Full wwPDB X-ray Structure Validation Report (i)

Apr 28, 2024 – 03:19 pm BST

PDB ID	:	2WAL
Title	:	Crystal Structure of human GADD45gamma
Authors	:	Bhattacharya, S.; Mueller, J.J.; Roske, Y.; Turnbull, A.P.; Quedenau, C.; Goetz, F.; Buessow, K.; Heinemann, U.
Deposited on	:	2009-02-09
Resolution	:	2.40 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

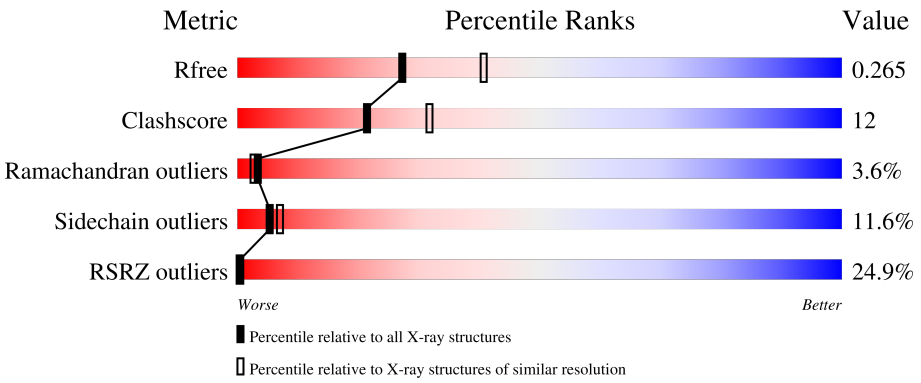
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 >=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 <=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	159	21% 69% 19% 5% 6%
1	B	159	23% 60% 21% • • 14%

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:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MLA	A	1160	-	X	-	-
2	MLA	A	1161	-	X	-	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2210 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.

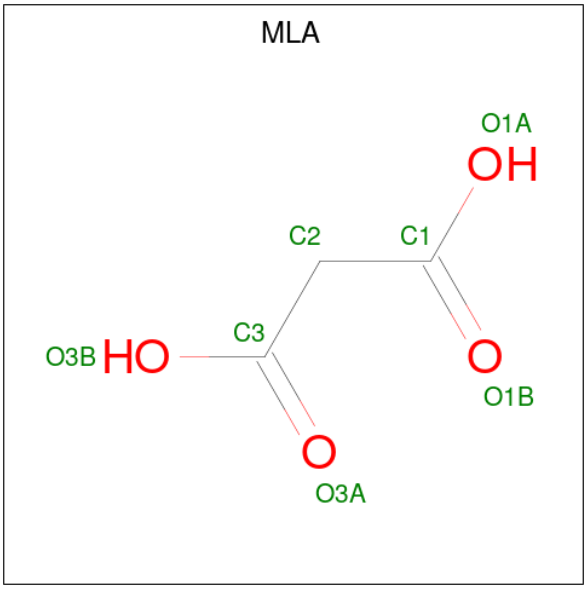
- Molecule 1 is a protein called GROWTH ARREST AND DNA-DAMAGE-INDUCIBLE PROTEIN GADD45 GAMMA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	149	1115	695	188	225	7	0	0	0
1	B	136	1027	645	174	201	7	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	58	ALA	THR	engineered mutation	UNP O95257
B	58	ALA	THR	engineered mutation	UNP O95257

- Molecule 2 is MALONIC ACID (three-letter code: MLA) (formula: $C_3H_4O_4$).



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

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

- Molecule 3 is water.

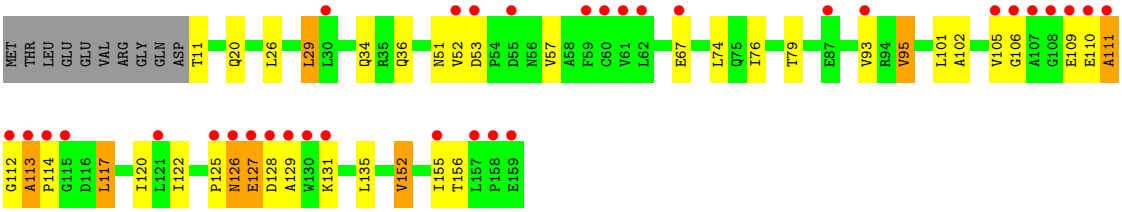
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	33	33	33	0	0
			Total	O		
3	B	21	21	21	0	0

3 Residue-property plots (i)

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

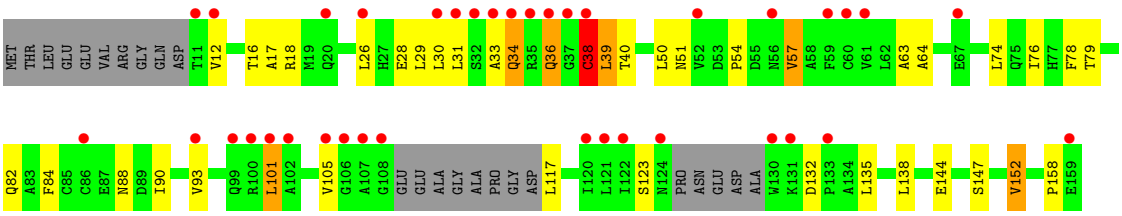
• Molecule 1: GROWTH ARREST AND DNA-DAMAGE-INDUCIBLE PROTEIN GADD45 GAMMA

Chain A:



• Molecule 1: GROWTH ARREST AND DNA-DAMAGE-INDUCIBLE PROTEIN GADD45 GAMMA

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	108.72Å 108.72Å 73.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.81 – 2.40 19.78 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.81-2.40) 99.5 (19.78-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.68 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.215 , 0.251 0.231 , 0.265	Depositor DCC
R_{free} test set	988 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	54.6	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 82.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	2210	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MLA

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	0.55	0/1133	0.68	0/1543
1	B	0.50	0/1041	0.63	0/1413
All	All	0.53	0/2174	0.66	0/2956

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
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	127	GLU	Peptide
1	B	34	GLN	Peptide

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	1115	0	1083	27	0
1	B	1027	0	1014	28	0
2	A	14	0	4	1	0
3	A	33	0	0	0	0
3	B	21	0	0	0	0
All	All	2210	0	2101	51	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 12.

All (51) 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:105:VAL:HG21	1:A:117:LEU:HD21	1.41	1.02
1:A:29:LEU:HD13	1:A:122:ILE:HD13	1.43	1.00
1:A:117:LEU:HD22	1:A:120:ILE:HD11	1.60	0.83
1:A:117:LEU:HD12	1:A:117:LEU:H	1.57	0.69
1:A:105:VAL:CG2	1:A:117:LEU:HD21	2.21	0.67
1:B:16:THR:HG22	1:B:18:ARG:H	1.61	0.65
1:A:76:ILE:HD12	1:B:76:ILE:HD12	1.79	0.64
1:A:93:VAL:HG21	1:A:155:ILE:HD12	1.82	0.62
1:B:34:GLN:HE22	1:B:105:VAL:HG23	1.65	0.62
1:A:105:VAL:HG21	1:A:117:LEU:CD2	2.23	0.61
1:A:79:THR:HG1	1:B:79:THR:HG1	1.41	0.60
1:A:79:THR:HG21	1:B:76:ILE:HD13	1.84	0.57
1:A:102:ALA:CA	1:A:117:LEU:HD11	2.35	0.57
1:B:50:LEU:HD22	1:B:84:PHE:CE2	2.43	0.54
1:B:39:LEU:O	1:B:39:LEU:HD23	2.08	0.53
1:A:152:VAL:HG22	1:A:152:VAL:O	2.09	0.52
1:B:50:LEU:HD22	1:B:84:PHE:CD2	2.44	0.52
1:A:112:GLY:O	1:A:113:ALA:HB3	2.10	0.51
1:B:64:ALA:HB2	1:B:78:PHE:HE2	1.76	0.51
1:B:57:VAL:HA	1:B:123:SER:HA	1.93	0.50
1:A:102:ALA:HA	1:A:117:LEU:HD11	1.94	0.49
1:A:110:GLU:O	1:A:111:ALA:HB2	2.12	0.49
1:B:30:LEU:O	1:B:33:ALA:O	2.31	0.48
1:A:95:VAL:HG13	2:A:1160:MLA:O3B	2.14	0.48
1:B:64:ALA:HB2	1:B:78:PHE:CE2	2.50	0.47
1:B:16:THR:HG22	1:B:17:ALA:N	2.30	0.47
1:A:36:GLN:NE2	1:A:129:ALA:HB1	2.30	0.46
1:A:127:GLU:HB3	1:A:129:ALA:HB2	1.96	0.46
1:B:34:GLN:NE2	1:B:105:VAL:HG23	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:THR:HG22	1:B:18:ARG:N	2.27	0.46
1:B:26:LEU:HA	1:B:138:LEU:HD11	1.98	0.45
1:B:54:PRO:HA	1:B:57:VAL:HG12	1.98	0.45
1:A:127:GLU:O	1:A:128:ASP:HB2	2.16	0.45
1:B:152:VAL:O	1:B:152:VAL:HG22	2.19	0.43
1:B:33:ALA:O	1:B:34:GLN:CB	2.67	0.43
1:A:125:PRO:O	1:A:126:ASN:C	2.56	0.43
1:B:82:GLN:NE2	1:B:152:VAL:HG11	2.34	0.42
1:B:88:ASN:O	1:B:90:ILE:HD12	2.19	0.42
1:A:101:LEU:HD23	1:A:117:LEU:HD13	2.00	0.42
1:B:39:LEU:O	1:B:40:THR:C	2.58	0.42
1:B:63:ALA:CB	1:B:101:LEU:HD12	2.50	0.42
1:B:36:GLN:HB3	1:B:38:CYS:SG	2.60	0.41
1:B:33:ALA:O	1:B:34:GLN:HB3	2.21	0.41
1:A:79:THR:OG1	1:B:79:THR:OG1	2.21	0.41
1:A:52:VAL:HG12	1:A:53:ASP:N	2.36	0.41
1:A:102:ALA:N	1:A:117:LEU:HD11	2.36	0.41
1:B:54:PRO:HA	1:B:57:VAL:CG1	2.52	0.40
1:A:34:GLN:OE1	1:A:106:GLY:N	2.54	0.40
1:A:36:GLN:HE21	1:A:129:ALA:HB1	1.86	0.40
1:A:117:LEU:CD2	1:A:120:ILE:HD11	2.41	0.40
1:B:144:GLU:O	1:B:147:SER:HB2	2.21	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	147/159 (92%)	136 (92%)	6 (4%)	5 (3%)	3 3
1	B	130/159 (82%)	119 (92%)	6 (5%)	5 (4%)	3 2
All	All	277/318 (87%)	255 (92%)	12 (4%)	10 (4%)	3 3

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	38	CYS
1	A	126	ASN
1	B	132	ASP
1	A	111	ALA
1	A	114	PRO
1	B	36	GLN
1	A	109	GLU
1	B	158	PRO
1	A	113	ALA
1	B	12	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	120/129 (93%)	106 (88%)	14 (12%)	5 7
1	B	112/129 (87%)	99 (88%)	13 (12%)	5 7
All	All	232/258 (90%)	205 (88%)	27 (12%)	5 7

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	THR
1	A	20	GLN
1	A	26	LEU
1	A	29	LEU
1	A	51	ASN
1	A	57	VAL
1	A	67	GLU
1	A	74	LEU
1	A	95	VAL
1	A	117	LEU
1	A	131	LYS
1	A	135	LEU
1	A	152	VAL

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Mol	Chain	Res	Type
1	A	156	THR
1	B	28	GLU
1	B	29	LEU
1	B	31	LEU
1	B	38	CYS
1	B	39	LEU
1	B	51	ASN
1	B	57	VAL
1	B	74	LEU
1	B	93	VAL
1	B	101	LEU
1	B	117	LEU
1	B	135	LEU
1	B	152	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	36	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 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	MLA	A	1160	-	6,6,6	1.97	2 (33%)	7,7,7	1.47	2 (28%)
2	MLA	A	1161	-	6,6,6	2.01	3 (50%)	7,7,7	1.01	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	MLA	A	1160	-	-	4/4/4/4	-
2	MLA	A	1161	-	-	4/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1160	MLA	C2-C3	3.39	1.56	1.51
2	A	1161	MLA	C2-C3	3.23	1.56	1.51
2	A	1161	MLA	C2-C1	2.40	1.54	1.51
2	A	1161	MLA	O1A-C1	-2.03	1.23	1.30
2	A	1160	MLA	O3B-C3	-2.02	1.23	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1160	MLA	O3A-C3-C2	-2.19	115.69	122.08
2	A	1160	MLA	O3B-C3-C2	2.12	121.30	114.54

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1161	MLA	O1A-C1-C2-C3
2	A	1161	MLA	O1B-C1-C2-C3
2	A	1161	MLA	C1-C2-C3-O3B
2	A	1161	MLA	C1-C2-C3-O3A

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Mol	Chain	Res	Type	Atoms
2	A	1160	MLA	O1B-C1-C2-C3
2	A	1160	MLA	O1A-C1-C2-C3
2	A	1160	MLA	C1-C2-C3-O3A
2	A	1160	MLA	C1-C2-C3-O3B

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1160	MLA	1	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, 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	149/159 (93%)	1.56	34 (22%)	0	0	47, 56, 80, 92	0
1	B	136/159 (85%)	1.51	37 (27%)	0	0	40, 56, 74, 90	0
All	All	285/318 (89%)	1.54	71 (24%)	0	0	40, 56, 78, 92	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	111	ALA	11.6
1	A	113	ALA	11.4
1	A	128	ASP	11.3
1	A	108	GLY	10.6
1	A	109	GLU	9.8
1	B	106	GLY	8.9
1	B	107	ALA	8.5
1	A	114	PRO	8.3
1	A	127	GLU	7.8
1	B	11	THR	7.4
1	B	124	ASN	7.2
1	B	108	GLY	7.0
1	A	110	GLU	6.9
1	A	112	GLY	6.7
1	A	126	ASN	6.4
1	A	107	ALA	6.3
1	A	52	VAL	5.8
1	A	159	GLU	5.6
1	B	130	TRP	5.5
1	B	131	LYS	5.3
1	A	105	VAL	5.1
1	A	129	ALA	4.9
1	B	12	VAL	4.9
1	B	159	GLU	4.9

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Mol	Chain	Res	Type	RSRZ
1	B	101	LEU	4.8
1	A	106	GLY	4.8
1	A	130	TRP	4.4
1	B	133	PRO	4.4
1	B	34	GLN	4.1
1	A	55	ASP	4.1
1	B	30	LEU	4.0
1	B	99	GLN	4.0
1	B	122	ILE	3.7
1	B	36	GLN	3.5
1	A	131	LYS	3.3
1	A	115	GLY	3.3
1	A	125	PRO	3.2
1	B	37	GLY	3.2
1	A	61	VAL	3.1
1	B	38	CYS	3.1
1	B	59	PHE	3.1
1	A	53	ASP	3.0
1	B	61	VAL	3.0
1	A	60	CYS	2.9
1	B	105	VAL	2.9
1	B	93	VAL	2.8
1	A	67	GLU	2.7
1	A	158	PRO	2.7
1	A	30	LEU	2.7
1	B	26	LEU	2.7
1	B	100	ARG	2.7
1	B	67	GLU	2.7
1	B	31	LEU	2.7
1	B	121	LEU	2.6
1	A	155	ILE	2.6
1	A	93	VAL	2.6
1	B	60	CYS	2.5
1	B	120	ILE	2.4
1	B	32	SER	2.4
1	B	35	ARG	2.4
1	B	86	CYS	2.3
1	B	33	ALA	2.3
1	B	52	VAL	2.3
1	A	59	PHE	2.3
1	A	62	LEU	2.2
1	A	87	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	56	ASN	2.1
1	A	121	LEU	2.1
1	B	102	ALA	2.1
1	A	157	LEU	2.0
1	B	20	GLN	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, 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(Å^2)	Q<0.9
2	MLA	A	1161	7/7	0.86	0.37	80,81,84,86	7
2	MLA	A	1160	7/7	0.88	0.26	54,61,68,72	7

6.5 Other polymers (i)

There are no such residues in this entry.