



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 09:39 pm BST

PDB ID : 3PMI
Title : PWWP Domain of Human Mutated Melanoma-Associated Antigen 1
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Edwards, A.M.; Bochkarev, A.; Min, J.; Wu, H.; Structural Genomics Consor-
tium (SGC)
Deposited on : 2010-11-17
Resolution : 2.82 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 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

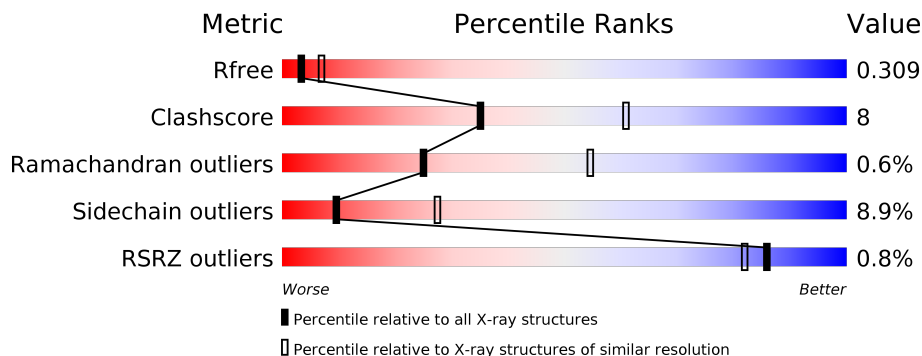
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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 ≥ 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	134	
1	B	134	
1	C	134	
1	D	134	

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
3	PEG	A	10	-	-	-	X

2 Entry composition [i](#)

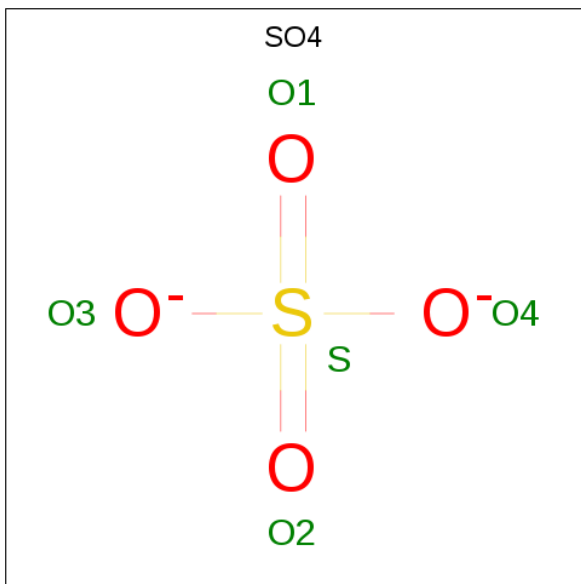
There are 4 unique types of molecules in this entry. The entry contains 3838 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PWWP domain-containing protein MUM1.

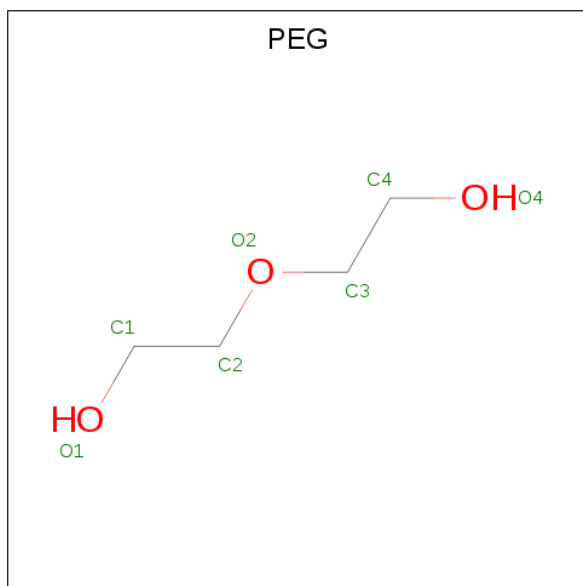
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	122	Total 948	C 613	N 158	O 171	S 3	Se 3	0	0	0
1	B	122	Total 967	C 624	N 165	O 172	S 3	Se 3	0	0	0
1	C	121	Total 947	C 615	N 158	O 168	S 3	Se 3	0	0	0
1	D	120	Total 935	C 602	N 158	O 169	S 3	Se 3	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 4 is UNKNOWN LIGAND (three-letter code: UNL) (formula:).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total	C	N	O	0	0
			12	6	4	2		
4	C	1	Total	C	N	O	0	0
			12	6	4	2		

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: PWWP domain-containing protein MUM1

Chain A: 



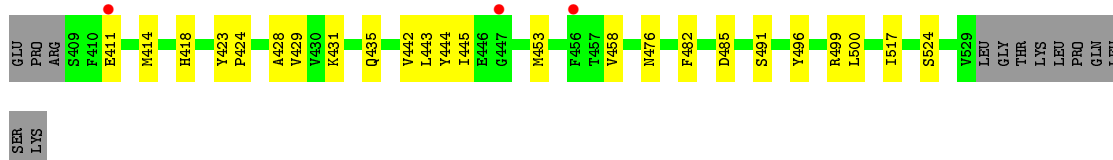
- Molecule 1: PWWP domain-containing protein MUM1

Chain B: 




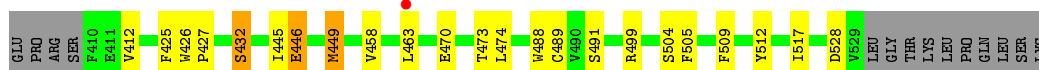
- Molecule 1: PWWP domain-containing protein MUM1

Chain C: 



- Molecule 1: PWWP domain-containing protein MUM1

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.04Å 57.72Å 187.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.82 47.44 – 2.82	Depositor EDS
% Data completeness (in resolution range)	99.2 (50.00-2.82) 99.2 (47.44-2.82)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.81Å)	Xtrriage
Refinement program	REFMAC refmac_5.6.0081	Depositor
R, R_{free}	0.228 , 0.304 0.228 , 0.309	Depositor DCC
R_{free} test set	751 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	57.8	Xtrriage
Anisotropy	0.539	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.027 for k,h,-l	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3838	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, UNL, 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	0.59	0/969	0.64	0/1310
1	B	0.55	0/990	0.60	0/1336
1	C	0.56	0/970	0.64	0/1309
1	D	0.54	0/957	0.58	0/1295
All	All	0.56	0/3886	0.62	0/5250

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 [i](#)

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	948	0	894	26	0
1	B	967	0	923	13	0
1	C	947	0	902	10	0
1	D	935	0	875	13	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	7	0	10	0	0
4	C	12	0	0	1	0
4	D	12	0	0	0	0
All	All	3838	0	3604	61	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 8.

All (61) 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:518:SER:HB2	1:A:522:ARG:HE	1.21	1.01
1:A:527:GLN:O	1:A:530:LEU:HB2	1.70	0.92
1:A:411:GLU:H	1:A:414:MSE:SE	2.12	0.83
1:B:454:LYS:HE3	1:B:455:GLY:H	1.45	0.81
1:B:411:GLU:HB2	1:B:414:MSE:HE3	1.61	0.81
1:C:499:ARG:HD3	1:C:517:ILE:HD12	1.63	0.80
1:A:518:SER:HB2	1:A:522:ARG:NE	1.97	0.79
1:D:499:ARG:HG2	1:D:499:ARG:HH11	1.51	0.75
1:A:454:LYS:HE2	1:A:454:LYS:H	1.57	0.70
1:A:512:TYR:O	1:A:518:SER:OG	2.11	0.68
1:A:445:ILE:HG23	1:A:449:MSE:HE1	1.75	0.67
1:D:427:PRO:HG2	1:D:489:CYS:SG	2.40	0.62
1:D:505:PHE:CD1	1:D:517:ILE:HD11	2.35	0.62
4:C:20:UNL:NH1	4:C:20:UNL:CG	2.63	0.62
1:D:488:TRP:HZ2	1:D:528:ASP:HB3	1.65	0.61
1:A:518:SER:CB	1:A:522:ARG:HE	2.07	0.60
1:B:475:LEU:O	1:B:479:ARG:HG3	2.03	0.59
1:A:479:ARG:HG2	1:A:486:ILE:HG21	1.86	0.57
1:A:529:VAL:O	1:A:530:LEU:HD13	2.05	0.55
1:C:428:ALA:HB1	1:C:442:VAL:HG13	1.87	0.55
1:D:499:ARG:NH1	1:D:499:ARG:HG2	2.20	0.55
1:B:411:GLU:H	1:B:414:MSE:SE	2.40	0.54
1:D:509:PHE:O	1:D:512:TYR:HB3	2.09	0.53
1:A:445:ILE:HG23	1:A:449:MSE:CE	2.39	0.52
1:C:443:LEU:HD13	1:C:453:MSE:HE2	1.92	0.52
1:A:482:PHE:HE1	1:B:453:MSE:HE3	1.75	0.51
1:A:488:TRP:CD1	1:A:525:ILE:HG12	2.46	0.51
1:D:426:TRP:CD1	1:D:427:PRO:HD2	2.45	0.50
1:D:505:PHE:CD1	1:D:517:ILE:CD1	2.94	0.50
1:B:529:VAL:O	1:B:530:LEU:HD23	2.12	0.49
1:A:505:PHE:CD1	1:A:517:ILE:CD1	2.96	0.48
1:C:482:PHE:HB3	1:C:485:ASP:HB2	1.94	0.48
1:A:528:ASP:C	1:A:530:LEU:H	2.17	0.47
1:D:445:ILE:HG23	1:D:449:MSE:HE1	1.97	0.46
1:A:505:PHE:CD1	1:A:517:ILE:HD11	2.50	0.46
1:A:495:ASP:HB3	1:A:512:TYR:OH	2.16	0.46
1:A:446:GLU:H	1:A:449:MSE:HE2	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:LEU:HD22	1:A:445:ILE:HD11	1.98	0.45
1:B:442:VAL:H	1:B:456:PHE:H	1.63	0.45
1:A:500:LEU:HD12	1:A:500:LEU:HA	1.85	0.44
1:B:496:TYR:O	1:B:499:ARG:HB2	2.18	0.44
1:C:418:HIS:HD2	1:C:444:TYR:OH	2.00	0.44
1:A:475:LEU:O	1:A:479:ARG:HG3	2.18	0.43
1:D:446:GLU:HG2	1:D:449:MSE:HB3	1.99	0.43
1:B:519:TYR:HB3	1:B:520:PRO:HD3	1.99	0.43
1:D:425:PHE:HB2	1:D:474:LEU:HD22	1.99	0.43
1:B:479:ARG:HG2	1:B:486:ILE:HG21	2.00	0.43
1:D:412:VAL:HG22	1:D:432:SER:HA	2.00	0.42
1:A:526:GLN:NE2	1:B:529:VAL:HB	2.34	0.42
1:C:411:GLU:H	1:C:414:MSE:SE	2.53	0.42
1:A:479:ARG:HG2	1:A:486:ILE:CG2	2.49	0.42
1:C:429:VAL:HB	1:C:445:ILE:HD11	2.02	0.42
1:A:417:TRP:CD1	1:A:466:PHE:HA	2.55	0.41
1:A:486:ILE:HD12	1:A:486:ILE:HA	1.91	0.41
1:B:454:LYS:HE3	1:B:455:GLY:N	2.23	0.41
1:C:496:TYR:CZ	1:C:500:LEU:HD22	2.55	0.41
1:C:423:TYR:HB3	1:C:424:PRO:HD2	2.03	0.41
1:A:439:LYS:HA	1:A:459:SER:HA	2.02	0.41
1:D:458:VAL:HG21	1:D:463:LEU:HD21	2.03	0.41
1:C:418:HIS:CE1	1:C:458:VAL:HG11	2.57	0.40
1:B:493:ILE:O	1:B:497:ARG:HG3	2.22	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	120/134 (90%)	112 (93%)	7 (6%)	1 (1%)	19 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	120/134 (90%)	114 (95%)	4 (3%)	2 (2%)	9	27
1	C	119/134 (89%)	111 (93%)	8 (7%)	0	100	100
1	D	118/134 (88%)	109 (92%)	9 (8%)	0	100	100
All	All	477/536 (89%)	446 (94%)	28 (6%)	3 (1%)	25	54

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	529	VAL
1	B	455	GLY
1	B	529	VAL

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	97/117 (83%)	86 (89%)	11 (11%)	6	17
1	B	101/117 (86%)	89 (88%)	12 (12%)	5	15
1	C	98/117 (84%)	93 (95%)	5 (5%)	24	54
1	D	96/117 (82%)	89 (93%)	7 (7%)	14	37
All	All	392/468 (84%)	357 (91%)	35 (9%)	9	27

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	411	GLU
1	A	442	VAL
1	A	449	MSE
1	A	454	LYS
1	A	477	GLN
1	A	500	LEU
1	A	504	SER
1	A	518	SER

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Mol	Chain	Res	Type
1	A	523	LYS
1	A	527	GLN
1	A	530	LEU
1	B	409	SER
1	B	431	LYS
1	B	432	SER
1	B	436	ARG
1	B	442	VAL
1	B	454	LYS
1	B	463	LEU
1	B	485	ASP
1	B	491	SER
1	B	499	ARG
1	B	504	SER
1	B	524	SER
1	C	431	LYS
1	C	435	GLN
1	C	476	ASN
1	C	491	SER
1	C	524	SER
1	D	432	SER
1	D	446	GLU
1	D	449	MSE
1	D	470	GLU
1	D	473	THR
1	D	491	SER
1	D	504	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	526	GLN
1	A	527	GLN
1	B	418	HIS
1	C	418	HIS
1	C	420	HIS
1	C	472	GLN
1	C	476	ASN
1	D	476	ASN
1	D	484	GLN

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](#)

Of 5 ligands modelled in this entry, 2 are unknown - leaving 3 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	$\# Z > 2$
2	SO4	A	2	-	4,4,4	0.38	0	6,6,6	0.14	0
2	SO4	B	1	-	4,4,4	0.38	0	6,6,6	0.16	0
3	PEG	A	10	-	6,6,6	0.64	0	5,5,5	0.50	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	10	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	10	PEG	O2-C3-C4-O4
3	A	10	PEG	C1-C2-O2-C3

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, 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	119/134 (88%)	-0.13	0	100 100	27, 49, 86, 104	0
1	B	119/134 (88%)	-0.20	0	100 100	26, 52, 91, 113	0
1	C	118/134 (88%)	-0.14	3 (2%)	57 47	34, 51, 79, 100	0
1	D	117/134 (87%)	0.09	1 (0%)	84 80	33, 54, 102, 141	0
All	All	473/536 (88%)	-0.10	4 (0%)	86 82	26, 52, 91, 141	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	447	GLY	2.8
1	D	463	LEU	2.5
1	C	456	PHE	2.2
1	C	411	GLU	2.1

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. 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(\AA^2)	Q<0.9
3	PEG	A	10	7/7	0.69	0.44	84,93,106,108	0
4	UNL	D	21	12/-	0.83	0.27	76,85,88,89	0
4	UNL	C	20	12/-	0.86	0.26	77,89,103,104	0
2	SO4	A	2	5/5	0.93	0.18	99,103,110,110	0
2	SO4	B	1	5/5	0.94	0.23	101,104,108,109	0

6.5 Other polymers [i](#)

There are no such residues in this entry.