



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2024 – 07:55 AM EDT

PDB ID : 3WPS  
Title : crystal structure of the GAP domain of MgcRacGAP(S387D)  
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Deposited on : 2014-01-15  
Resolution : 2.70 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

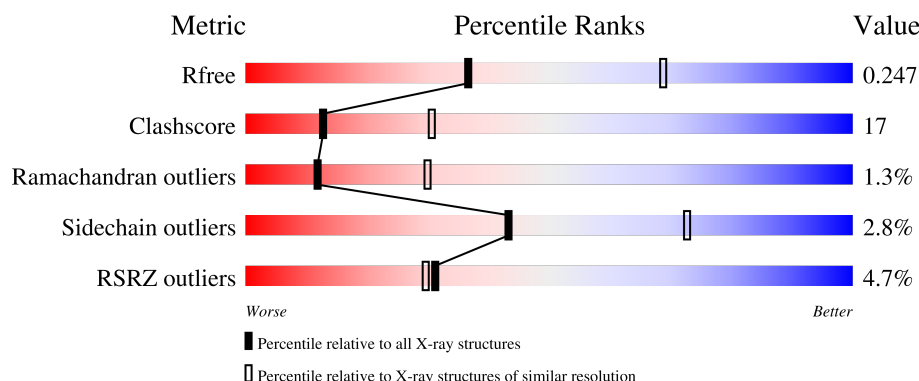
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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  $\geq 3$ , 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	208	<div> <div>5%</div> <div>67%</div> <div>28%</div> <div>• •</div> </div>
1	B	208	<div> <div>4%</div> <div>64%</div> <div>29%</div> <div>• •</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3279 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 Rac GTPase-activating protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	Se	0	0	0
			1588	1010	272	294	3	9			
1	B	201	Total	C	N	O	S	Se	0	0	0
			1588	1010	272	294	3	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	GLY	-	expression tag	UNP Q9H0H5
A	340	SER	-	expression tag	UNP Q9H0H5
A	341	SER	-	expression tag	UNP Q9H0H5
A	342	GLY	-	expression tag	UNP Q9H0H5
A	343	SER	-	expression tag	UNP Q9H0H5
A	344	SER	-	expression tag	UNP Q9H0H5
A	345	GLY	-	expression tag	UNP Q9H0H5
A	387	ASP	SER	engineered mutation	UNP Q9H0H5
A	518	SER	LEU	SEE REMARK 999	UNP Q9H0H5
B	339	GLY	-	expression tag	UNP Q9H0H5
B	340	SER	-	expression tag	UNP Q9H0H5
B	341	SER	-	expression tag	UNP Q9H0H5
B	342	GLY	-	expression tag	UNP Q9H0H5
B	343	SER	-	expression tag	UNP Q9H0H5
B	344	SER	-	expression tag	UNP Q9H0H5
B	345	GLY	-	expression tag	UNP Q9H0H5
B	387	ASP	SER	engineered mutation	UNP Q9H0H5
B	518	SER	LEU	SEE REMARK 999	UNP Q9H0H5

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	49	Total	O	0	0
			49	49		
3	B	44	Total	O	0	0
			44	44		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.32Å 77.32Å 108.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.66 – 2.70 38.66 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.66-2.70) 99.9 (38.66-2.71)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.13	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.50 (at 2.72Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.210 , 0.259 0.207 , 0.247	Depositor DCC
$R_{free}$ test set	960 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.7	Xtriage
Anisotropy	0.470	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -h,-k,l 0.047 for h,-h-k,-l 0.023 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3279	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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: SO4

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	0.35	0/1608	0.61	0/2163
1	B	0.34	0/1608	0.59	0/2163
All	All	0.35	0/3216	0.60	0/4326

There are no bond length outliers.

There are no bond angle outliers.

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	1588	0	1625	50	0
1	B	1588	0	1625	58	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	49	0	0	2	0
3	B	44	0	0	2	0
All	All	3279	0	3250	107	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 17.

All (107) 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:B:408:LEU:HB2	1:B:411:LYS:HE3	1.46	0.96
1:A:492:ASP:H	1:A:495:ASN:HD22	1.25	0.85
1:A:396:LEU:O	1:A:400:PHE:HB2	1.82	0.79
1:B:405:THR:O	1:B:407:PRO:HD3	1.82	0.78
1:B:372:GLU:HG3	1:B:409:LEU:HB2	1.65	0.78
1:A:440:ARG:HD3	1:A:443:MSE:CE	2.15	0.77
1:B:443:MSE:HE2	1:B:523:ARG:HB2	1.66	0.76
1:A:440:ARG:HD3	1:A:443:MSE:HE2	1.70	0.73
1:B:475:PHE:CD1	1:B:543:MSE:HE1	2.24	0.71
1:A:406:VAL:HB	1:A:407:PRO:HD3	1.71	0.71
1:B:347:GLY:O	1:B:348:GLU:HG3	1.92	0.70
1:A:346:ILE:N	1:A:430:LYS:HA	2.08	0.68
1:A:385:ARG:HB2	1:A:491:MSE:CE	2.26	0.66
1:A:424:ASP:HA	1:A:427:ARG:HH12	1.60	0.66
1:A:424:ASP:HA	1:A:427:ARG:NH1	2.11	0.66
1:B:458:MSE:HE2	1:B:530:ARG:HB2	1.78	0.66
1:B:545:VAL:O	1:B:546:GLU:HB2	1.95	0.65
1:B:492:ASP:OD1	1:B:495:ASN:ND2	2.31	0.64
1:A:477:MSE:CE	1:A:480:LEU:HD12	2.28	0.64
1:B:396:LEU:HD11	1:B:412:VAL:HG21	1.80	0.63
1:A:440:ARG:O	1:A:444:GLU:HG3	1.99	0.62
1:B:475:PHE:HD1	1:B:543:MSE:HE1	1.65	0.62
1:A:477:MSE:HG3	1:A:532:LEU:HD23	1.84	0.60
1:A:488:HIS:O	1:A:490:LYS:HG3	2.01	0.60
1:B:400:PHE:HE2	1:B:409:LEU:HD11	1.66	0.60
1:B:377:GLY:O	1:B:383:LEU:HD11	2.02	0.59
1:A:517:MSE:HE2	3:A:719:HOH:O	2.01	0.59
1:A:405:THR:O	1:A:408:LEU:HB2	2.02	0.59
1:B:475:PHE:HD1	1:B:543:MSE:CE	2.16	0.58
1:A:398:GLU:C	1:A:400:PHE:H	2.07	0.58
1:A:374:GLU:OE2	1:A:482:ARG:HD2	2.04	0.58
1:B:432:PRO:HB3	1:B:506:ALA:HB3	1.86	0.57
1:A:544:MSE:HE1	1:B:536:LEU:HD21	1.86	0.57
1:B:398:GLU:HG2	1:B:399:LYS:N	2.19	0.57
1:A:357:GLN:HA	1:A:361:MSE:HE3	1.87	0.56
1:B:470:ARG:HD2	1:B:542:PHE:CE1	2.40	0.56
1:B:492:ASP:H	1:B:495:ASN:HD22	1.53	0.56
1:B:512:PRO:HG2	1:B:517:MSE:HE2	1.88	0.56
1:B:415:ILE:HD12	1:B:418:ILE:HD12	1.88	0.56
1:A:372:GLU:HG3	1:A:409:LEU:HB3	1.89	0.55
1:B:431:GLU:HB2	1:B:432:PRO:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:420:SER:O	1:B:424:ASP:HB2	2.08	0.54
1:A:507:HIS:CG	1:A:512:PRO:HG2	2.42	0.54
1:A:385:ARG:HB2	1:A:491:MSE:HE1	1.89	0.54
1:B:443:MSE:CE	1:B:523:ARG:HB2	2.35	0.53
1:B:357:GLN:HG3	3:B:739:HOH:O	2.08	0.52
1:B:399:LYS:HZ1	1:B:411:LYS:HZ3	1.56	0.52
1:B:439:ASN:O	1:B:443:MSE:HG3	2.09	0.52
1:A:477:MSE:HE2	1:A:480:LEU:HD12	1.92	0.52
1:A:390:ASP:O	1:A:394:LYS:HG3	2.09	0.52
1:B:522:LYS:HD3	3:B:744:HOH:O	2.10	0.52
1:B:545:VAL:O	1:B:546:GLU:CB	2.58	0.51
1:B:359:SER:HB2	1:B:546:GLU:OE2	2.11	0.51
1:B:466:PRO:HB2	1:B:469:ASN:OD1	2.11	0.50
1:B:458:MSE:O	1:B:462:VAL:HG23	2.12	0.50
1:A:434:LEU:O	1:A:435:THR:OG1	2.28	0.49
1:B:372:GLU:HG3	1:B:409:LEU:CB	2.40	0.49
1:B:499:VAL:O	1:B:502:PRO:HD2	2.11	0.49
1:B:477:MSE:O	1:B:481:GLN:HG3	2.13	0.49
1:A:395:GLU:O	1:A:398:GLU:N	2.45	0.49
1:A:442:PHE:CE1	1:A:458:MSE:HE1	2.47	0.49
1:A:442:PHE:CD1	1:A:458:MSE:HE1	2.47	0.49
1:B:538:TYR:O	1:B:541:GLN:HG2	2.12	0.49
1:B:470:ARG:HD2	1:B:542:PHE:CZ	2.48	0.48
1:A:443:MSE:HE3	1:A:523:ARG:NH1	2.28	0.48
1:B:427:ARG:NH1	1:B:502:PRO:O	2.47	0.48
1:A:440:ARG:HA	1:A:443:MSE:HE2	1.95	0.47
1:B:545:VAL:O	1:B:545:VAL:HG12	2.15	0.47
1:A:440:ARG:CZ	1:A:443:MSE:HE1	2.45	0.47
1:B:399:LYS:NZ	1:B:411:LYS:NZ	2.62	0.47
1:A:546:GLU:OXT	1:A:546:GLU:HG2	2.15	0.46
1:A:380:GLU:O	1:A:489:THR:HA	2.15	0.46
1:A:443:MSE:HE3	1:A:523:ARG:CZ	2.46	0.46
1:A:477:MSE:O	1:A:481:GLN:HG3	2.15	0.46
1:A:481:GLN:O	1:A:485:GLN:HG3	2.15	0.46
1:B:524:GLN:HB2	1:B:525:PRO:HD3	1.98	0.46
1:B:383:LEU:O	1:B:384:TYR:HB2	2.16	0.46
1:A:484:ALA:HA	1:A:491:MSE:O	2.17	0.45
1:B:399:LYS:HG2	1:B:407:PRO:HG3	1.98	0.45
1:B:399:LYS:O	1:B:407:PRO:HG3	2.16	0.45
1:B:407:PRO:O	1:B:409:LEU:N	2.50	0.45
1:A:383:LEU:O	1:A:384:TYR:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:MSE:CE	1:A:477:MSE:HA	2.46	0.44
1:A:450:ASP:HB3	1:A:453:ASN:HB2	1.98	0.44
1:A:372:GLU:HG3	1:A:409:LEU:CB	2.48	0.43
1:B:433:LEU:HD21	1:B:473:LEU:HA	1.99	0.43
1:B:377:GLY:C	1:B:383:LEU:HD11	2.39	0.43
1:B:458:MSE:HG3	1:B:530:ARG:HB3	2.00	0.43
1:A:517:MSE:O	1:A:521:ILE:HG13	2.19	0.43
1:A:470:ARG:CD	1:A:542:PHE:CE1	3.02	0.42
1:B:459:TYR:CZ	1:B:534:LEU:HD22	2.54	0.42
1:A:477:MSE:HA	1:A:477:MSE:HE3	2.02	0.42
1:B:350:MSE:SE	1:B:430:LYS:HD3	2.69	0.42
1:B:474:ALA:HA	1:B:539:TRP:CE3	2.55	0.42
1:A:432:PRO:HB3	1:A:506:ALA:HB3	2.02	0.42
1:A:480:LEU:HD22	1:A:496:LEU:HD22	2.01	0.42
1:B:399:LYS:HZ1	1:B:411:LYS:NZ	2.16	0.41
1:B:403:VAL:O	1:B:403:VAL:HG12	2.21	0.41
1:A:398:GLU:O	1:A:400:PHE:N	2.53	0.41
1:A:439:ASN:OD1	1:A:440:ARG:NH1	2.54	0.41
1:B:376:ARG:HG3	1:B:376:ARG:HH11	1.85	0.41
1:B:399:LYS:NZ	1:B:411:LYS:HZ3	2.18	0.41
1:A:544:MSE:HG3	3:A:738:HOH:O	2.19	0.41
1:B:402:ARG:HG3	1:B:403:VAL:N	2.35	0.40
1:B:524:GLN:N	1:B:525:PRO:CD	2.84	0.40
1:B:439:ASN:HB3	1:B:508:ALA:HB2	2.02	0.40
1:A:403:VAL:HG13	1:A:408:LEU:HG	2.02	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	199/208 (96%)	181 (91%)	15 (8%)	3 (2%)	10 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	199/208 (96%)	184 (92%)	13 (6%)	2 (1%)	15	37
All	All	398/416 (96%)	365 (92%)	28 (7%)	5 (1%)	12	30

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	408	LEU
1	B	408	LEU
1	A	399	LYS
1	B	545	VAL
1	A	356	SER

### 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	179/174 (103%)	176 (98%)	3 (2%)	60	84
1	B	179/174 (103%)	172 (96%)	7 (4%)	32	61
All	All	358/348 (103%)	348 (97%)	10 (3%)	43	73

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	357	GLN
1	A	460	GLN
1	A	477	MSE
1	B	376	ARG
1	B	398	GLU
1	B	400	PHE
1	B	402	ARG
1	B	413	ASP
1	B	424	ASP
1	B	460	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	357	GLN
1	A	460	GLN
1	A	495	ASN
1	A	511	ASN
1	A	524	GLN
1	B	371	ASN
1	B	375	GLN
1	B	453	ASN
1	B	460	GLN
1	B	495	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

2 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	SO4	B	601	-	4,4,4	0.39	0	6,6,6	0.08	0
2	SO4	A	600	-	4,4,4	0.38	0	6,6,6	0.12	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	192/208 (92%)	-0.13	10 (5%) 27 25	26, 41, 91, 107	0
1	B	192/208 (92%)	-0.05	8 (4%) 36 35	28, 48, 89, 106	0
All	All	384/416 (92%)	-0.09	18 (4%) 31 30	26, 45, 91, 107	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	348	GLU	4.8
1	A	346	ILE	4.8
1	B	347	GLY	4.1
1	A	546	GLU	3.8
1	B	403	VAL	3.7
1	B	402	ARG	3.7
1	A	347	GLY	3.6
1	A	403	VAL	3.2
1	A	406	VAL	3.1
1	A	402	ARG	3.1
1	B	346	ILE	3.0
1	B	546	GLU	2.7
1	A	404	LYS	2.7
1	A	405	THR	2.7
1	B	400	PHE	2.6
1	B	406	VAL	2.6
1	B	404	LYS	2.2
1	A	487	PRO	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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( $\text{\AA}^2$ )	Q<0.9
2	SO4	B	601	5/5	0.84	0.23	118,119,119,120	0
2	SO4	A	600	5/5	0.95	0.12	105,105,106,106	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.