



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 26, 2016 – 02:17 PM EDT

PDB ID : 5CGG
Title : Yeast 20S proteasome beta5-G48C mutant in complex with alpha-chloroacetamide 1
Authors : Dubiella, C.; Groll, M.
Deposited on : 2015-07-09
Resolution : 2.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027257
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027257

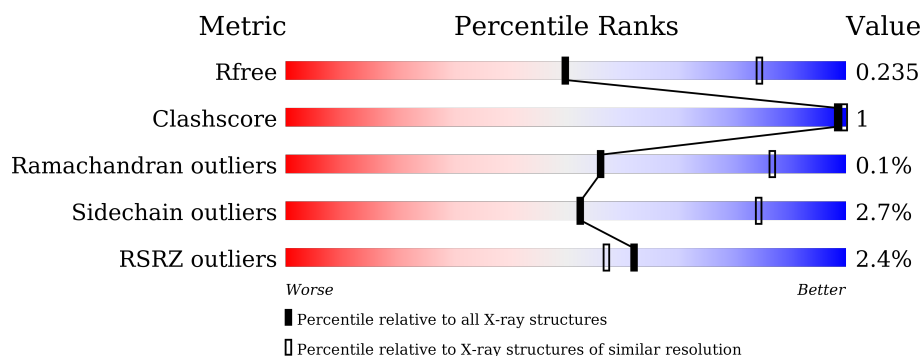
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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. 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	250	<div> <div>3%</div> <div> <div></div> <div>98%</div> <div>.</div> </div> </div>
1	O	250	<div> <div>3%</div> <div> <div></div> <div>99%</div> <div>.</div> </div> </div>
2	B	258	<div> <div>3%</div> <div> <div></div> <div>90%</div> <div>.. 5%</div> </div> </div>
2	P	258	<div> <div>5%</div> <div> <div></div> <div>90%</div> <div>.. 5%</div> </div> </div>
3	C	254	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>.. 6%</div> </div> </div>
3	Q	254	<div> <div>7%</div> <div> <div></div> <div>90%</div> <div>.. 6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	D	260	
4	R	260	
5	E	234	
5	S	234	
6	F	288	
6	T	288	
7	G	252	
7	U	252	
8	H	232	
8	V	232	
9	I	205	
9	W	205	
10	J	198	
10	X	198	
11	K	212	
11	Y	212	
12	L	222	
12	Z	222	
13	M	246	
13	a	246	
14	N	196	
14	b	196	
15	g	6	
15	h	6	

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
16	MG	J	201	-	-	-	X
16	MG	N	201	-	-	-	X
16	MG	V	301	-	-	-	X
18	MES	K	302	-	-	-	X
18	MES	Y	302	-	-	-	X

2 Entry composition

There are 19 unique types of molecules in this entry. The entry contains 49712 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 Proteasome subunit alpha type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	0	0
			1914	1219	315	376	4			
1	O	250	Total	C	N	O	S	0	0	0
			1914	1219	315	376	4			

- Molecule 2 is a protein called Proteasome subunit alpha type-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			
2	P	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			

- Molecule 3 is a protein called Proteasome subunit alpha type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	240	Total	C	N	O	S	0	0	0
			1881	1176	329	372	4			
3	Q	240	Total	C	N	O	S	0	0	0
			1881	1176	329	372	4			

- Molecule 4 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	235	Total	C	N	O	S	0	0	0
			1813	1136	304	366	7			
4	R	235	Total	C	N	O	S	0	0	0
			1813	1136	304	366	7			

- Molecule 5 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	231	Total	C	N	O	S	0	0	0
			1772	1114	307	347	4			
5	S	231	Total	C	N	O	S	0	0	0
			1772	1114	307	347	4			

- Molecule 6 is a protein called Probable proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	243	Total	C	N	O	S	0	0	0
			1892	1203	329	356	4			
6	T	243	Total	C	N	O	S	0	0	0
			1892	1203	329	356	4			

- Molecule 7 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	241	Total	C	N	O	S	0	0	0
			1907	1214	320	365	8			
7	U	241	Total	C	N	O	S	0	0	0
			1906	1214	320	364	8			

- Molecule 8 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			
8	V	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			

- Molecule 9 is a protein called Proteasome subunit beta type-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	204	Total	C	N	O	S	0	0	0
			1580	1010	258	304	8			
9	W	204	Total	C	N	O	S	0	0	0
			1580	1010	258	304	8			

- Molecule 10 is a protein called Proteasome subunit beta type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	X	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			

- Molecule 11 is a protein called Proteasome subunit beta type-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	212	Total	C	N	O	S	0	0	0
			1645	1046	280	311	8			
11	Y	212	Total	C	N	O	S	0	0	0
			1645	1046	280	311	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	48	CYS	GLY	engineered mutation	UNP P30656
Y	48	CYS	GLY	engineered mutation	UNP P30656

- Molecule 12 is a protein called Proteasome subunit beta type-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	222	Total	C	N	O	S	0	0	0
			1756	1115	303	334	4			
12	Z	222	Total	C	N	O	S	0	0	0
			1756	1115	303	334	4			

- Molecule 13 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	233	Total	C	N	O	S	0	0	0
			1823	1154	312	350	7			
13	a	233	Total	C	N	O	S	0	0	0
			1823	1154	312	350	7			

- Molecule 14 is a protein called Proteasome subunit beta type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	196	Total	C	N	O	S	0	0	0
			1511	955	250	299	7			
14	b	196	Total	C	N	O	S	0	0	0
			1511	955	250	299	7			

- Molecule 15 is a protein called carfilzomib alpha-chloroacetamide 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	g	6	Total	C	N	O	0	0	0
			46	34	6	6			
15	h	6	Total	C	N	O	0	0	0
			46	34	6	6			

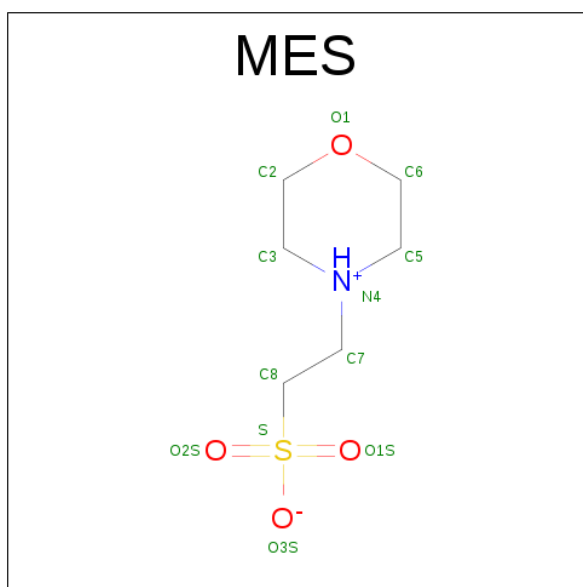
- Molecule 16 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	G	1	Total	Mg	0	0
			1	1		
16	J	1	Total	Mg	0	0
			1	1		
16	K	1	Total	Mg	0	0
			1	1		
16	I	1	Total	Mg	0	0
			1	1		
16	V	1	Total	Mg	0	0
			1	1		
16	Z	1	Total	Mg	0	0
			1	1		
16	N	1	Total	Mg	0	0
			1	1		
16	Y	1	Total	Mg	0	0
			1	1		

- Molecule 17 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	G	2	Total	Cl	0	0
			2	2		
17	N	1	Total	Cl	0	0
			1	1		
17	U	1	Total	Cl	0	0
			1	1		

- Molecule 18 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
18	K	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
18	Y	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 19 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	A	10	Total	O	0	0
			10	10		
19	B	17	Total	O	0	0
			17	17		
19	C	9	Total	O	0	0
			9	9		
19	D	5	Total	O	0	0
			5	5		
19	E	7	Total	O	0	0
			7	7		
19	F	16	Total	O	0	0
			16	16		
19	G	13	Total	O	0	0
			13	13		
19	H	14	Total	O	0	0
			14	14		
19	I	9	Total	O	0	0
			9	9		
19	J	19	Total	O	0	0
			19	19		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	K	9	Total 9	O 9	0	0
19	L	10	Total 10	O 10	0	0
19	M	12	Total 12	O 12	0	0
19	N	13	Total 13	O 13	0	0
19	O	8	Total 8	O 8	0	0
19	P	4	Total 4	O 4	0	0
19	Q	10	Total 10	O 10	0	0
19	R	4	Total 4	O 4	0	0
19	S	6	Total 6	O 6	0	0
19	T	12	Total 12	O 12	0	0
19	U	15	Total 15	O 15	0	0
19	V	13	Total 13	O 13	0	0
19	W	8	Total 8	O 8	0	0
19	X	8	Total 8	O 8	0	0
19	Y	8	Total 8	O 8	0	0
19	Z	10	Total 10	O 10	0	0
19	a	12	Total 12	O 12	0	0
19	b	17	Total 17	O 17	0	0
19	h	1	Total 1	O 1	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 errors 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: Proteasome subunit alpha type-2



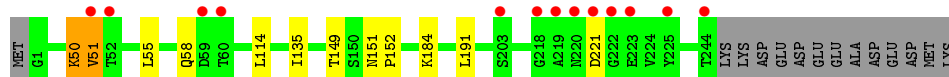
- Molecule 1: Proteasome subunit alpha type-2



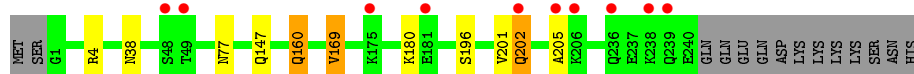
- Molecule 2: Proteasome subunit alpha type-3



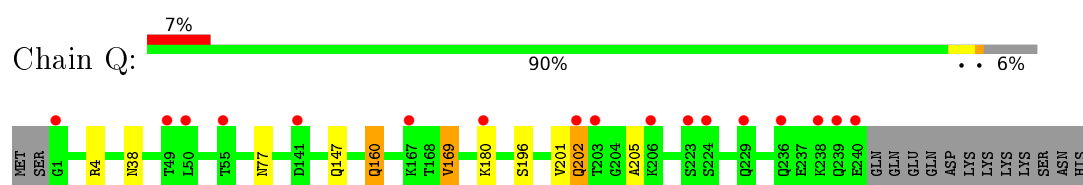
- Molecule 2: Proteasome subunit alpha type-3



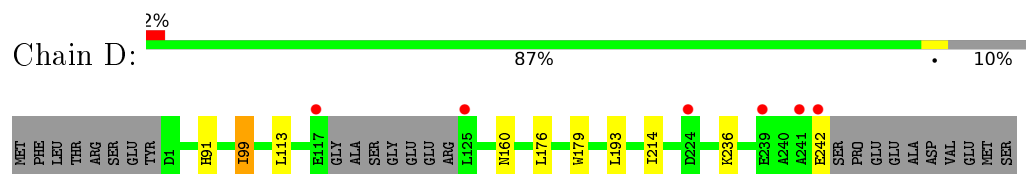
- Molecule 3: Proteasome subunit alpha type-4



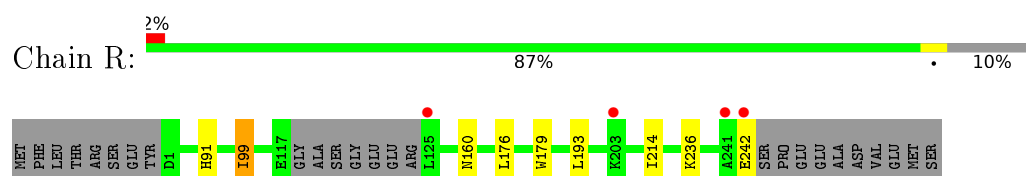
- Molecule 3: Proteasome subunit alpha type-4



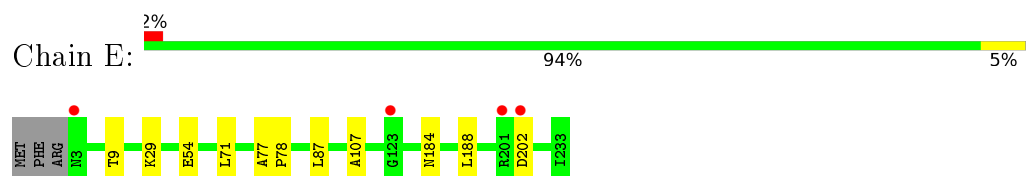
- Molecule 4: Proteasome subunit alpha type-5



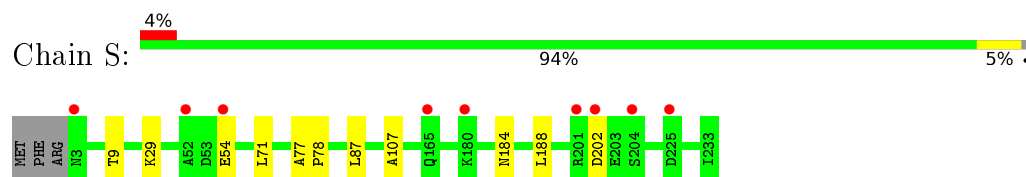
- Molecule 4: Proteasome subunit alpha type-5



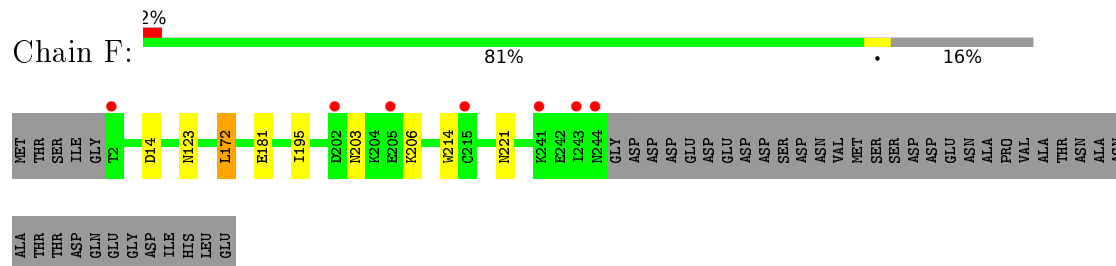
- Molecule 5: Proteasome subunit alpha type-6



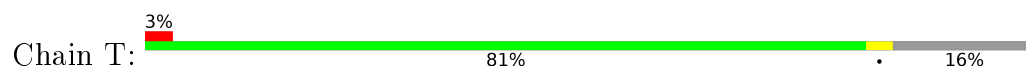
- Molecule 5: Proteasome subunit alpha type-6

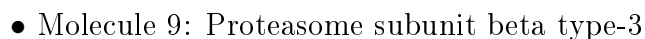
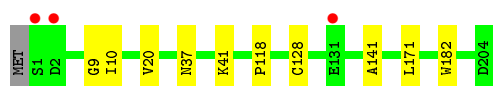
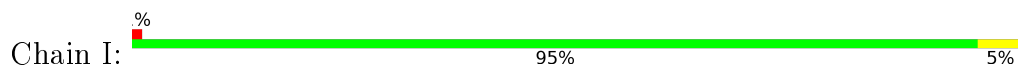
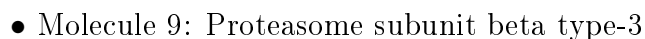
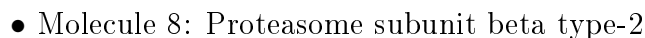
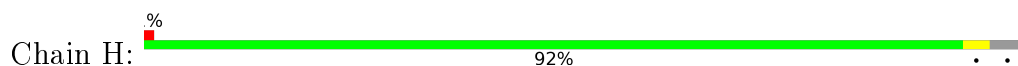
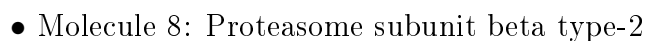
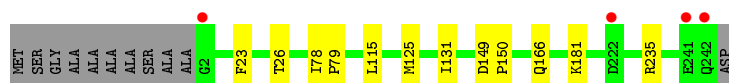
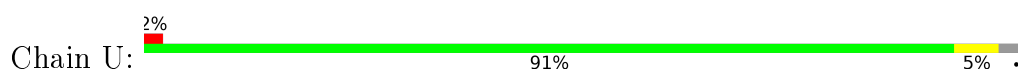
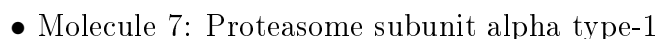
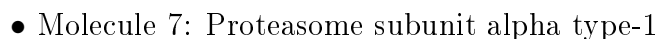


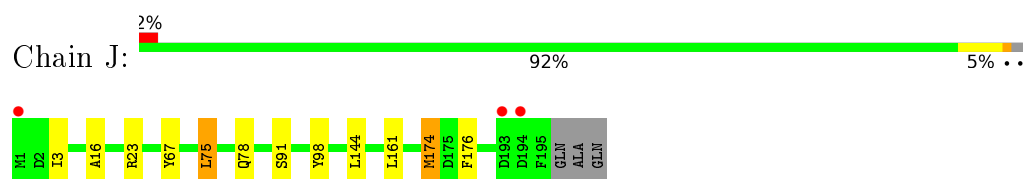
- Molecule 6: Probable proteasome subunit alpha type-7



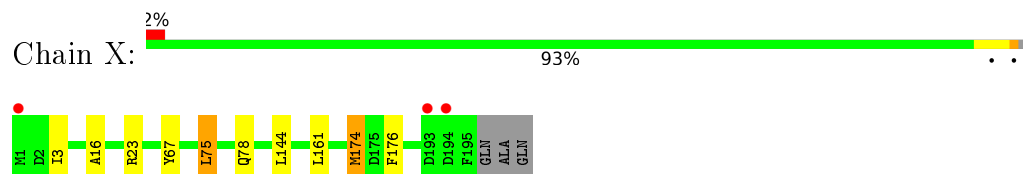
- Molecule 6: Probable proteasome subunit alpha type-7



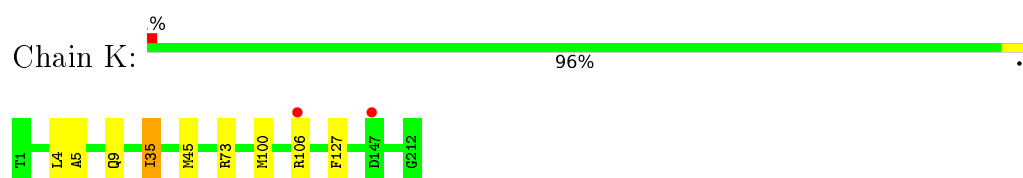




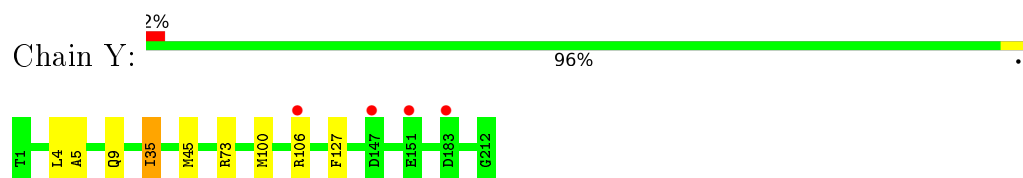
- Molecule 10: Proteasome subunit beta type-4



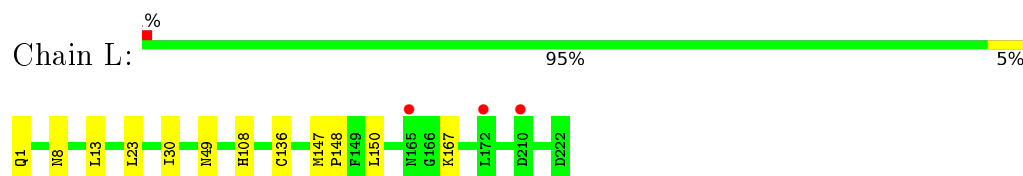
- Molecule 11: Proteasome subunit beta type-5



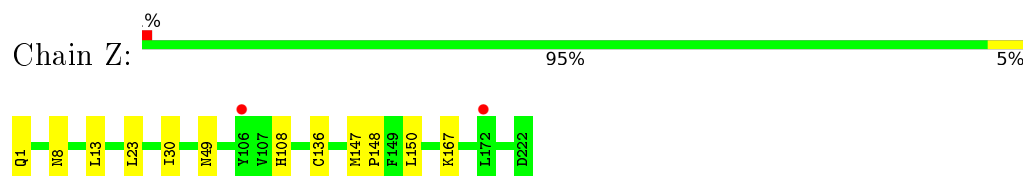
- Molecule 11: Proteasome subunit beta type-5



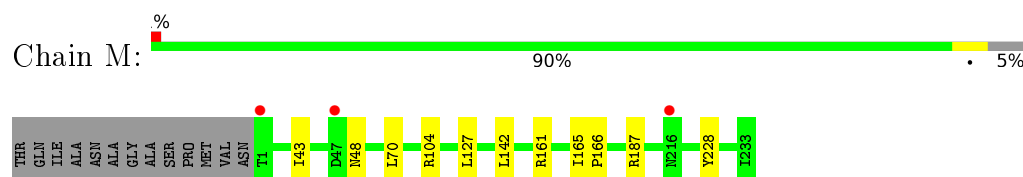
- Molecule 12: Proteasome subunit beta type-6



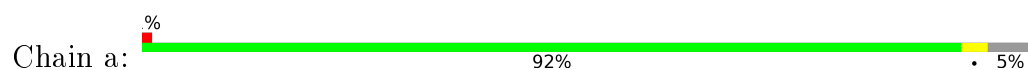
- Molecule 12: Proteasome subunit beta type-6



- Molecule 13: Proteasome subunit beta type-7



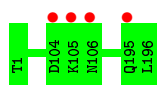
- Molecule 13: Proteasome subunit beta type-7



- Molecule 14: Proteasome subunit beta type-1



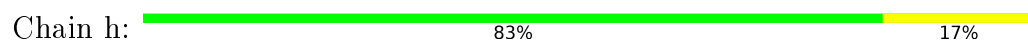
- Molecule 14: Proteasome subunit beta type-1



- Molecule 15: carfilzomib alpha-chloroacetamide 1



- Molecule 15: carfilzomib alpha-chloroacetamide 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	134.44Å 300.47Å 144.74Å 90.00° 113.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 15.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.7 (15.00-2.90) 95.7 (15.00-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.204 , 0.231 0.210 , 0.235	Depositor DCC
R_{free} test set	11055 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	58.1	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 38.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	49712	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, DPP, 00E, CL, 02N, MES, ACY

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	0.27	0/1951	0.46	0/2642
1	O	0.27	0/1951	0.46	0/2642
2	B	0.27	0/1934	0.48	0/2618
2	P	0.27	0/1934	0.48	0/2618
3	C	0.27	0/1910	0.50	0/2586
3	Q	0.27	0/1910	0.50	0/2586
4	D	0.27	0/1837	0.47	0/2475
4	R	0.26	0/1837	0.46	0/2475
5	E	0.27	0/1799	0.46	0/2433
5	S	0.27	0/1799	0.46	0/2433
6	F	0.27	0/1932	0.45	0/2609
6	T	0.27	0/1932	0.45	0/2609
7	G	0.27	0/1945	0.47	0/2634
7	U	0.27	0/1944	0.46	0/2632
8	H	0.26	0/1715	0.47	0/2326
8	V	0.26	0/1715	0.47	0/2326
9	I	0.27	0/1610	0.47	0/2174
9	W	0.27	0/1610	0.47	0/2174
10	J	0.26	0/1589	0.48	0/2142
10	X	0.26	0/1589	0.48	0/2142
11	K	0.27	0/1682	0.47	0/2277
11	Y	0.27	0/1682	0.47	0/2277
12	L	0.27	0/1794	0.46	0/2420
12	Z	0.27	0/1794	0.46	0/2420
13	M	0.28	0/1854	0.51	0/2514
13	a	0.27	0/1854	0.51	0/2514
14	N	0.26	0/1540	0.46	0/2087
14	b	0.26	0/1540	0.46	0/2087
15	g	2.13	1/19 (5.3%)	1.19	0/24
15	h	1.98	1/19 (5.3%)	1.04	0/24
All	All	0.27	2/50221 (0.0%)	0.47	0/67920

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	g	5	PHE	CB-CG	-8.56	1.36	1.51
15	h	5	PHE	CB-CG	-8.07	1.37	1.51

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	1914	0	1929	1	0
1	O	1914	0	1929	0	0
2	B	1904	0	1904	5	0
2	P	1904	0	1904	4	0
3	C	1881	0	1895	4	0
3	Q	1881	0	1895	4	0
4	D	1813	0	1797	3	0
4	R	1813	0	1797	2	0
5	E	1772	0	1775	2	0
5	S	1772	0	1775	2	0
6	F	1892	0	1883	1	0
6	T	1892	0	1883	1	0
7	G	1907	0	1901	4	0
7	U	1906	0	1901	4	0
8	H	1684	0	1688	3	0
8	V	1684	0	1688	3	0
9	I	1580	0	1574	4	0
9	W	1580	0	1574	3	0
10	J	1561	0	1569	5	0
10	X	1561	0	1569	4	0
11	K	1645	0	1596	6	0
11	Y	1645	0	1596	5	0
12	L	1756	0	1711	3	0
12	Z	1756	0	1711	3	0
13	M	1823	0	1832	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	a	1823	0	1832	0	0
14	N	1511	0	1481	1	0
14	b	1511	0	1481	0	0
15	g	46	0	43	0	0
15	h	46	0	43	0	0
16	G	1	0	0	0	0
16	I	1	0	0	0	0
16	J	1	0	0	0	0
16	K	1	0	0	0	0
16	N	1	0	0	0	0
16	V	1	0	0	0	0
16	Y	1	0	0	0	0
16	Z	1	0	0	0	0
17	G	2	0	0	0	0
17	N	1	0	0	0	0
17	U	1	0	0	0	0
18	K	12	0	13	0	0
18	Y	12	0	13	0	0
19	A	10	0	0	1	0
19	B	17	0	0	0	0
19	C	9	0	0	0	0
19	D	5	0	0	1	0
19	E	7	0	0	0	0
19	F	16	0	0	0	0
19	G	13	0	0	1	0
19	H	14	0	0	0	0
19	I	9	0	0	0	0
19	J	19	0	0	0	0
19	K	9	0	0	0	0
19	L	10	0	0	0	0
19	M	12	0	0	0	0
19	N	13	0	0	0	0
19	O	8	0	0	0	0
19	P	4	0	0	0	0
19	Q	10	0	0	0	0
19	R	4	0	0	0	0
19	S	6	0	0	0	0
19	T	12	0	0	0	0
19	U	15	0	0	0	0
19	V	13	0	0	0	0
19	W	8	0	0	0	0
19	X	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
19	Y	8	0	0	0	0
19	Z	10	0	0	0	0
19	a	12	0	0	0	0
19	b	17	0	0	0	0
19	h	1	0	0	0	0
All	All	49712	0	49182	75	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 1.

The worst 5 of 75 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)
12:Z:13:LEU:HD13	12:Z:150:LEU:HD21	1.83	0.60
12:L:13:LEU:HD13	12:L:150:LEU:HD21	1.84	0.59
11:Y:5:ALA:HB3	11:Y:100:MET:HE2	1.84	0.59
11:K:35:ILE:HB	11:K:45:MET:CE	2.35	0.55
11:Y:35:ILE:HB	11:Y:45:MET:CE	2.36	0.55

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	248/250 (99%)	241 (97%)	7 (3%)	0	100	100
1	O	248/250 (99%)	241 (97%)	7 (3%)	0	100	100
2	B	242/258 (94%)	234 (97%)	6 (2%)	2 (1%)	24	60
2	P	242/258 (94%)	234 (97%)	6 (2%)	2 (1%)	24	60
3	C	238/254 (94%)	230 (97%)	6 (2%)	2 (1%)	24	60

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	Q	238/254 (94%)	230 (97%)	6 (2%)	2 (1%)	24	60
4	D	231/260 (89%)	225 (97%)	6 (3%)	0	100	100
4	R	231/260 (89%)	225 (97%)	6 (3%)	0	100	100
5	E	229/234 (98%)	223 (97%)	6 (3%)	0	100	100
5	S	229/234 (98%)	223 (97%)	6 (3%)	0	100	100
6	F	241/288 (84%)	235 (98%)	6 (2%)	0	100	100
6	T	241/288 (84%)	234 (97%)	7 (3%)	0	100	100
7	G	239/252 (95%)	235 (98%)	4 (2%)	0	100	100
7	U	239/252 (95%)	235 (98%)	4 (2%)	0	100	100
8	H	220/232 (95%)	213 (97%)	7 (3%)	0	100	100
8	V	220/232 (95%)	214 (97%)	6 (3%)	0	100	100
9	I	202/205 (98%)	195 (96%)	7 (4%)	0	100	100
9	W	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
10	J	193/198 (98%)	189 (98%)	4 (2%)	0	100	100
10	X	193/198 (98%)	189 (98%)	4 (2%)	0	100	100
11	K	210/212 (99%)	206 (98%)	4 (2%)	0	100	100
11	Y	210/212 (99%)	206 (98%)	4 (2%)	0	100	100
12	L	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
12	Z	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
13	M	231/246 (94%)	221 (96%)	10 (4%)	0	100	100
13	a	231/246 (94%)	221 (96%)	9 (4%)	1 (0%)	39	74
14	N	194/196 (99%)	188 (97%)	6 (3%)	0	100	100
14	b	194/196 (99%)	188 (97%)	6 (3%)	0	100	100
All	All	6276/6614 (95%)	6099 (97%)	168 (3%)	9 (0%)	56	87

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL
3	C	202	GLN
2	P	51	VAL
3	Q	202	GLN
3	C	205	ALA

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	209/209 (100%)	206 (99%)	3 (1%)	74	93
1	O	209/209 (100%)	206 (99%)	3 (1%)	74	93
2	B	203/216 (94%)	197 (97%)	6 (3%)	48	83
2	P	203/216 (94%)	197 (97%)	6 (3%)	48	83
3	C	212/226 (94%)	205 (97%)	7 (3%)	45	80
3	Q	212/226 (94%)	205 (97%)	7 (3%)	45	80
4	D	194/215 (90%)	188 (97%)	6 (3%)	47	82
4	R	194/215 (90%)	188 (97%)	6 (3%)	47	82
5	E	190/193 (98%)	183 (96%)	7 (4%)	41	77
5	S	190/193 (98%)	183 (96%)	7 (4%)	41	77
6	F	201/239 (84%)	193 (96%)	8 (4%)	38	74
6	T	201/239 (84%)	193 (96%)	8 (4%)	38	74
7	G	206/210 (98%)	201 (98%)	5 (2%)	57	86
7	U	206/210 (98%)	201 (98%)	5 (2%)	57	86
8	H	181/190 (95%)	177 (98%)	4 (2%)	60	88
8	V	181/190 (95%)	177 (98%)	4 (2%)	60	88
9	I	172/173 (99%)	169 (98%)	3 (2%)	68	91
9	W	172/173 (99%)	169 (98%)	3 (2%)	68	91
10	J	173/175 (99%)	168 (97%)	5 (3%)	50	83
10	X	173/175 (99%)	168 (97%)	5 (3%)	50	83
11	K	170/170 (100%)	165 (97%)	5 (3%)	50	83
11	Y	170/170 (100%)	165 (97%)	5 (3%)	50	83
12	L	185/185 (100%)	179 (97%)	6 (3%)	46	81
12	Z	185/185 (100%)	179 (97%)	6 (3%)	46	81
13	M	199/208 (96%)	193 (97%)	6 (3%)	48	83
13	a	199/208 (96%)	193 (97%)	6 (3%)	48	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	162/162 (100%)	161 (99%)	1 (1%)	90	97
14	b	162/162 (100%)	162 (100%)	0	100	100
15	g	2/2 (100%)	1 (50%)	1 (50%)	0	0
15	h	2/2 (100%)	2 (100%)	0	100	100
All	All	5318/5546 (96%)	5174 (97%)	144 (3%)	52	84

5 of 144 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
13	M	43	ILE
2	P	191	LEU
12	Z	49	ASN
13	M	70	LEU
1	O	61	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 69 such sidechains are listed below:

Mol	Chain	Res	Type
13	M	194	ASN
3	Q	77	ASN
12	Z	79	HIS
13	M	213	GLN
2	P	123	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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
15	00E	g	1	15	9,9,10	0.68	0	9,10,12	3.40	3 (33%)
15	DPP	g	2	15	2,5,6	0.97	0	1,5,7	2.26	1 (100%)
15	00E	h	1	15	9,9,10	0.46	0	9,10,12	2.32	5 (55%)
15	DPP	h	2	15	2,5,6	1.20	0	1,5,7	2.05	1 (100%)

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
15	00E	g	1	15	-	0/2/11/12	0/1/1/1
15	DPP	g	2	15	-	0/2/4/6	0/0/0/0
15	00E	h	1	15	-	0/2/11/12	0/1/1/1
15	DPP	h	2	15	-	0/2/4/6	0/0/0/0

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	g	1	00E	C-CA-NB	-9.15	99.86	113.02
15	h	1	00E	C-CA-NB	-4.37	106.73	113.02
15	h	1	00E	CA-NB-CD1	-2.83	107.28	110.53
15	g	1	00E	CA-NB-CD2	-2.83	107.28	110.53
15	h	1	00E	CA-NB-CD2	-2.46	107.71	110.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 12 are monoatomic - leaving 2 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	$\# Z > 2$
18	MES	K	302	-	12,12,12	2.10	1 (8%)	15,16,16	1.31	2 (13%)
18	MES	Y	302	-	12,12,12	2.10	1 (8%)	15,16,16	1.44	2 (13%)

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
18	MES	K	302	-	-	0/6/14/14	0/1/1/1
18	MES	Y	302	-	-	0/6/14/14	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	K	302	MES	C8-S	-6.97	1.67	1.77
18	Y	302	MES	C8-S	-6.96	1.67	1.77

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	K	302	MES	O3S-S-C8	2.56	110.31	104.99
18	Y	302	MES	O3S-S-C8	2.95	111.12	104.99
18	K	302	MES	O1S-S-C8	3.21	109.14	106.87
18	Y	302	MES	O1S-S-C8	3.62	109.43	106.87

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
15	h	1
15	g	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	h	3:ACY	C	4:LEU	N	4.86
1	g	3:ACY	C	4:LEU	N	4.48

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	250/250 (100%)	-0.23	7 (2%) 56 50	36, 51, 91, 148	0
1	O	250/250 (100%)	-0.14	8 (3%) 51 43	41, 60, 100, 144	0
2	B	244/258 (94%)	-0.10	9 (3%) 45 38	40, 60, 105, 158	0
2	P	244/258 (94%)	-0.08	13 (5%) 30 23	41, 62, 106, 159	0
3	C	240/254 (94%)	-0.05	10 (4%) 40 33	36, 60, 122, 159	0
3	Q	240/254 (94%)	0.18	17 (7%) 19 13	47, 75, 152, 181	0
4	D	235/260 (90%)	-0.23	6 (2%) 59 54	41, 62, 92, 142	0
4	R	235/260 (90%)	-0.15	4 (1%) 73 70	47, 67, 106, 148	0
5	E	231/234 (98%)	-0.10	4 (1%) 73 70	43, 65, 98, 138	0
5	S	231/234 (98%)	-0.00	9 (3%) 43 36	43, 71, 119, 147	0
6	F	243/288 (84%)	-0.27	7 (2%) 55 49	39, 59, 108, 133	0
6	T	243/288 (84%)	-0.13	10 (4%) 41 34	40, 70, 125, 161	0
7	G	241/252 (95%)	-0.28	6 (2%) 61 55	35, 53, 91, 142	0
7	U	241/252 (95%)	-0.28	4 (1%) 73 70	36, 56, 91, 135	0
8	H	222/232 (95%)	-0.15	2 (0%) 85 84	39, 52, 87, 116	0
8	V	222/232 (95%)	-0.15	2 (0%) 85 84	41, 56, 88, 123	0
9	I	204/205 (99%)	-0.46	3 (1%) 76 74	34, 50, 78, 103	0
9	W	204/205 (99%)	-0.44	3 (1%) 76 74	33, 50, 79, 105	0
10	J	195/198 (98%)	-0.41	3 (1%) 76 74	34, 51, 79, 119	0
10	X	195/198 (98%)	-0.32	3 (1%) 76 74	35, 54, 81, 132	0
11	K	212/212 (100%)	-0.38	2 (0%) 85 84	36, 53, 83, 101	0
11	Y	212/212 (100%)	-0.28	4 (1%) 70 66	41, 55, 90, 113	0
12	L	222/222 (100%)	-0.36	3 (1%) 78 76	32, 53, 85, 112	0
12	Z	222/222 (100%)	-0.37	2 (0%) 85 84	35, 52, 87, 110	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/246 (94%)	-0.42	3 (1%) 79 78	31, 51, 76, 92	0
13	a	233/246 (94%)	-0.39	3 (1%) 79 78	32, 51, 76, 89	0
14	N	196/196 (100%)	-0.36	1 (0%) 91 90	34, 47, 75, 97	0
14	b	196/196 (100%)	-0.36	4 (2%) 68 64	36, 49, 78, 101	0
15	g	2/6 (33%)	-0.66	0 100 100	47, 47, 47, 47	0
15	h	2/6 (33%)	-0.15	0 100 100	51, 51, 51, 51	0
All	All	6340/6626 (95%)	-0.23	152 (2%) 62 57	31, 56, 102, 181	0

The worst 5 of 152 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	9.3
3	Q	50	LEU	6.2
2	B	220	ASN	6.1
2	P	220	ASN	5.5
10	X	1	MET	5.3

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
15	00E	h	1	9/10	0.75	0.58	-	56,59,60,64	0
15	DPP	h	2	6/7	0.79	0.22	-	52,54,55,57	0
15	00E	g	1	9/10	0.77	0.60	-	53,54,57,59	0
15	DPP	g	2	6/7	0.91	0.21	-	48,50,52,56	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
16	MG	J	201	1/1	0.97	0.24	7.12	30,30,30,30	0
18	MES	K	302	12/12	0.95	0.24	3.82	40,43,52,58	0
16	MG	V	301	1/1	0.76	0.21	2.59	67,67,67,67	0
16	MG	N	201	1/1	0.93	0.22	2.54	42,42,42,42	0
18	MES	Y	302	12/12	0.96	0.22	2.26	40,42,55,55	0
16	MG	Y	301	1/1	0.97	0.13	-0.32	42,42,42,42	0
16	MG	I	301	1/1	0.99	0.15	-0.47	61,61,61,61	0
16	MG	G	301	1/1	0.96	0.12	-0.59	42,42,42,42	0
16	MG	Z	301	1/1	0.97	0.09	-2.33	52,52,52,52	0
16	MG	K	301	1/1	0.94	0.09	-2.37	43,43,43,43	0
17	CL	N	202	1/1	0.95	0.13	-	30,30,30,30	0
17	CL	G	302	1/1	0.99	0.12	-	42,42,42,42	0
17	CL	U	301	1/1	0.98	0.19	-	44,44,44,44	0
17	CL	G	303	1/1	0.94	0.19	-	30,30,30,30	0

6.5 Other polymers

There are no such residues in this entry.