



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 08:00 am BST

PDB ID : 1M21
Title : Crystal structure analysis of the peptide amidase PAM in complex with the competitive inhibitor chymostatin
Authors : Labahn, J.; Neumann, S.; Buldt, G.; Kula, M.-R.; Granzin, J.
Deposited on : 2002-06-21
Resolution : 1.80 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

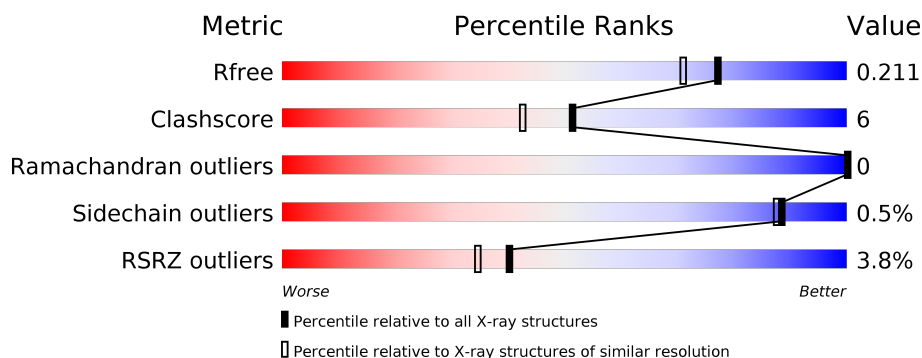
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	503	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>10%</div> <div></div> </div> </div>
1	B	503	<div> <div>4%</div> <div> <div></div> <div>86%</div> <div>10%</div> <div></div> </div> </div>
2	C	4	<div> <div>25%</div> <div> <div>50%</div> <div>25%</div> <div>25%</div> </div> </div>
2	D	4	<div> <div>25%</div> <div> <div>50%</div> <div>25%</div> <div>25%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7821 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 Peptide Amidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	487	Total	C	N	O	S	0	0	0
			3651	2283	667	691	10			
1	B	487	Total	C	N	O	S	0	0	0
			3651	2283	667	691	10			

- Molecule 2 is a protein called CHYMOSTATIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	0	0	0
			44	31	7	6			
2	D	4	Total	C	N	O	0	0	0
			44	31	7	6			

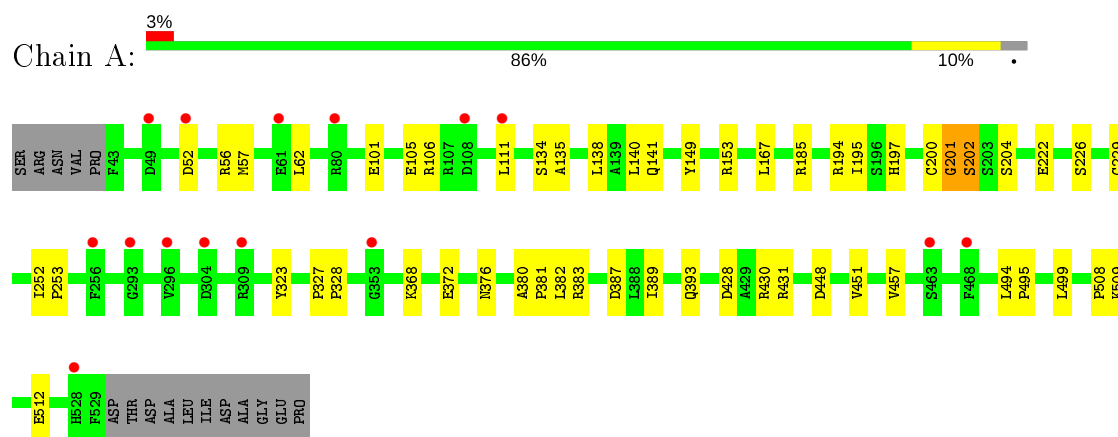
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	212	Total	O	0	0
			212	212		
3	B	219	Total	O	0	0
			219	219		

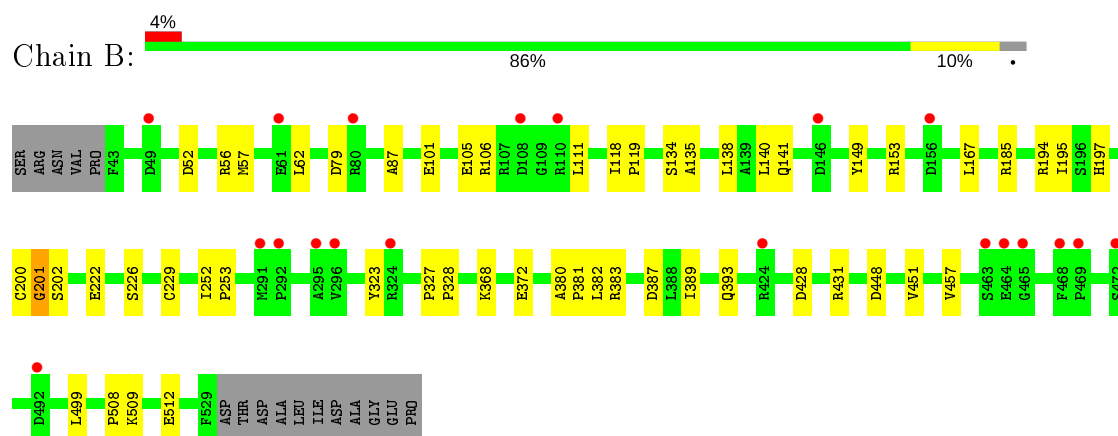
3 Residue-property plots [i](#)

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.

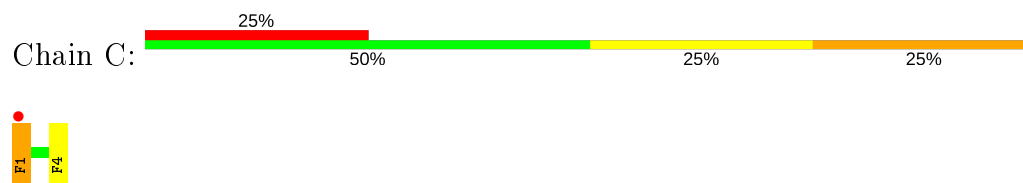
• Molecule 1: Peptide Amidase



• Molecule 1: Peptide Amidase



• Molecule 2: CHYMOSTATIN



• Molecule 2: CHYMOSTATIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.60 Å 62.56 Å 102.38 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.29 – 1.80 60.29 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.4 (60.29-1.80) 98.0 (60.29-1.78)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 1.78 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.203 , 0.217 0.196 , 0.211	Depositor DCC
R_{free} test set	4379 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	13.4	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 26.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.429 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7821	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PHA, CSI

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	0.41	1/3724 (0.0%)	0.66	4/5070 (0.1%)
1	B	0.34	0/3724	0.61	1/5070 (0.0%)
2	C	2.50	1/19 (5.3%)	2.10	0/21
2	D	2.50	1/19 (5.3%)	2.09	0/21
All	All	0.42	3/7486 (0.0%)	0.65	5/10182 (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	B	0	1
2	C	1	0
2	D	1	0
All	All	2	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	201	GLY	C-N	13.49	1.65	1.34
2	C	1	PHE	C-O	-7.00	1.10	1.23
2	D	1	PHE	C-O	-6.98	1.10	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	GLY	C-N-CA	12.40	152.70	121.70
1	A	201	GLY	CA-C-N	-10.48	94.15	117.20
1	A	201	GLY	O-C-N	8.71	136.64	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	201	GLY	C-N-CA	6.57	138.13	121.70
1	A	202	SER	CB-CA-C	5.29	120.15	110.10

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	2	CSI	CB
2	D	2	CSI	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	201	GLY	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	3651	0	3645	42	0
1	B	3651	0	3644	39	0
2	C	44	0	38	5	0
2	D	44	0	38	5	0
3	A	212	0	0	4	0
3	B	219	0	0	2	0
All	All	7821	0	7365	83	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 6.

All (83) 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)
2:C:4:PHA:O	2:C:4:PHA:C	1.63	1.44
2:D:4:PHA:C	2:D:4:PHA:O	1.63	1.43
1:B:428:ASP:HA	1:B:431:ARG:HH12	1.29	0.95
1:A:428:ASP:HA	1:A:431:ARG:HH12	1.29	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:ASP:HA	1:A:431:ARG:NH1	2.00	0.76
1:A:141:GLN:HG2	1:A:185:ARG:HD2	1.68	0.75
1:B:141:GLN:HG2	1:B:185:ARG:HD2	1.68	0.75
1:B:428:ASP:HA	1:B:431:ARG:NH1	2.00	0.75
1:A:448:ASP:HB3	1:A:509:LYS:HD3	1.74	0.69
1:A:430:ARG:HD3	3:A:1088:HOH:O	1.92	0.67
1:B:448:ASP:HB3	1:B:509:LYS:HD3	1.74	0.67
1:B:431:ARG:NH1	1:B:431:ARG:HB3	2.10	0.66
1:A:431:ARG:HB3	1:A:431:ARG:NH1	2.11	0.65
1:A:393:GLN:NE2	3:A:1029:HOH:O	2.28	0.65
1:B:431:ARG:HB3	1:B:431:ARG:HH11	1.64	0.63
1:B:327:PRO:HB2	1:B:328:PRO:HD3	1.81	0.62
1:A:327:PRO:HB2	1:A:328:PRO:HD3	1.82	0.62
1:B:194:ARG:NH1	3:B:2098:HOH:O	2.33	0.62
1:A:101:GLU:O	1:A:105:GLU:HG3	2.01	0.60
1:A:431:ARG:HB3	1:A:431:ARG:HH11	1.64	0.60
1:B:194:ARG:HD2	1:B:197:HIS:CE1	2.37	0.59
1:A:226:SER:OG	2:C:4:PHA:C	2.51	0.59
1:B:138:LEU:HA	1:B:141:GLN:HG3	1.84	0.59
1:B:101:GLU:O	1:B:105:GLU:HG3	2.02	0.59
1:B:200:CYS:SG	1:B:229:CYS:SG	2.96	0.59
1:A:138:LEU:HA	1:A:141:GLN:HG3	1.84	0.58
1:A:194:ARG:HD2	1:A:197:HIS:CE1	2.38	0.58
1:A:226:SER:HG	2:C:4:PHA:HC	1.68	0.58
1:B:226:SER:OG	2:D:4:PHA:C	2.51	0.58
1:B:106:ARG:HG3	1:B:111:LEU:HD21	1.86	0.57
1:A:222:GLU:HA	1:A:226:SER:HB2	1.86	0.57
1:B:222:GLU:HA	1:B:226:SER:HB2	1.85	0.57
1:A:106:ARG:HG3	1:A:111:LEU:HD21	1.87	0.56
1:B:226:SER:OG	2:D:4:PHA:HC	2.06	0.56
1:B:226:SER:HG	2:D:4:PHA:HC	1.71	0.56
1:A:226:SER:OG	2:C:4:PHA:HC	2.06	0.54
1:A:508:PRO:O	1:A:512:GLU:HG3	2.07	0.54
1:A:430:ARG:CD	3:A:1088:HOH:O	2.52	0.53
1:B:508:PRO:O	1:B:512:GLU:HG3	2.08	0.52
1:A:57:MET:HA	1:A:62:LEU:O	2.10	0.52
1:A:323:TYR:OH	1:A:457:VAL:HG22	2.10	0.52
1:B:389:ILE:O	1:B:393:GLN:HG3	2.10	0.52
1:B:431:ARG:CB	1:B:431:ARG:HH11	2.22	0.51
1:B:323:TYR:OH	1:B:457:VAL:HG22	2.10	0.51
1:A:431:ARG:CB	1:A:431:ARG:HH11	2.23	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:CYS:SG	1:A:229:CYS:SG	2.96	0.51
1:A:389:ILE:O	1:A:393:GLN:HG3	2.11	0.50
1:B:57:MET:HA	1:B:62:LEU:O	2.10	0.50
1:B:202:SER:HB3	2:D:4:PHA:HC	1.94	0.49
1:B:431:ARG:NH1	3:B:2104:HOH:O	2.45	0.49
1:A:200:CYS:SG	1:A:201:GLY:N	2.83	0.47
1:A:135:ALA:HB2	1:A:252:ILE:HD12	1.97	0.47
1:A:368:LYS:O	1:A:372:GLU:HG3	2.15	0.47
1:A:376:ASN:HB2	3:A:961:HOH:O	2.14	0.46
1:B:149:TYR:CE2	1:B:153:ARG:HD2	2.50	0.46
1:B:382:LEU:HD22	1:B:387:ASP:HB3	1.97	0.46
1:A:149:TYR:CE2	1:A:153:ARG:HD2	2.51	0.46
1:A:194:ARG:HH11	1:A:194:ARG:HG3	1.81	0.46
1:A:382:LEU:HD22	1:A:387:ASP:HB3	1.98	0.46
1:A:194:ARG:HD2	1:A:197:HIS:ND1	2.31	0.46
1:B:135:ALA:HB2	1:B:252:ILE:HD12	1.97	0.46
1:B:368:LYS:O	1:B:372:GLU:HG3	2.16	0.45
1:B:167:LEU:HD23	1:B:167:LEU:N	2.31	0.45
1:B:194:ARG:HD2	1:B:197:HIS:ND1	2.31	0.44
1:A:52:ASP:O	1:A:56:ARG:HG3	2.18	0.44
1:B:194:ARG:HH11	1:B:194:ARG:HG3	1.83	0.44
1:B:451:VAL:HA	1:B:499:LEU:O	2.18	0.44
1:A:167:LEU:N	1:A:167:LEU:HD23	2.33	0.44
1:A:134:SER:HA	1:A:140:LEU:HB2	2.00	0.44
1:A:195:ILE:C	1:A:195:ILE:HD12	2.38	0.44
1:A:451:VAL:HA	1:A:499:LEU:O	2.17	0.43
1:B:380:ALA:HB1	1:B:381:PRO:HD2	2.00	0.43
1:B:134:SER:HA	1:B:140:LEU:HB2	2.00	0.43
1:B:52:ASP:O	1:B:56:ARG:HG3	2.18	0.43
1:B:195:ILE:C	1:B:195:ILE:HD12	2.38	0.43
1:A:226:SER:HG	2:C:4:PHA:C	2.29	0.43
1:A:380:ALA:HB1	1:A:381:PRO:HD2	2.00	0.42
1:A:382:LEU:O	1:A:383:ARG:HD3	2.20	0.42
1:B:118:ILE:HA	1:B:119:PRO:HD3	1.88	0.41
1:B:79:ASP:OD2	1:B:87:ALA:N	2.42	0.41
1:B:382:LEU:O	1:B:383:ARG:HD3	2.19	0.41
1:A:202:SER:O	1:A:204:SER:N	2.50	0.40
1:A:494:LEU:HA	1:A:495:PRO:HD3	1.95	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	485/503 (96%)	467 (96%)	18 (4%)	0	100	100
1	B	485/503 (96%)	468 (96%)	17 (4%)	0	100	100
2	C	1/4 (25%)	0	1 (100%)	0	100	100
2	D	1/4 (25%)	0	1 (100%)	0	100	100
All	All	972/1014 (96%)	935 (96%)	37 (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	366/379 (97%)	365 (100%)	1 (0%)	92	91
1	B	366/379 (97%)	365 (100%)	1 (0%)	92	91
2	C	2/2 (100%)	1 (50%)	1 (50%)	0	0
2	D	2/2 (100%)	1 (50%)	1 (50%)	0	0
All	All	736/762 (97%)	732 (100%)	4 (0%)	88	87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	PRO
1	B	253	PRO

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Mol	Chain	Res	Type
2	C	1	PHE
2	D	1	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	446	GLN
1	B	393	GLN
1	B	446	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	PHA	C	4	2	10,11,11	3.83	2 (20%)	10,13,13	0.95	0
2	PHA	D	4	2	10,11,11	3.83	2 (20%)	10,13,13	0.95	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	PHA	C	4	2	-	0/5/6/6	0/1/1/1
2	PHA	D	4	2	-	0/5/6/6	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	4	PHA	O-C	11.01	1.63	1.19
2	C	4	PHA	O-C	11.00	1.63	1.19
2	C	4	PHA	CD2-CG	2.90	1.45	1.38
2	D	4	PHA	CD2-CG	2.87	1.45	1.38

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	4	PHA	5	0
2	D	4	PHA	5	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1
2	D	1
2	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	1:PHE	C	2:CSI	N	4.39
1	D	1:PHE	C	2:CSI	N	4.39
1	A	201:GLY	C	202:SER	N	1.65

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	487/503 (96%)	0.37	15 (3%)	49	43	5, 12, 24, 32	0
1	B	487/503 (96%)	0.40	20 (4%)	37	31	5, 12, 24, 32	0
2	C	2/4 (50%)	3.16	1 (50%)	0	0	20, 20, 20, 32	0
2	D	2/4 (50%)	3.42	1 (50%)	0	0	20, 20, 20, 32	0
All	All	978/1014 (96%)	0.40	37 (3%)	40	35	5, 12, 24, 32	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	1	PHE	5.4
2	C	1	PHE	5.3
1	A	463	SER	5.1
1	B	292	PRO	4.5
1	B	295	ALA	4.5
1	B	465	GLY	4.1
1	A	296	VAL	4.0
1	A	528	HIS	3.8
1	B	468	PHE	3.8
1	A	468	PHE	3.3
1	B	472	SER	2.9
1	A	293	GLY	2.9
1	A	61	GLU	2.7
1	A	108	ASP	2.7
1	B	108	ASP	2.6
1	A	80	ARG	2.6
1	B	324	ARG	2.6
1	B	464	GLU	2.5
1	A	256	PHE	2.5
1	B	291	MET	2.5
1	A	309	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	49	ASP	2.4
1	A	49	ASP	2.4
1	B	424	ARG	2.4
1	A	353	GLY	2.4
1	B	146	ASP	2.3
1	B	110	ARG	2.3
1	B	463	SER	2.3
1	B	156	ASP	2.2
1	B	492	ASP	2.2
1	B	469	PRO	2.2
1	A	304	ASP	2.2
1	B	61	GLU	2.2
1	B	80	ARG	2.2
1	A	52	ASP	2.1
1	B	296	VAL	2.1
1	A	111	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	PHA	C	4	11/11	0.66	0.25	22,25,27,27	0
2	PHA	D	4	11/11	0.78	0.23	22,25,27,27	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.