8.97	29,35,44,59	0
4	EDO	B	403[B]	4/4	0.98	0.17	8.82	10,12,14,20	4
4	EDO	B	403[A]	4/4	0.98	0.17	8.70	10,10,10,13	4
5	BU2	B	408[A]	6/6	0.96	0.12	8.09	21,24,31,43	6
5	BU2	B	408[B]	6/6	0.96	0.12	7.75	25,27,28,34	6
5	BU2	D	406	6/6	0.93	0.15	7.25	44,56,64,82	0
4	EDO	A	403[A]	4/4	0.96	0.15	6.82	6,7,9,12	4
5	BU2	B	409	6/6	0.92	0.12	6.46	31,39,44,51	0
4	EDO	A	403[B]	4/4	0.96	0.15	6.44	15,16,21,28	4
4	EDO	A	406	4/4	0.97	0.11	6.01	24,29,29,60	0
5	BU2	D	405	6/6	0.83	0.15	4.84	32,43,71,110	0
4	EDO	D	404[A]	4/4	0.93	0.12	4.36	22,23,25,27	4
4	EDO	D	404[B]	4/4	0.93	0.12	4.09	28,29,30,34	4
5	BU2	A	409	6/6	0.81	0.13	4.00	38,51,56,60	0
4	EDO	C	405	4/4	0.86	0.14	3.47	33,41,42,58	0
5	BU2	C	406	6/6	0.77	0.14	3.10	32,41,48,50	0
4	EDO	A	407	4/4	0.94	0.12	1.90	32,37,46,55	0
6	DMS	A	410	4/4	0.91	0.13	1.78	51,54,67,79	0
5	BU2	B	410	6/6	0.86	0.10	1.52	32,39,50,57	0
3	BV8	D	402[B]	80/80	0.98	0.07	0.25	8,12,19,24	80
3	BV8	D	402[C]	80/80	0.98	0.07	0.25	9,17,24,36	80
3	BV8	A	402[B]	80/80	0.98	0.07	0.12	5,9,34,46	80
3	BV8	A	402[A]	80/80	0.98	0.07	0.12	13,19,33,39	80
3	BV8	B	402[A]	80/80	0.99	0.07	0.09	14,19,38,51	80
3	BV8	B	402[B]	80/80	0.99	0.07	0.09	7,11,33,40	80
3	BV8	C	402[D]	80/80	0.98	0.07	0.01	7,18,24,28	80
3	BV8	C	402[A]	80/80	0.98	0.07	0.01	8,13,24,32	80
4	EDO	B	406	4/4	0.98	0.06	-0.09	26,27,29,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	D	401	1/1	1.00	0.04	-1.45	17,17,17,17	0
2	ZN	A	401	1/1	1.00	0.05	-1.53	15,15,15,15	0
2	ZN	C	401	1/1	1.00	0.04	-1.98	19,19,19,19	0
2	ZN	B	401	1/1	1.00	0.04	-2.07	16,16,16,16	0
4	EDO	C	404	4/4	0.77	0.17	-	55,57,68,73	0
6	DMS	A	411	4/4	0.93	0.12	-	56,71,83,86	0
4	EDO	B	407	4/4	0.86	0.18	-	51,55,55,92	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.