



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:14 PM UTC

PDB ID : 5O3W / pdb_00005o3w
Title : Structural characterization of the fast and promiscuous macrocyclase from plant - PCY1-S562A bound to Presegetalin A1
Authors : Ludewig, H.; Czekster, C.M.; Bent, A.F.; Naismith, J.H.
Deposited on : 2017-05-25
Resolution : 2.00 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

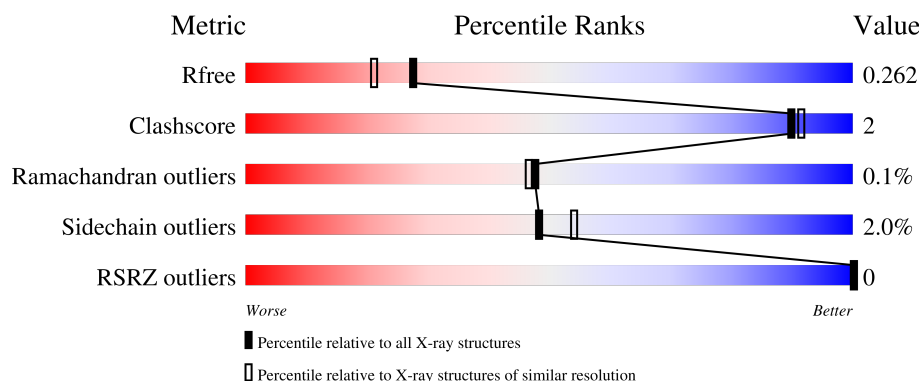
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	724	
1	B	724	
1	C	724	
1	D	724	
2	W	32	

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Mol	Chain	Length	Quality of chain
2	X	32	16% 81%
2	Y	32	19% 81%
2	Z	32	25% 75%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 23929 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 cyclase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	709	Total	C	N	O	S	0	1	0
			5703	3643	966	1067	27			
1	B	717	Total	C	N	O	S	0	2	0
			5773	3685	985	1077	26			
1	C	716	Total	C	N	O	S	0	2	0
			5765	3681	984	1074	26			
1	D	702	Total	C	N	O	S	0	1	0
			5651	3612	956	1056	27			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	562	ALA	SER	engineered mutation	UNP R4P353
B	562	ALA	SER	engineered mutation	UNP R4P353
C	562	ALA	SER	engineered mutation	UNP R4P353
D	562	ALA	SER	engineered mutation	UNP R4P353

- Molecule 2 is a protein called Presegetalin A1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	W	7	Total	C	N	O	0	0	0
			48	28	8	12			
2	X	6	Total	C	N	O	0	0	0
			39	23	7	9			
2	Y	6	Total	C	N	O	0	0	0
			39	23	7	9			
2	Z	8	Total	C	N	O	0	0	0
			55	33	9	13			

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	190	Total O 190 190	0	0
5	B	220	Total O 220 220	0	0
5	C	238	Total O 238 238	0	0
5	D	178	Total O 178 178	0	0
5	W	5	Total O 5 5	0	0
5	X	3	Total O 3 3	0	0

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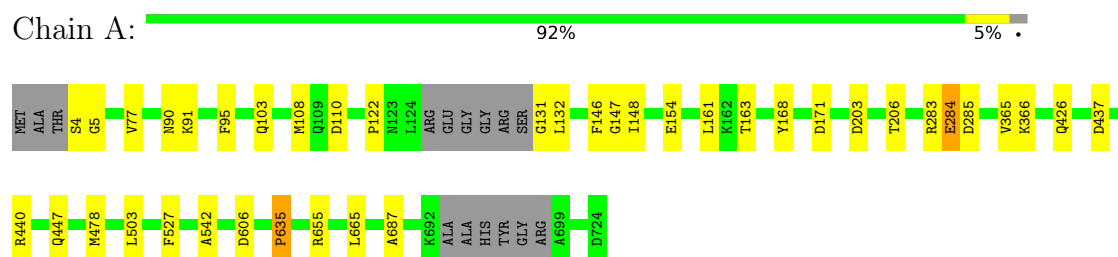
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	Y	5	Total	O	0	0
			5	5		
5	Z	5	Total	O	0	0
			5	5		

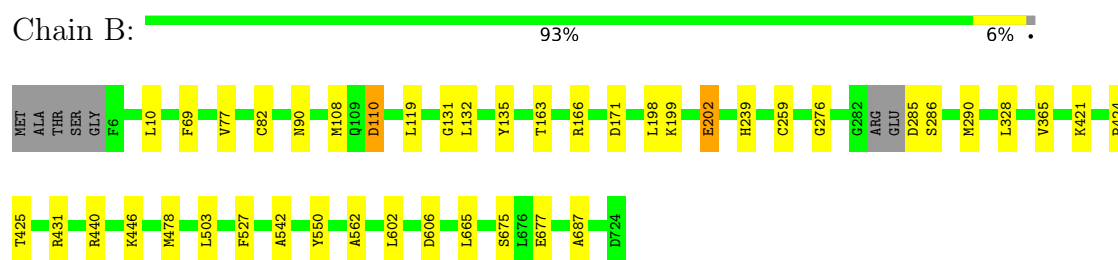
3 Residue-property plots

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.

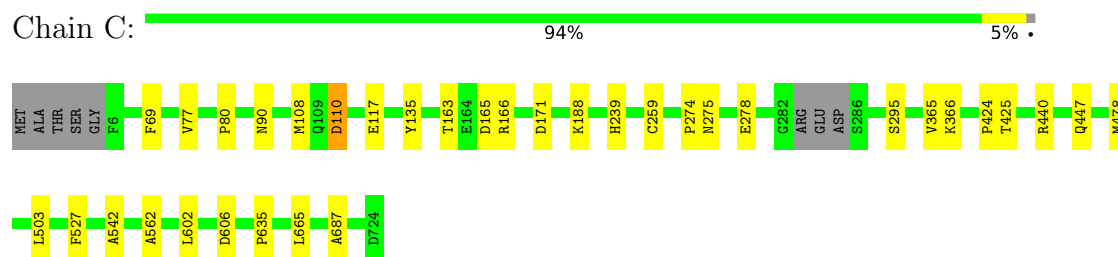
• Molecule 1: Peptide cyclase 1



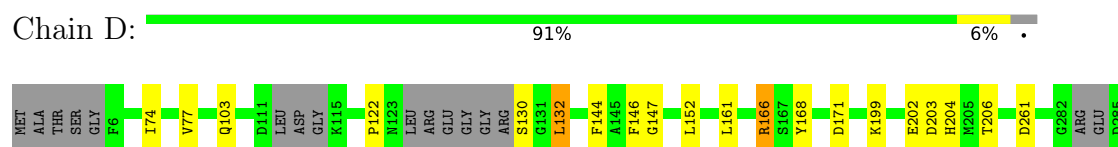
• Molecule 1: Peptide cyclase 1



• Molecule 1: Peptide cyclase 1

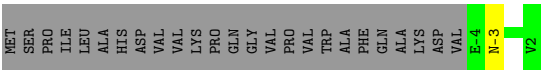


• Molecule 1: Peptide cyclase 1

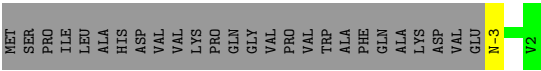




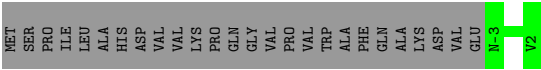
• Molecule 2: Presegetalin A1



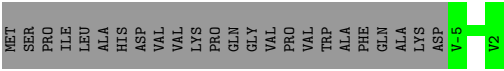
• Molecule 2: Presegetalin A1



• Molecule 2: Presegetalin A1



• Molecule 2: Presegetalin A1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.08Å 88.60Å 144.95Å 93.00° 99.32° 90.05°	Depositor
Resolution (Å)	77.07 – 2.00 77.07 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.0 (77.07-2.00) 97.0 (77.07-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.221 , 0.256 0.227 , 0.262	Depositor DCC
R_{free} test set	9515 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.690	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 25.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.085 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	23929	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: SO4, MG

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	0.89	0/5852	0.97	5/7923 (0.1%)
1	B	0.92	2/5928 (0.0%)	0.96	5/8025 (0.1%)
1	C	0.92	2/5920 (0.0%)	0.96	3/8014 (0.0%)
1	D	0.88	2/5798 (0.0%)	0.96	4/7848 (0.1%)
2	W	0.68	0/48	0.67	0/64
2	X	0.77	0/39	0.75	0/52
2	Y	0.63	0/39	0.79	0/52
2	Z	0.92	0/55	0.84	0/74
All	All	0.90	6/23679 (0.0%)	0.96	17/32052 (0.1%)

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	562	ALA	CA-C	6.62	1.59	1.52
1	C	562	ALA	CA-C	6.19	1.58	1.52
1	C	80	PRO	C-O	5.96	1.30	1.23
1	D	562	ALA	CA-C	5.58	1.58	1.52
1	D	261	ASP	C-O	5.14	1.30	1.23
1	B	131	GLY	N-CA	5.12	1.49	1.44

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	284	GLU	N-CA-C	-7.22	102.47	110.91
1	C	635	PRO	N-CA-C	6.97	123.92	113.81
1	D	635	PRO	N-CA-C	6.86	123.76	113.81
1	B	424	PRO	N-CA-C	6.82	126.52	112.47
1	A	635	PRO	N-CA-C	6.61	123.39	113.81
1	C	424	PRO	N-CA-C	6.37	125.59	112.47
1	B	677	GLU	N-CA-C	6.22	120.06	112.23
1	A	5	GLY	N-CA-C	5.91	120.04	112.83
1	A	147	GLY	N-CA-C	5.55	120.61	110.95
1	D	147	GLY	N-CA-C	5.46	120.46	110.95
1	B	425	THR	N-CA-C	5.32	119.62	113.18
1	A	285	ASP	N-CA-C	5.23	118.45	111.75
1	C	425	THR	N-CA-C	5.21	119.48	113.18
1	D	144	PHE	N-CA-C	5.11	117.39	108.76
1	B	285	ASP	CA-C-N	5.05	131.19	121.54
1	B	285	ASP	C-N-CA	5.05	131.19	121.54
1	D	132	LEU	N-CA-C	5.01	117.57	109.06

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	284	GLU	Peptide

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	5703	0	5543	21	0
1	B	5773	0	5615	26	0
1	C	5765	0	5611	22	0
1	D	5651	0	5490	21	0
2	W	48	0	45	1	0
2	X	39	0	39	1	0
2	Y	39	0	39	0	0
2	Z	55	0	54	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	A	190	0	0	0	0
5	B	220	0	0	2	0
5	C	238	0	0	1	0
5	D	178	0	0	1	0
5	W	5	0	0	0	0
5	X	3	0	0	0	0
5	Y	5	0	0	0	0
5	Z	5	0	0	0	0
All	All	23929	0	22436	81	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 2.

All (81) 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:B:421:LYS:CE	1:C:278:GLU:HG2	1.83	1.08
1:B:166[B]:ARG:HG3	1:B:166[B]:ARG:HH11	1.14	1.07
1:B:421:LYS:HE2	1:C:278:GLU:CG	1.90	1.01
1:B:446:LYS:NZ	1:B:550:TYR:OH	1.95	0.98
1:B:421:LYS:HE3	1:C:278:GLU:HG2	1.47	0.92
1:B:421:LYS:CE	1:C:278:GLU:CG	2.48	0.91
1:B:421:LYS:HE2	1:C:278:GLU:HG3	1.56	0.88
1:A:203:ASP:HB3	1:A:206:THR:HG22	1.58	0.85
1:A:146:PHE:HE2	1:A:148:ILE:HD11	1.43	0.82
1:B:166[B]:ARG:HG3	1:B:166[B]:ARG:NH1	1.89	0.82
1:B:421:LYS:HE2	1:C:278:GLU:HG2	1.55	0.81
1:D:290:MET:HE1	1:D:328:LEU:HD22	1.69	0.75
1:A:132:LEU:CD2	1:A:148:ILE:HD13	2.17	0.74
1:B:166[B]:ARG:HH11	1:B:166[B]:ARG:CG	1.98	0.73
1:C:295:SER:HB3	1:D:332:SER:HB2	1.70	0.72
1:A:146:PHE:CE2	1:A:148:ILE:HD11	2.23	0.71
1:B:290:MET:HE1	1:B:328:LEU:HD22	1.72	0.70
1:D:122:PRO:HB3	1:D:132:LEU:HD13	1.76	0.66
1:A:103:GLN:OE1	1:A:131:GLY:HA2	1.96	0.66
1:A:91:LYS:CD	1:A:108:MET:HE2	2.28	0.64
1:D:366:LYS:NZ	1:D:447:GLN:OE1	2.32	0.63
1:C:295:SER:HB3	1:D:332:SER:CB	2.29	0.61
1:A:132:LEU:CD2	1:A:148:ILE:CD1	2.78	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:ARG:CG	1:D:166:ARG:HH11	2.14	0.60
1:B:440:ARG:HD3	5:B:1095:HOH:O	2.01	0.60
1:A:132:LEU:HD21	1:A:148:ILE:CD1	2.33	0.59
1:B:82:CYS:SG	1:B:431:ARG:NH1	2.76	0.59
1:C:275:ASN:O	1:C:278:GLU:HB3	2.02	0.59
1:A:91:LYS:HD3	1:A:108:MET:HE2	1.83	0.58
1:D:166:ARG:HH11	1:D:166:ARG:HG3	1.69	0.57
1:A:122:PRO:HG3	1:A:132:LEU:HD12	1.86	0.57
1:A:108:MET:HE1	1:A:163:THR:HB	1.86	0.56
1:B:82:CYS:HB3	5:B:902:HOH:O	2.05	0.56
1:B:198:LEU:HG	1:B:202:GLU:HB3	1.88	0.54
1:A:90:ASN:O	1:A:110:ASP:O	2.25	0.53
1:A:122:PRO:CG	1:A:132:LEU:HD12	2.38	0.53
1:A:366:LYS:NZ	1:A:447:GLN:OE1	2.41	0.53
1:C:90:ASN:O	1:C:110:ASP:O	2.27	0.53
1:D:450:VAL:CG1	5:D:884:HOH:O	2.55	0.52
1:D:74:ILE:HD11	1:D:701:MET:HE1	1.92	0.52
1:C:366:LYS:NZ	1:C:447:GLN:OE1	2.42	0.52
1:A:122:PRO:HB3	1:A:132:LEU:CD1	2.41	0.51
1:B:90:ASN:O	1:B:110:ASP:O	2.30	0.50
1:C:108:MET:HE1	1:C:163:THR:HB	1.93	0.50
1:B:665:LEU:HD22	1:B:687:ALA:HB2	1.94	0.49
1:C:665:LEU:HD22	1:C:687:ALA:HB2	1.94	0.49
1:B:132:LEU:O	2:X:-3:ASN:ND2	2.39	0.49
1:C:117:GLU:OE1	1:C:166[B]:ARG:NH2	2.46	0.49
1:C:274:PRO:O	1:C:275:ASN:OD1	2.31	0.48
1:B:290:MET:HE2	1:B:328:LEU:HB3	1.96	0.48
1:C:188:LYS:HB2	5:C:924:HOH:O	2.13	0.48
1:D:290:MET:HE2	1:D:328:LEU:HB3	1.96	0.47
1:A:132:LEU:HD23	1:A:148:ILE:HD13	1.97	0.47
1:D:585:ASN:O	1:D:586[A]:CYS:SG	2.74	0.46
1:D:634:GLN:HB2	1:D:637:ASN:OD1	2.17	0.45
1:B:421:LYS:CE	1:C:278:GLU:HG3	2.31	0.45
1:C:90:ASN:O	1:C:90:ASN:CG	2.58	0.45
1:B:239:HIS:CD2	1:B:602:LEU:HD22	2.52	0.45
1:D:132:LEU:HD23	1:D:146:PHE:CD2	2.52	0.45
1:D:132:LEU:HD23	1:D:146:PHE:HD2	1.82	0.45
1:D:665:LEU:HD22	1:D:687:ALA:HB2	2.00	0.44
1:A:161:LEU:HD23	1:A:168:TYR:HA	2.01	0.43
1:D:161:LEU:HD23	1:D:168:TYR:HA	2.01	0.43
1:A:154:GLU:OE2	1:A:655:ARG:NH2	2.42	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:PRO:O	1:C:275:ASN:CG	2.62	0.43
1:D:103:GLN:NE2	1:D:130:SER:O	2.51	0.42
1:A:95:PHE:CE2	2:W:-3:ASN:HB2	2.55	0.42
1:B:478:MET:HE1	1:B:542:ALA:HB2	2.01	0.42
1:A:665:LEU:HD22	1:A:687:ALA:HB2	2.02	0.41
1:D:166:ARG:CG	1:D:166:ARG:NH1	2.80	0.41
1:D:203:ASP:HB3	1:D:206:THR:OG1	2.20	0.41
1:B:199:LYS:O	1:B:202:GLU:HB2	2.21	0.41
1:B:290:MET:CE	1:B:328:LEU:HB3	2.50	0.41
1:C:239:HIS:CD2	1:C:602:LEU:HD22	2.56	0.41
1:D:478:MET:HE1	1:D:542:ALA:HB2	2.02	0.41
1:A:478:MET:HE1	1:A:542:ALA:HB2	2.03	0.40
1:C:478:MET:HE1	1:C:542:ALA:HB2	2.02	0.40
1:B:108:MET:HE1	1:B:163:THR:HB	2.02	0.40
1:B:119:LEU:HD12	1:B:166[A]:ARG:CD	2.51	0.40
1:C:165:ASP:C	1:C:166[B]:ARG:HD2	2.47	0.40
1:D:199:LYS:O	1:D:202:GLU:HG2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	704/724 (97%)	674 (96%)	29 (4%)	1 (0%)	48	46
1	B	715/724 (99%)	684 (96%)	29 (4%)	2 (0%)	36	35
1	C	714/724 (99%)	684 (96%)	30 (4%)	0	100	100
1	D	693/724 (96%)	667 (96%)	26 (4%)	0	100	100
2	W	5/32 (16%)	4 (80%)	1 (20%)	0	100	100
2	X	4/32 (12%)	4 (100%)	0	0	100	100
2	Y	4/32 (12%)	4 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	Z	6/32 (19%)	6 (100%)	0	0	100	100
All	All	2845/3024 (94%)	2727 (96%)	115 (4%)	3 (0%)	48	46

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	283	ARG
1	B	286	SER
1	B	276	GLY

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	620/628 (99%)	609 (98%)	11 (2%)	51	58
1	B	625/628 (100%)	612 (98%)	13 (2%)	47	52
1	C	624/628 (99%)	613 (98%)	11 (2%)	51	58
1	D	615/628 (98%)	601 (98%)	14 (2%)	44	49
2	W	5/26 (19%)	5 (100%)	0	100	100
2	X	4/26 (15%)	4 (100%)	0	100	100
2	Y	4/26 (15%)	4 (100%)	0	100	100
2	Z	6/26 (23%)	6 (100%)	0	100	100
All	All	2503/2616 (96%)	2454 (98%)	49 (2%)	48	54

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	77	VAL
1	A	171	ASP
1	A	365	VAL
1	A	426	GLN

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Mol	Chain	Res	Type
1	A	437	ASP
1	A	440	ARG
1	A	503	LEU
1	A	527	PHE
1	A	606	ASP
1	A	635	PRO
1	B	10	LEU
1	B	69	PHE
1	B	77	VAL
1	B	110	ASP
1	B	135	TYR
1	B	171	ASP
1	B	202	GLU
1	B	259	CYS
1	B	365	VAL
1	B	503	LEU
1	B	527	PHE
1	B	606	ASP
1	B	675	SER
1	C	69	PHE
1	C	77	VAL
1	C	110	ASP
1	C	135	TYR
1	C	171	ASP
1	C	259	CYS
1	C	365	VAL
1	C	440	ARG
1	C	503	LEU
1	C	527	PHE
1	C	606	ASP
1	D	77	VAL
1	D	152	LEU
1	D	166	ARG
1	D	171	ASP
1	D	204	HIS
1	D	365	VAL
1	D	426	GLN
1	D	437	ASP
1	D	440	ARG
1	D	503	LEU
1	D	527	PHE
1	D	606	ASP

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Mol	Chain	Res	Type
1	D	635	PRO
1	D	678	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	ASN
1	A	353	ASN
1	A	356	GLN
1	A	422	ASN
1	A	426	GLN
1	A	532	GLN
1	A	574	GLN
1	A	595	HIS
1	B	41	ASN
1	B	71	GLN
1	B	90	ASN
1	B	94	HIS
1	B	107	GLN
1	B	353	ASN
1	B	356	GLN
1	B	422	ASN
1	B	532	GLN
1	B	574	GLN
1	C	235	ASN
1	C	353	ASN
1	C	356	GLN
1	C	417	GLN
1	C	422	ASN
1	C	532	GLN
1	C	574	GLN
1	D	90	ASN
1	D	235	ASN
1	D	353	ASN
1	D	356	GLN
1	D	417	GLN
1	D	422	ASN
1	D	426	GLN
1	D	532	GLN
1	D	574	GLN
1	D	634	GLN
1	D	670	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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	$\# Z > 2$
4	SO4	C	802	-	4,4,4	0.33	0	6,6,6	0.31	0
4	SO4	B	802	-	4,4,4	0.31	0	6,6,6	0.23	0

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 ⓘ

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	709/724 (97%)	-1.12	0 100 100	20, 36, 69, 128	1 (0%)
1	B	717/724 (99%)	-1.21	0 100 100	16, 33, 65, 118	2 (0%)
1	C	716/724 (98%)	-1.18	0 100 100	16, 33, 71, 130	2 (0%)
1	D	702/724 (96%)	-1.12	0 100 100	18, 36, 67, 107	1 (0%)
2	W	7/32 (21%)	-1.09	0 100 100	37, 41, 50, 52	0
2	X	6/32 (18%)	-1.29	0 100 100	31, 33, 40, 40	0
2	Y	6/32 (18%)	-1.28	0 100 100	29, 30, 37, 41	0
2	Z	8/32 (25%)	-0.93	0 100 100	35, 43, 47, 47	0
All	All	2871/3024 (94%)	-1.16	0 100 100	16, 35, 68, 130	6 (0%)

There are no RSRZ outliers to report.

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	MG	C	801	1/1	0.99	0.04	35,35,35,35	0
4	SO4	B	802	5/5	0.99	0.04	37,37,38,38	0
3	MG	B	801	1/1	1.00	0.03	38,38,38,38	0
4	SO4	C	802	5/5	1.00	0.04	36,36,37,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.