



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 07:55 am BST

PDB ID : 5MFG
Title : Designed armadillo repeat protein YIIIM5AII in complex with peptide (RR)₄
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Deposited on : 2016-11-18
Resolution : 1.90 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
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

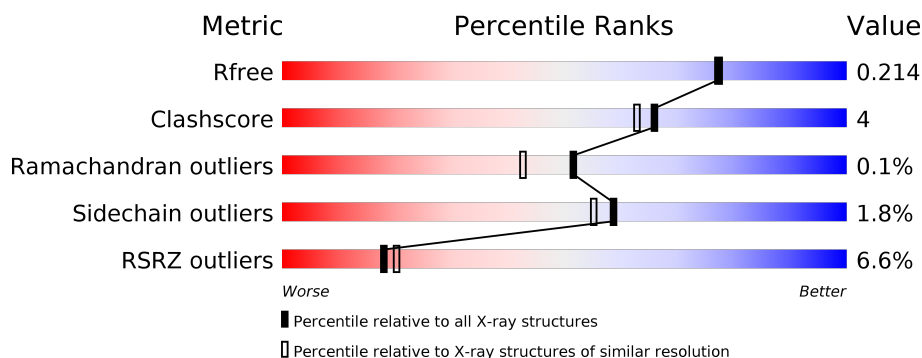
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	286	<div> <div>7%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div></div> </div> </div>
1	B	286	<div> <div>7%</div> <div> <div></div> <div>86%</div> <div>8%</div> <div>6%</div> </div> </div>
1	C	286	<div> <div>8%</div> <div> <div></div> <div>89%</div> <div>5%</div> <div>6%</div> </div> </div>
1	D	286	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>8%</div> <div>16%</div> </div> </div>
2	E	10	<div> <div>10%</div> <div> <div></div> <div>20%</div> <div>60%</div> <div>20%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16563 atoms, of which 8066 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 YIIIM5AII.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	282	Total	C	H	N	O	S	25	3	0
			4230	1315	2121	360	433	1			
1	B	269	Total	C	H	N	O		23	1	0
			3998	1243	2005	340	410				
1	C	270	Total	C	H	N	O	S	24	2	0
			4054	1266	2033	345	409	1			
1	D	241	Total	C	H	N	O	S	21	1	0
			3539	1105	1775	298	360	1			

- Molecule 2 is a protein called (RR)4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	10	Total	C	H	N	O	0	0	0
			243	60	132	40	11			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	5	Total	Ca	0	0
			5	5		
3	A	7	Total	Ca	0	0
			7	7		
3	D	1	Total	Ca	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	116	Total	O	0	0
			116	116		
4	B	101	Total	O	0	0
			101	101		

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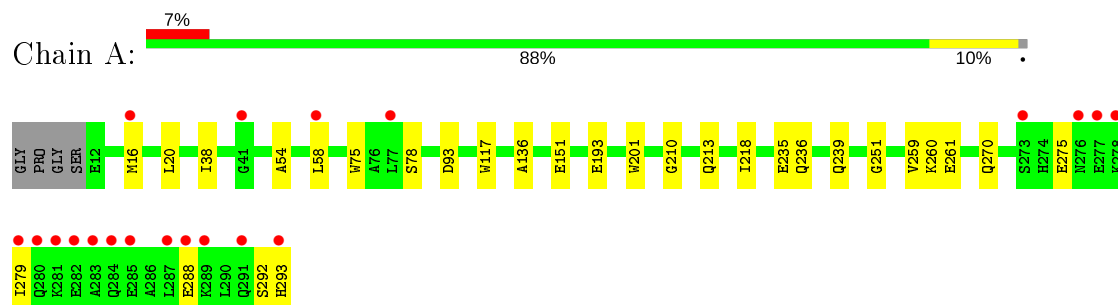
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	123	Total 123	O 123	0	0
4	D	130	Total 130	O 130	0	0
4	E	16	Total 16	O 16	0	0

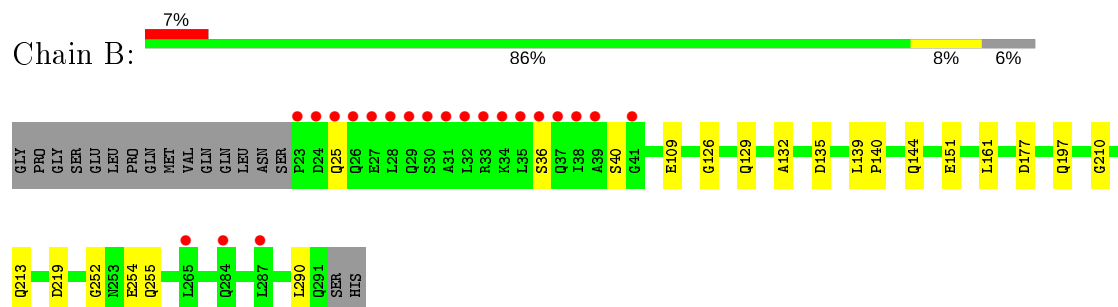
3 Residue-property plots [i](#)

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.

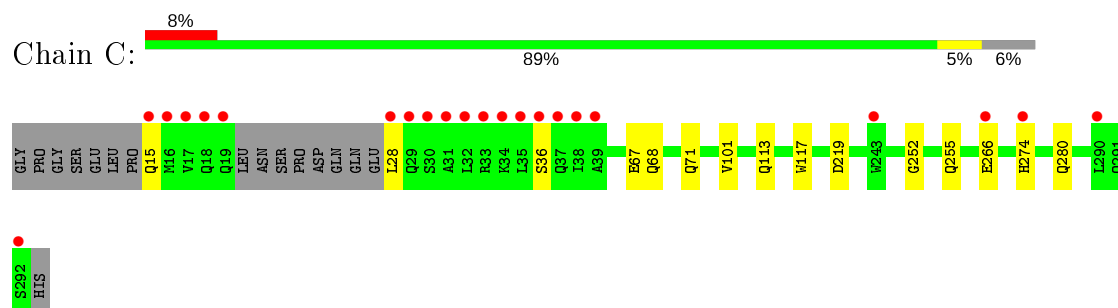
• Molecule 1: YIIIM5AII



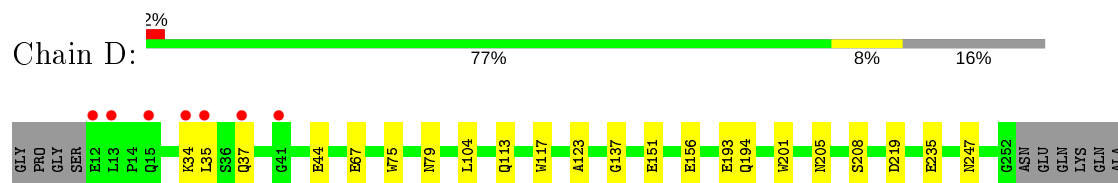
• Molecule 1: YIIIM5AII



• Molecule 1: YIIIM5AII

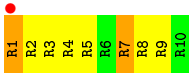


• Molecule 1: YIIIM5AII



VAL
LYS
GLU
ALA
GLY
ALA
LEU
GLU
LYS
LEU
GLN
GLN
GLN
SER
HIS
GLU
ASN
GLU
LYS
ILE
GLN
LYS
GLU
ALA
GLN
GLU
ALA
LEU
GLU
LYS
LEU
GLN
SER
HIS

● Molecule 2: (RR)4



4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α , β , γ	168.54 Å 168.54 Å 80.58 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.65 – 1.90 48.65 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.65-1.90) 100.0 (48.65-1.90)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 1.90 Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.182 , 0.213 0.183 , 0.214	Depositor DCC
R_{free} test set	5055 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 28.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.229 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	16563	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: CA

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	0.26	0/2138	0.38	0/2911
1	B	0.25	0/2014	0.37	0/2743
1	C	0.26	0/2044	0.38	0/2784
1	D	0.25	0/1788	0.37	0/2446
2	E	0.49	0/110	0.76	0/137
All	All	0.26	0/8094	0.38	0/11021

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
2	E	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	1	ARG	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	2109	2121	2122	25	2
1	B	1993	2005	2004	12	0
1	C	2021	2033	2030	8	2
1	D	1764	1775	1776	19	0
2	E	111	132	132	12	0
3	A	7	0	0	0	0
3	B	5	0	0	0	0
3	D	1	0	0	0	0
4	A	116	0	0	6	2
4	B	101	0	0	5	3
4	C	123	0	0	4	1
4	D	130	0	0	11	0
4	E	16	0	0	0	0
All	All	8497	8066	8064	64	5

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 (64) 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:D:34:LYS:NZ	1:D:37:GLN:OE1	1.96	0.98
1:D:113:GLN:O	4:D:401:HOH:O	1.90	0.89
1:B:219:ASP:OD2	4:B:401:HOH:O	1.95	0.85
1:D:156:GLU:OE1	4:D:401:HOH:O	1.96	0.82
1:D:117:TRP:N	4:D:401:HOH:O	2.15	0.79
1:D:205:ASN:OD1	2:E:9:ARG:NH2	2.15	0.78
1:C:101:VAL:O	4:C:401:HOH:O	2.01	0.78
1:A:270:GLN:NE2	1:D:156:GLU:OE2	2.16	0.78
1:A:236:GLN:OE1	4:A:401:HOH:O	2.04	0.75
1:D:219:ASP:OD2	4:D:402:HOH:O	2.04	0.74
1:C:274:HIS:O	1:C:280:GLN:NE2	2.24	0.70
1:B:177:ASP:OD2	4:B:402:HOH:O	2.08	0.70
1:A:275:GLU:O	1:A:279:ILE:N	2.26	0.68
1:A:93:ASP:OD1	4:A:402:HOH:O	2.12	0.68
1:A:193:GLU:OE2	4:A:403:HOH:O	2.12	0.68
1:B:135:ASP:OD2	4:B:403:HOH:O	2.11	0.68
1:B:132:ALA:O	4:B:404:HOH:O	2.12	0.67
1:D:208:SER:OG	4:D:404:HOH:O	2.13	0.67
1:D:67:GLU:OE2	4:D:403:HOH:O	2.12	0.67
1:D:123:ALA:O	4:D:405:HOH:O	2.15	0.65
1:C:219:ASP:OD2	4:C:402:HOH:O	2.15	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:75:TRP:O	1:D:79:ASN:ND2	2.31	0.62
1:B:25:GLN:NE2	4:B:408:HOH:O	2.31	0.58
1:A:235:GLU:OE2	1:A:235:GLU:N	2.36	0.58
1:D:104:LEU:O	4:D:406:HOH:O	2.18	0.56
1:B:210:GLY:N	1:B:213:GLN:OE1	2.38	0.53
1:D:235:GLU:OE2	1:D:235:GLU:N	2.35	0.52
1:D:247:ASN:OD1	2:E:7:ARG:NH2	2.31	0.52
1:A:201:TRP:CH2	2:E:8:ARG:HB3	2.45	0.52
1:A:218:ILE:O	4:A:404:HOH:O	2.19	0.52
1:B:126:GLY:N	1:B:129:GLN:OE1	2.43	0.52
1:D:137:GLY:O	4:D:407:HOH:O	2.19	0.51
1:C:117:TRP:CZ3	2:E:3:ARG:HB3	2.46	0.50
1:C:67:GLU:OE2	1:C:67:GLU:N	2.39	0.50
1:D:193:GLU:OE2	4:D:408:HOH:O	2.20	0.50
1:D:201:TRP:CH2	2:E:9:ARG:HB3	2.46	0.50
1:B:252:GLY:N	1:B:255:GLN:OE1	2.45	0.49
1:C:68:GLN:NE2	4:C:421:HOH:O	2.47	0.48
1:A:210:GLY:N	1:A:213:GLN:OE1	2.42	0.47
1:B:109:GLU:OE1	1:B:109:GLU:N	2.44	0.46
2:E:2:ARG:HD2	2:E:2:ARG:N	2.29	0.46
1:A:54:ALA:O	1:A:58:LEU:HD13	2.16	0.46
1:C:113:GLN:NE2	4:C:415:HOH:O	2.42	0.45
1:A:75:TRP:CZ3	2:E:2:ARG:HB3	2.52	0.44
1:A:193:GLU:N	1:A:193:GLU:OE1	2.44	0.44
1:A:20:LEU:HD13	1:A:58:LEU:CD1	2.47	0.44
1:D:194:GLN:NE2	4:D:420:HOH:O	2.45	0.44
1:A:117:TRP:CH2	2:E:4:ARG:HB3	2.52	0.44
1:B:139:LEU:N	1:B:140:PRO:HD2	2.33	0.44
1:C:252:GLY:N	1:C:255:GLN:OE1	2.47	0.43
1:A:259:VAL:O	1:A:261:GLU:N	2.51	0.43
1:A:75:TRP:CZ2	2:E:4:ARG:HD3	2.53	0.43
1:A:16:MET:HE1	1:A:38:ILE:HD12	2.01	0.43
1:A:75:TRP:CH2	2:E:4:ARG:HG2	2.54	0.43
1:A:136:ALA:O	4:A:405:HOH:O	2.22	0.42
1:A:78[B]:SER:HB2	2:E:4:ARG:HH22	1.85	0.42
1:A:292:SER:O	1:A:293:HIS:C	2.58	0.42
1:D:151:GLU:N	1:D:151:GLU:OE2	2.52	0.41
1:A:20:LEU:HD13	1:A:58:LEU:HD12	2.02	0.41
1:A:151:GLU:OE2	1:A:151:GLU:N	2.47	0.40
1:A:251:GLY:N	4:A:421:HOH:O	2.54	0.40
1:A:78[B]:SER:CB	2:E:4:ARG:HH22	2.34	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:GLU:OE2	1:B:151:GLU:N	2.53	0.40
1:B:36:SER:O	1:B:40:SER:OG	2.26	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:514:HOH:O	4:B:489:HOH:O[6_665]	1.93	0.27
1:A:288:GLU:OE1	1:C:36:SER:HG[1_554]	1.36	0.24
4:B:499:HOH:O	4:C:510:HOH:O[1_554]	2.07	0.13
4:A:492:HOH:O	4:B:489:HOH:O[6_665]	2.07	0.13
1:A:288:GLU:OE1	1:C:36:SER:OG[1_554]	2.10	0.10

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	283/286 (99%)	279 (99%)	3 (1%)	1 (0%)	34	24
1	B	268/286 (94%)	267 (100%)	1 (0%)	0	100	100
1	C	268/286 (94%)	264 (98%)	4 (2%)	0	100	100
1	D	240/286 (84%)	238 (99%)	2 (1%)	0	100	100
2	E	8/10 (80%)	8 (100%)	0	0	100	100
All	All	1067/1154 (92%)	1056 (99%)	10 (1%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	LYS

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	225/224 (100%)	224 (100%)	1 (0%)	91	91
1	B	210/224 (94%)	205 (98%)	5 (2%)	49	43
1	C	212/224 (95%)	208 (98%)	4 (2%)	57	53
1	D	188/224 (84%)	186 (99%)	2 (1%)	73	73
2	E	10/10 (100%)	7 (70%)	3 (30%)	0	0
All	All	845/906 (93%)	830 (98%)	15 (2%)	59	55

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	239	GLN
1	B	144	GLN
1	B	161	LEU
1	B	197	GLN
1	B	254	GLU
1	B	290	LEU
1	C	15	GLN
1	C	28	LEU
1	C	71	GLN
1	C	266	GLU
1	D	35	LEU
1	D	44	GLU
2	E	1	ARG
2	E	5	ARG
2	E	7	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 13 ligands modelled in this entry, 13 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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, 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	282/286 (98%)	0.54	20 (7%) 16 17	29, 40, 74, 92	0
1	B	269/286 (94%)	0.48	21 (7%) 13 14	31, 42, 77, 102	0
1	C	270/286 (94%)	0.62	22 (8%) 12 13	29, 37, 78, 92	0
1	D	241/286 (84%)	0.25	7 (2%) 51 54	31, 38, 61, 75	0
2	E	10/10 (100%)	0.98	1 (10%) 7 8	48, 54, 61, 67	0
All	All	1072/1154 (92%)	0.48	71 (6%) 18 20	29, 39, 74, 102	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	35	LEU	18.1
1	C	35	LEU	11.9
1	C	32	LEU	10.1
1	B	31	ALA	9.2
1	C	31	ALA	8.5
1	C	17	VAL	8.1
1	A	281	LYS	7.8
1	C	38	ILE	7.6
1	C	18	GLN	7.4
1	B	38	ILE	7.2
1	C	30	SER	6.9
1	B	28	LEU	6.7
1	A	279	ILE	6.5
1	A	283	ALA	6.5
1	A	276	ASN	6.4
1	B	34	LYS	6.2
1	A	282	GLU	6.0
1	B	23	PRO	5.9
1	B	26	GLN	5.4
1	B	30	SER	5.3

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Mol	Chain	Res	Type	RSRZ
1	B	29	GLN	5.2
1	B	37	GLN	5.2
1	A	280	GLN	4.8
1	A	285	GLU	4.6
1	C	243[A]	TRP	4.6
1	C	28	LEU	4.6
1	B	33	ARG	4.6
1	B	24	ASP	4.4
1	C	29	GLN	4.4
1	A	293	HIS	4.3
1	D	41	GLY	4.1
1	A	284	GLN	4.1
1	B	27	GLU	4.0
1	B	39	ALA	4.0
1	B	32	LEU	3.8
1	A	16	MET	3.8
1	C	19	GLN	3.8
1	C	34	LYS	3.7
1	A	41	GLY	3.6
1	B	25	GLN	3.5
1	D	12	GLU	3.4
1	C	36	SER	3.4
1	A	288	GLU	3.4
1	C	37	GLN	3.2
1	A	278	LYS	3.1
1	C	15	GLN	3.1
1	D	13	LEU	3.1
1	A	58	LEU	3.0
1	C	274	HIS	3.0
1	C	16	MET	3.0
1	A	273	SER	2.9
1	C	33	ARG	2.9
1	C	39	ALA	2.8
1	B	36	SER	2.8
1	B	284	GLN	2.8
1	D	34	LYS	2.7
1	A	277	GLU	2.7
1	B	265	LEU	2.7
1	D	37	GLN	2.6
1	D	35	LEU	2.6
1	A	291	GLN	2.6
1	B	41	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	77	LEU	2.5
1	B	287	LEU	2.5
1	C	290	LEU	2.4
1	A	289	LYS	2.4
1	D	15	GLN	2.4
1	A	287	LEU	2.2
1	C	266	GLU	2.1
1	C	292	SER	2.1
2	E	1	ARG	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 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(Å ²)	Q<0.9
3	CA	B	304	1/1	0.92	0.06	60,60,60,60	1
3	CA	A	301	1/1	0.92	0.05	53,53,53,53	0
3	CA	A	305	1/1	0.94	0.17	47,47,47,47	0
3	CA	D	301	1/1	0.95	0.06	59,59,59,59	1
3	CA	A	307	1/1	0.97	0.16	54,54,54,54	0
3	CA	A	302	1/1	0.97	0.10	37,37,37,37	0
3	CA	B	301	1/1	0.97	0.10	44,44,44,44	0
3	CA	A	304	1/1	0.97	0.15	38,38,38,38	0
3	CA	A	306	1/1	0.98	0.03	43,43,43,43	0
3	CA	A	303	1/1	0.98	0.10	46,46,46,46	0
3	CA	B	303	1/1	0.99	0.08	40,40,40,40	0
3	CA	B	302	1/1	0.99	0.12	41,41,41,41	0
3	CA	B	305	1/1	0.99	0.13	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.