



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 06:44 PM GMT

PDB ID : 1C72  
Title : TYR115, GLN165 AND TRP209 CONTRIBUTE TO THE 1,2-EPOXY-3-(P-NITROPHENOXY)PROPANE CONJUGATING ACTIVITIES OF GLUTATHIONE S-TRANSFERASE CGSTM1-1  
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Deposited on : 2000-02-02  
Resolution : 2.80 Å(reported)

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

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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 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
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)	:	trunk26865

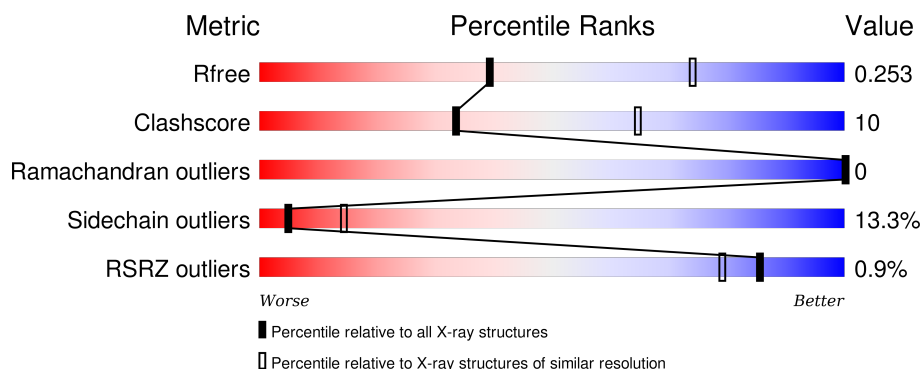
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	219	<div> <div>2%</div> <div>70%</div> <div>23%</div> <div>6%</div> <div>.</div> </div>
1	B	219	<div> <div>69%</div> <div>26%</div> <div>..</div> </div>
1	C	219	<div> <div>%</div> <div>73%</div> <div>22%</div> <div>..</div> </div>
1	D	219	<div> <div>%</div> <div>71%</div> <div>26%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EPY	A	220	X	-	-	X
2	EPY	B	221	X	-	-	X
2	EPY	C	222	X	-	-	X
2	EPY	D	223	-	-	-	X

## 2 Entry composition [i](#)

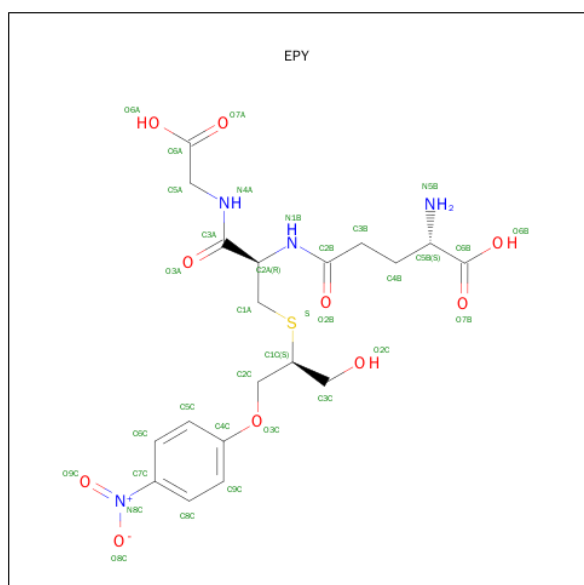
There are 3 unique types of molecules in this entry. The entry contains 7490 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (GLUTATHIONE S-TRANSFERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1799	1164	306	320	9			
1	B	217	Total	C	N	O	S	0	0	0
			1799	1164	306	320	9			
1	C	217	Total	C	N	O	S	0	0	0
			1799	1164	306	320	9			
1	D	217	Total	C	N	O	S	0	0	0
			1799	1164	306	320	9			

- Molecule 2 is 1-HYDROXY-2-S-GLUTATHIONYL-3-PARA-NITROPHENOXY-PROPANE (three-letter code: EPY) (formula:  $C_{19}H_{26}N_4O_{10}S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			34	19	4	10	1		
2	B	1	Total	C	N	O	S	0	0
			34	19	4	10	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			34	19	4	10	1		
2	D	1	Total	C	N	O	S	0	0
			34	19	4	10	1		

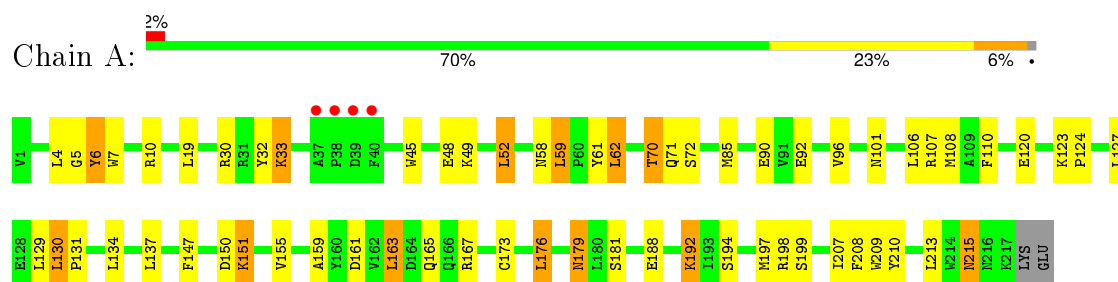
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total	O	0	0
			30	30		
3	B	42	Total	O	0	0
			42	42		
3	C	50	Total	O	0	0
			50	50		
3	D	36	Total	O	0	0
			36	36		

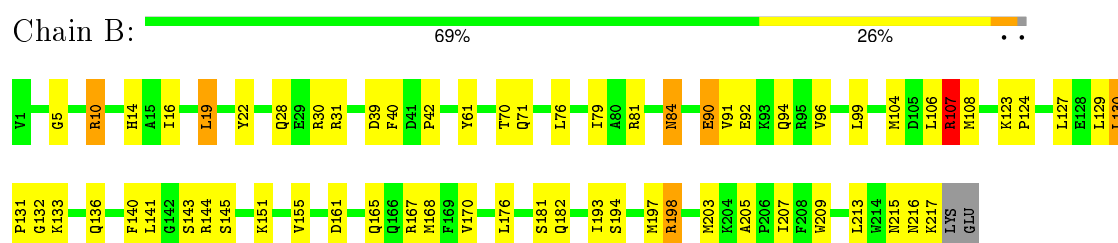
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

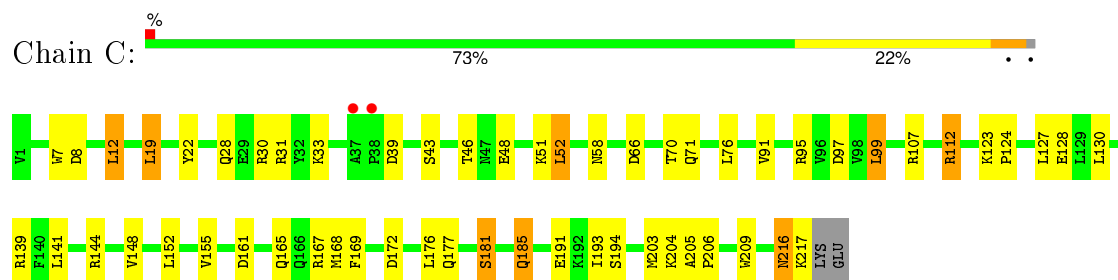
#### • Molecule 1: PROTEIN (GLUTATHIONE S-TRANSFERASE)



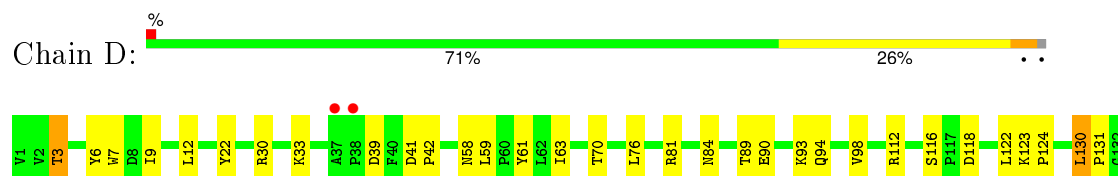
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K133	L134	Q135	Q136	L137	S138	R139
R144	S145	M146	F147	V148	G149	
T153	F154	V155		L158		
V170		S181		E191	K192	I193
S194		M197	R198	A205	P206	I207
F208	H209	T210	T211	N215	N216	T217
LYS	GLU					

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	125.60Å 136.65Å 125.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.00 – 2.80 31.05 – 2.81	Depositor EDS
% Data completeness (in resolution range)	86.6 (31.00-2.80) 81.8 (31.05-2.81)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.08 (at 2.81Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.190 , 0.280 0.171 , 0.253	Depositor DCC
$R_{free}$ test set	2163 reflections (9.90%)	DCC
Wilson B-factor (Å <sup>2</sup> )	42.1	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 62.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 23375 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7490	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: EPY

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	0.44	1/1847 (0.1%)	0.56	0/2499
1	B	0.40	1/1847 (0.1%)	0.57	2/2499 (0.1%)
1	C	0.39	0/1847	0.52	0/2499
1	D	0.40	0/1847	0.51	0/2499
All	All	0.41	2/7388 (0.0%)	0.54	2/9996 (0.0%)

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	2
1	B	0	2
1	D	0	1
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	165	GLN	CD-OE1	6.63	1.38	1.24
1	B	107	ARG	CD-NE	-5.71	1.36	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	ARG	CD-NE-CZ	-10.23	109.28	123.60
1	B	107	ARG	CG-CD-NE	6.03	124.45	111.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	32	TYR	Sidechain
1	A	6	TYR	Sidechain
1	B	107	ARG	Sidechain
1	B	198	ARG	Sidechain
1	D	61	TYR	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	1799	0	1784	39	0
1	B	1799	0	1784	32	0
1	C	1799	0	1784	32	0
1	D	1799	0	1784	43	0
2	A	34	0	23	6	0
2	B	34	0	23	6	0
2	C	34	0	22	4	0
2	D	34	0	24	13	0
3	A	30	0	0	0	0
3	B	42	0	0	0	0
3	C	50	0	0	0	0
3	D	36	0	0	0	0
All	All	7490	0	7228	149	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 10.

The worst 5 of 149 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:C:7:TRP:CZ2	2:C:222:EPY:H6C1	2.04	0.93
1:B:107:ARG:HD2	2:B:221:EPY:O8C	1.69	0.92
1:D:7:TRP:CZ2	2:D:223:EPY:H6C1	2.11	0.84
1:D:216:ASN:H	1:D:216:ASN:HD22	1.28	0.80
2:D:223:EPY:C2C	2:D:223:EPY:C2A	2.56	0.79

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	215/219 (98%)	209 (97%)	6 (3%)	0	100	100
1	B	215/219 (98%)	205 (95%)	10 (5%)	0	100	100
1	C	215/219 (98%)	208 (97%)	7 (3%)	0	100	100
1	D	215/219 (98%)	205 (95%)	10 (5%)	0	100	100
All	All	860/876 (98%)	827 (96%)	33 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	192/194 (99%)	166 (86%)	26 (14%)	5	14
1	B	192/194 (99%)	165 (86%)	27 (14%)	4	12
1	C	192/194 (99%)	165 (86%)	27 (14%)	4	12
1	D	192/194 (99%)	170 (88%)	22 (12%)	7	21
All	All	768/776 (99%)	666 (87%)	102 (13%)	5	14

5 of 102 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	181	SER
1	C	39	ASP
1	D	155	VAL
1	B	182	GLN
1	B	215	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	94	GLN
1	C	216	ASN
1	D	215	ASN
1	C	185	GLN
1	D	47	ASN

### 5.3.3 RNA [i](#)

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

4 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	EPY	A	220	-	23,34,34	1.96	5 (21%)	29,44,44	2.85	7 (24%)
2	EPY	B	221	-	23,34,34	1.21	2 (8%)	29,44,44	4.36	10 (34%)
2	EPY	C	222	-	23,34,34	2.12	6 (26%)	29,44,44	3.26	8 (27%)
2	EPY	D	223	-	23,34,34	2.38	5 (21%)	29,44,44	2.36	10 (34%)

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	EPY	A	220	-	1/1/8/10	0/32/38/38	0/1/1/1
2	EPY	B	221	-	1/1/8/10	0/32/38/38	0/1/1/1
2	EPY	C	222	-	1/1/8/10	0/32/38/38	0/1/1/1
2	EPY	D	223	-	-	0/32/38/38	0/1/1/1

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	223	EPY	C1A-C2A	-5.80	1.37	1.53
2	D	223	EPY	O2B-C2B	-5.35	1.12	1.23
2	C	222	EPY	C1A-C2A	-4.30	1.41	1.53
2	C	222	EPY	C3C-C1C	-4.04	1.42	1.52
2	D	223	EPY	C8C-C9C	-3.69	1.32	1.38

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	220	EPY	C2C-O3C-C4C	-9.93	96.54	117.89
2	A	220	EPY	C1A-S-C1C	-8.34	87.70	101.42
2	B	221	EPY	C1A-C2A-N1B	-6.94	93.75	110.83
2	D	223	EPY	C1A-C2A-N1B	-5.05	98.39	110.83
2	B	221	EPY	O3A-C3A-C2A	-4.98	109.35	120.36

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	221	EPY	C1C
2	A	220	EPY	C1C
2	C	222	EPY	C1C

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	220	EPY	6	0
2	B	221	EPY	6	0
2	C	222	EPY	4	0
2	D	223	EPY	13	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	217/219 (99%)	-0.60	4 (1%) 71 61	9, 29, 54, 67	0
1	B	217/219 (99%)	-0.67	0 100 100	10, 31, 52, 68	0
1	C	217/219 (99%)	-0.70	2 (0%) 85 79	9, 29, 51, 70	0
1	D	217/219 (99%)	-0.62	2 (0%) 85 79	8, 28, 54, 74	0
All	All	868/876 (99%)	-0.65	8 (0%) 85 79	8, 30, 53, 74	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	38	PRO	3.3
1	D	38	PRO	2.8
1	A	37	ALA	2.3
1	A	40	PHE	2.3
1	C	37	ALA	2.2

### 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( $\text{\AA}^2$ )	Q<0.9
2	EPY	C	222	34/34	0.84	0.23	3.28	14,35,53,56	0
2	EPY	A	220	34/34	0.86	0.22	2.50	25,33,51,55	0
2	EPY	B	221	34/34	0.86	0.22	2.42	21,36,51,59	0
2	EPY	D	223	34/34	0.84	0.23	2.24	28,39,51,54	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.