



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 09:57 pm BST

PDB ID : 6FKZ  
Title : Crystal structure of zebrafish Sirtuin 5 in complex with 3-(phenylthio)succinyl-CPS1 peptide  
Authors : Pannek, M.; Steegborn, C.  
Deposited on : 2018-01-25  
Resolution : 3.30 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](mailto: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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

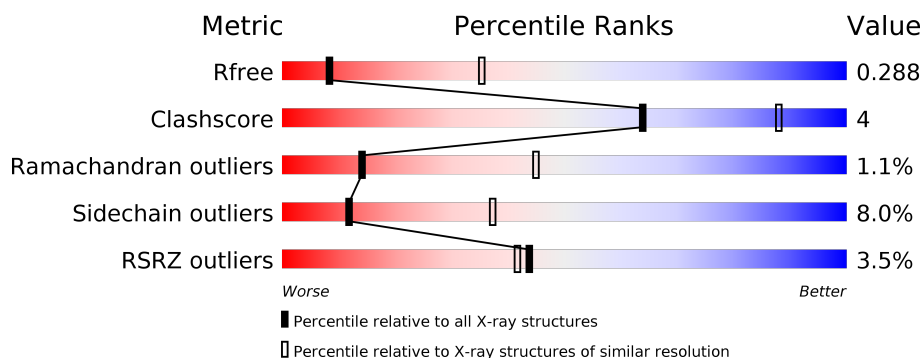
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 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	284	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 16%, green 75%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>75%</span> <span>16%</span> <span>• 7%</span> </div> </div>
1	B	284	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 5%, orange 1%, yellow 13%, green 80%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>5%</span> <span>80%</span> <span>13%</span> <span>• 6%</span> </div> </div>
2	E	8	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 13%, orange 1%, yellow 38%, green 63%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>13%</span> <span>63%</span> <span>38%</span> </div> </div>
3	H	8	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 13%, orange 1%, yellow 38%, green 63%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>13%</span> <span>63%</span> <span>38%</span> </div> </div>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4364 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 deacylase sirtuin-5, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	1	0
			2062	1302	373	372	15			
1	B	266	Total	C	N	O	S	0	1	0
			2071	1307	374	375	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	GLY	-	expression tag	UNP Q6DHI5
A	23	ILE	-	expression tag	UNP Q6DHI5
A	24	ASP	-	expression tag	UNP Q6DHI5
A	25	PRO	-	expression tag	UNP Q6DHI5
A	26	PHE	-	expression tag	UNP Q6DHI5
A	27	THR	-	expression tag	UNP Q6DHI5
B	22	GLY	-	expression tag	UNP Q6DHI5
B	23	ILE	-	expression tag	UNP Q6DHI5
B	24	ASP	-	expression tag	UNP Q6DHI5
B	25	PRO	-	expression tag	UNP Q6DHI5
B	26	PHE	-	expression tag	UNP Q6DHI5
B	27	THR	-	expression tag	UNP Q6DHI5

- Molecule 2 is a protein called 3(S)-(phenylthio)succinyl-CPS1 peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	8	Total	C	N	O	S	0	8	0
			82	57	9	15	1			

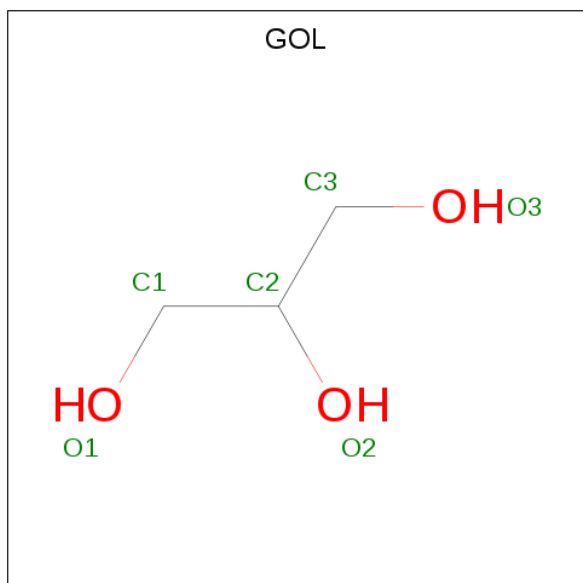
- Molecule 3 is a protein called 3(R)-(phenylthio)succinyl-CPS1 peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	8	Total	C	N	O	S	0	8	0
			82	57	9	15	1			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	1
			1	1		
4	A	1	Total	Zn	0	1
			1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C<sub>8</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	1
			15	8	2	4	1		

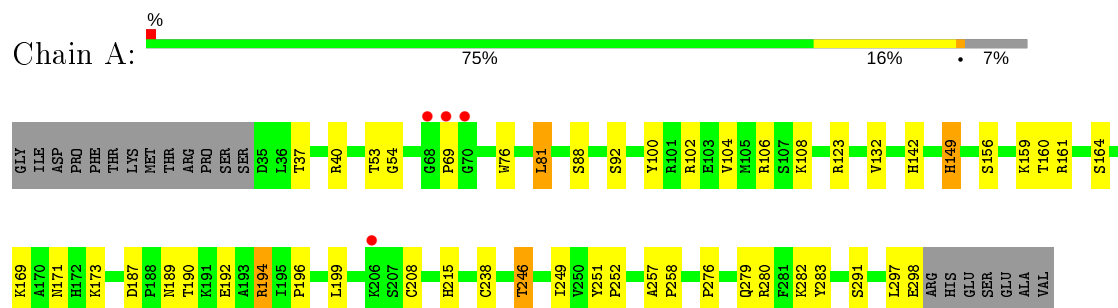
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	30	Total	O	0	0
			30	30		
7	B	13	Total	O	0	0
			13	13		
7	E	1	Total	O	0	0
			1	1		

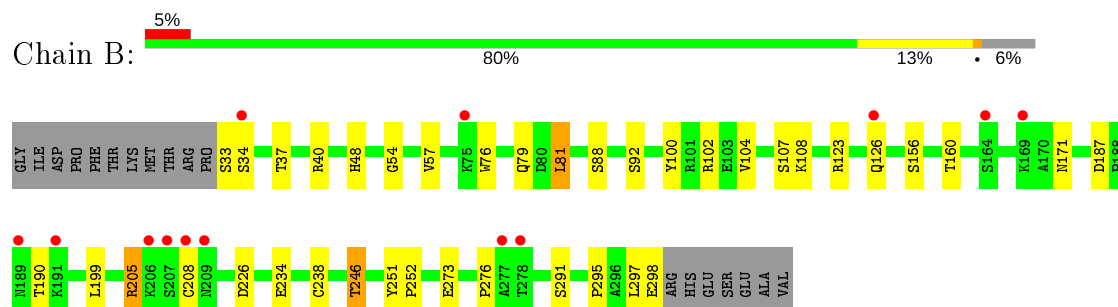
### 3 Residue-property plots [i](#)

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.

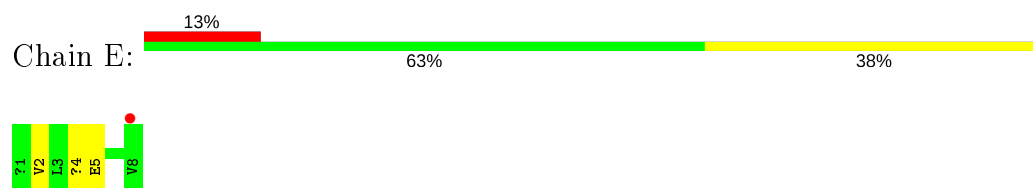
- Molecule 1: NAD-dependent protein deacylase sirtuin-5, mitochondrial



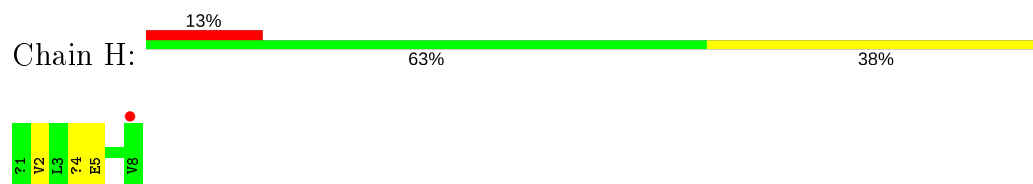
- Molecule 1: NAD-dependent protein deacylase sirtuin-5, mitochondrial



- Molecule 2: 3(S)-(phenylthio)succinyl-CPS1 peptide



- Molecule 3: 3(R)-(phenylthio)succinyl-CPS1 peptide



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.22Å 87.22Å 318.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.73 – 3.30 19.90 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.0 (48.73-3.30) 99.7 (19.90-3.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.20 (at 3.29Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.222 , 0.284 0.227 , 0.288	Depositor DCC
$R_{free}$ test set	573 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.5	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 76.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4364	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	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.*

<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: GOL, ZN, EPE, GZB, EYZ, DQK

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	0.62	0/2118	0.86	0/2873
1	B	0.61	0/2130	0.88	3/2889 (0.1%)
2	E	0.36	0/46	0.58	0/60
3	H	0.36	0/46	0.58	0/60
All	All	0.61	0/4340	0.87	3/5882 (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	B	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	226	ASP	CB-CG-OD2	-6.13	112.78	118.30
1	B	226	ASP	CB-CG-OD1	5.82	123.54	118.30
1	B	199	LEU	CA-CB-CG	5.53	128.01	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	33	SER	Peptide



## 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	2062	0	2024	20	0
1	B	2071	0	2035	13	0
2	E	82	0	46	1	0
3	H	82	0	46	4	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	6	0	8	0	0
6	B	15	0	18	0	0
7	A	30	0	0	0	0
7	B	13	0	0	0	0
7	E	1	0	0	0	0
All	All	4364	0	4177	36	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 (36) 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)
3:H:4[B]:EYZ:OAO	3:H:4[B]:EYZ:CAW	1.83	1.26
3:H:4[B]:EYZ:O	3:H:5[B]:GLU:HB2	1.97	0.64
2:E:4[A]:DQK:O	2:E:5[A]:GLU:HB2	2.02	0.59
1:B:54:GLY:HA3	1:B:246:THR:HB	1.84	0.59
1:A:54:GLY:HA3	1:A:246:THR:HB	1.87	0.56
1:A:196:PRO:HB2	1:A:199:LEU:HD12	1.85	0.56
1:A:298:GLU:OE2	1:B:107:SER:HB3	2.11	0.50
3:H:4[B]:EYZ:CAN	3:H:4[B]:EYZ:CAW	2.80	0.50
1:A:132:VAL:HG12	1:A:149[A]:HIS:CG	2.47	0.50
1:A:76:TRP:HB2	1:A:81:LEU:HD11	1.94	0.49
1:A:251:TYR:CE1	1:A:252:PRO:HG3	2.48	0.48
3:H:4[B]:EYZ:OAO	3:H:4[B]:EYZ:CAR	2.48	0.48
1:B:76:TRP:HB2	1:B:81:LEU:HD11	1.95	0.48
1:A:257:ALA:N	1:A:258:PRO:HD2	2.30	0.47
1:A:123:ARG:NH2	1:A:297:LEU:O	2.47	0.47
1:B:123:ARG:NH2	1:B:297:LEU:O	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ILE:HG22	1:A:249:ILE:O	2.16	0.45
1:A:192:GLU:OE1	1:A:194:ARG:NH2	2.40	0.45
1:B:48[B]:HIS:NE2	1:B:234:GLU:OE1	2.49	0.45
1:A:100:TYR:O	1:A:104:VAL:HG13	2.17	0.43
1:A:187:ASP:OD2	1:A:189:ASN:ND2	2.40	0.43
1:A:53:THR:HG21	1:A:142:HIS:CE1	2.55	0.42
1:B:100:TYR:O	1:B:104:VAL:HG13	2.19	0.42
1:A:104:VAL:O	1:A:108:LYS:HG2	2.20	0.41
1:B:205:ARG:CZ	1:B:205:ARG:HB3	2.50	0.41
1:B:104:VAL:O	1:B:108:LYS:HG2	2.21	0.41
1:B:102:ARG:NH2	1:B:171:ASN:OD1	2.54	0.41
1:B:295:PRO:O	1:B:298:GLU:HG2	2.20	0.41
1:A:102:ARG:NH2	1:A:171:ASN:OD1	2.53	0.41
1:B:37:THR:HA	1:B:40:ARG:HG3	2.03	0.41
1:A:37:THR:HA	1:A:40:ARG:HG3	2.01	0.41
1:A:106:ARG:HH22	1:B:123:ARG:HD2	1.85	0.41
1:B:251:TYR:HA	1:B:252:PRO:HA	1.87	0.40
1:A:161:ARG:NH1	1:A:215:HIS:ND1	2.66	0.40
1:A:53:THR:HG21	1:A:142:HIS:NE2	2.36	0.40
1:A:249:ILE:CG2	1:A:249:ILE:O	2.69	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	263/284 (93%)	240 (91%)	20 (8%)	3 (1%)	14	45
1	B	265/284 (93%)	239 (90%)	23 (9%)	3 (1%)	14	45
2	E	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
3	H	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
All	All	538/584 (92%)	487 (90%)	45 (8%)	6 (1%)	14	45

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	69	PRO
1	A	190	THR
1	B	34	SER
1	B	190	THR
1	B	276	PRO
1	A	276	PRO

### 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	220/237 (93%)	200 (91%)	20 (9%)	9	31
1	B	222/237 (94%)	207 (93%)	15 (7%)	16	44
2	E	5/5 (100%)	4 (80%)	1 (20%)	1	5
3	H	5/5 (100%)	4 (80%)	1 (20%)	1	5
All	All	452/484 (93%)	415 (92%)	37 (8%)	12	36

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	LEU
1	A	88	SER
1	A	92	SER
1	A	149[A]	HIS
1	A	149[B]	HIS
1	A	156	SER
1	A	159	LYS
1	A	160	THR
1	A	164	SER
1	A	169	LYS
1	A	173	LYS
1	A	194	ARG
1	A	208	CYS
1	A	238	CYS

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Mol	Chain	Res	Type
1	A	246	THR
1	A	279	GLN
1	A	280	ARG
1	A	282	LYS
1	A	283	TYR
1	A	291	SER
1	B	57	VAL
1	B	79	GLN
1	B	81	LEU
1	B	88	SER
1	B	92	SER
1	B	126	GLN
1	B	156	SER
1	B	160	THR
1	B	187	ASP
1	B	205	ARG
1	B	208	CYS
1	B	238	CYS
1	B	246	THR
1	B	273	GLU
1	B	291	SER
2	E	2[A]	VAL
3	H	2[B]	VAL

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

Mol	Chain	Res	Type
1	B	42	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	$\# Z  > 2$
2	GZB	E	1[A]	2	12,12,13	1.61	1 (8%)	14,14,16	0.91	0
3	GZB	H	1[B]	3	12,12,13	1.62	1 (8%)	14,14,16	0.88	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	GZB	E	1[A]	2	-	4/7/8/9	0/1/1/1
3	GZB	H	1[B]	3	-	4/7/8/9	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	1[B]	GZB	CAG-CAE	-5.41	1.38	1.50
2	E	1[A]	GZB	CAG-CAE	-5.37	1.39	1.50

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1[A]	GZB	N-CAE-CAG-CAL
2	E	1[A]	GZB	OAF-CAE-CAG-CAL
2	E	1[A]	GZB	OAF-CAE-CAG-CAH
3	H	1[B]	GZB	OAF-CAE-CAG-CAL
2	E	1[A]	GZB	N-CAE-CAG-CAH
3	H	1[B]	GZB	N-CAE-CAG-CAL
3	H	1[B]	GZB	N-CAE-CAG-CAH
3	H	1[B]	GZB	OAF-CAE-CAG-CAH

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z  > 2$
6	EPE	B	402[A]	-	15,15,15	1.92	1 (6%)	18,20,20	1.38	1 (5%)
5	GOL	A	402	-	5,5,5	0.52	0	5,5,5	0.42	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
6	EPE	B	402[A]	-	-	4/9/19/19	0/1/1/1
5	GOL	A	402	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	402[A]	EPE	C10-S	-6.94	1.67	1.77

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	402[A]	EPE	O2S-S-C10	4.51	112.34	106.92

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	402[A]	EPE	C9-C10-S-O3S
5	A	402	GOL	O1-C1-C2-O2
5	A	402	GOL	O2-C2-C3-O3
6	B	402[A]	EPE	C9-C10-S-O1S
6	B	402[A]	EPE	C9-C10-S-O2S
6	B	402[A]	EPE	C10-C9-N1-C2

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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	264/284 (92%)	-0.18	4 (1%) 73 72	36, 52, 86, 125	0
1	B	266/284 (93%)	0.05	13 (4%) 29 27	38, 64, 104, 120	1 (0%)
2	E	6/8 (75%)	0.33	1 (16%) 1 1	70, 72, 77, 78	6 (100%)
3	H	6/8 (75%)	0.30	1 (16%) 1 1	68, 69, 76, 76	6 (100%)
All	All	542/584 (92%)	-0.06	19 (3%) 44 42	36, 57, 100, 125	13 (2%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	207	SER	3.5
1	B	206	LYS	3.3
1	A	70	GLY	3.2
1	B	278	THR	3.2
1	A	69	PRO	3.1
1	B	75	LYS	2.8
1	A	206	LYS	2.7
1	B	126	GLN	2.6
1	A	68	GLY	2.5
1	B	34	SER	2.4
1	B	208	CYS	2.3
2	E	8[A]	VAL	2.2
3	H	8[B]	VAL	2.2
1	B	191	LYS	2.2
1	B	209	ASN	2.2
1	B	189	ASN	2.1
1	B	164	SER	2.1
1	B	169	LYS	2.0
1	B	277	ALA	2.0



## 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	GZB	E	1[A]	12/13	0.87	0.30	65,75,76,78	12
3	GZB	H	1[B]	12/13	0.87	0.30	64,73,74,76	12

## 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
5	GOL	A	402	6/6	0.76	0.36	60,63,64,64	0
6	EPE	B	402[A]	15/15	0.87	0.25	65,66,77,78	0
4	ZN	B	401[A]	1/1	0.96	0.10	92,92,92,92	0
4	ZN	A	401[A]	1/1	0.98	0.03	42,42,42,42	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.