Full wwPDB X-ray Structure Validation Report (i)

Apr 21, 2024 – 11:52 am BST

PDB ID	:	5O3W
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 (i) symbol.

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

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

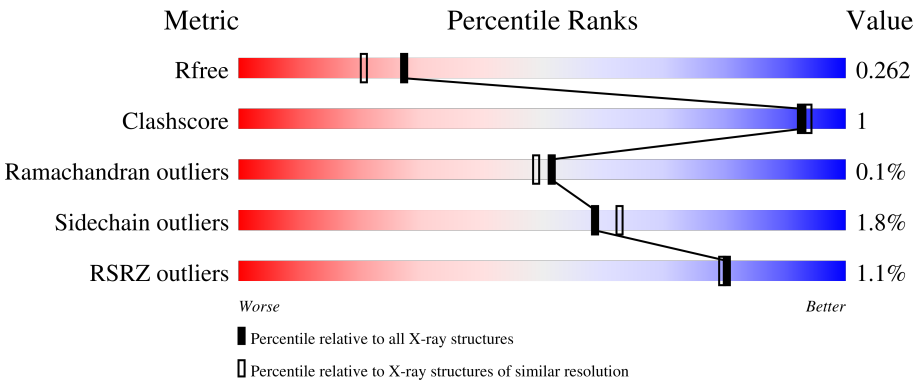
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 <=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	.% 93% 5% •
1	B	724	2% 94% 5% •
1	C	724	.% 95% • •
1	D	724	.% 92% • •
2	W	32	19% • 78%

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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 (i)

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

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

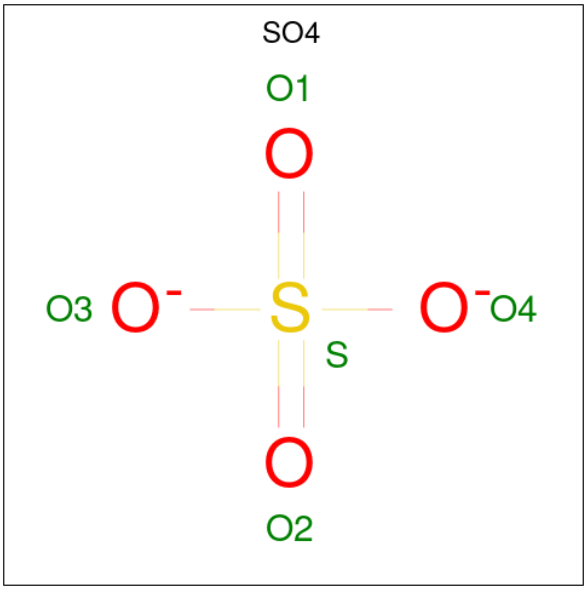
- Molecule 2 is a protein called Presegetalin A1.

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

- Molecule 3 is MAGNESIUM ION (three-letter code: 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 (three-letter code: SO4) (formula: O_4S).



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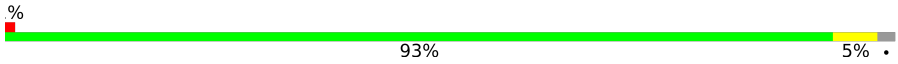
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	Y	5	5	5	0	0
			Total	O		
5	Z	5	5	5	0	0

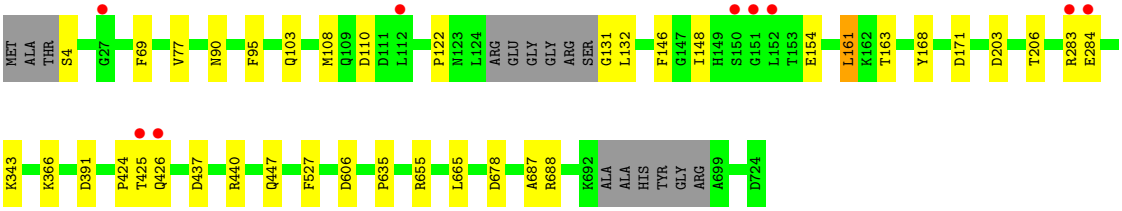
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Peptide cyclase 1

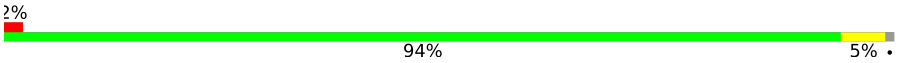
Chain A:

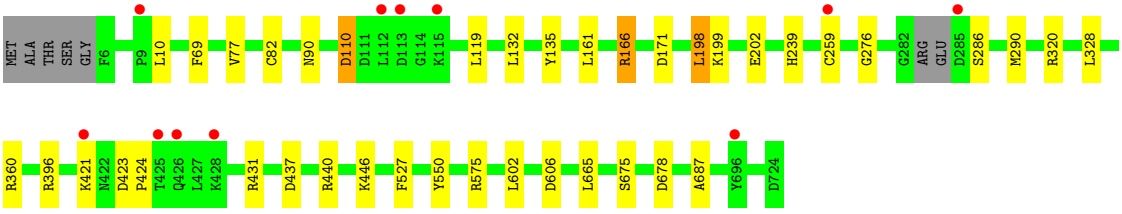




• Molecule 1: Peptide cyclase 1

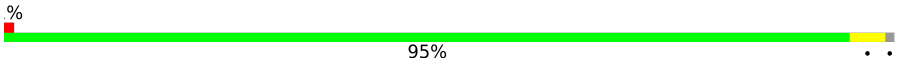
Chain B:





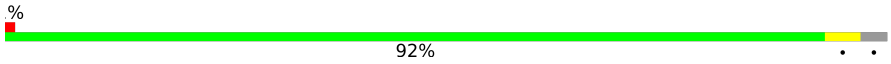
• Molecule 1: Peptide cyclase 1

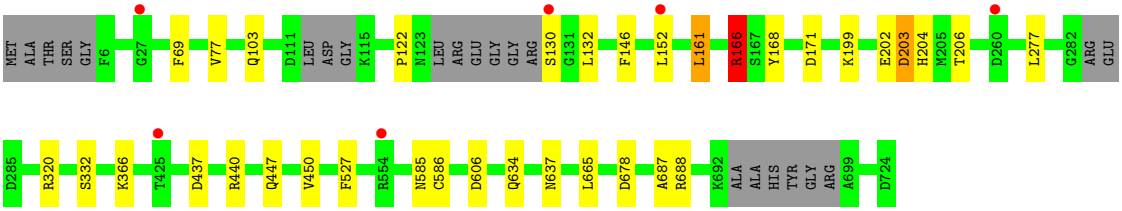
Chain C:



• Molecule 1: Peptide cyclase 1

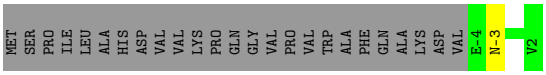
Chain D:





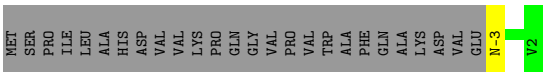
- Molecule 2: Presegetalin A1

Chain W:



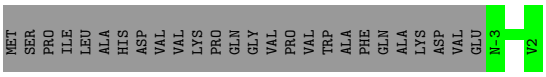
- Molecule 2: Presegetalin A1

Chain X:



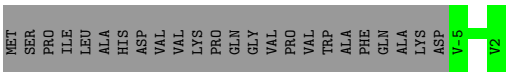
- Molecule 2: Presegetalin A1

Chain Y:



- Molecule 2: Presegetalin A1

Chain Z:



4 Data and refinement statistics (i)

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
$< I/\sigma(I) >$ [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)	0.31 , 25.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.085 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.94	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.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.67	0/5852	0.84	4/7923 (0.1%)
1	B	0.70	0/5928	0.85	10/8025 (0.1%)
1	C	0.70	0/5920	0.86	7/8014 (0.1%)
1	D	0.67	0/5798	0.86	7/7848 (0.1%)
2	W	0.60	0/48	0.59	0/64
2	X	0.60	0/39	0.55	0/52
2	Y	0.54	0/39	0.53	0/52
2	Z	0.73	0/55	0.65	0/74
All	All	0.69	0/23679	0.85	28/32052 (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

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	166	ARG	NE-CZ-NH1	8.41	124.50	120.30
1	B	166[A]	ARG	CG-CD-NE	-7.33	96.40	111.80
1	B	166[B]	ARG	CG-CD-NE	-7.33	96.40	111.80
1	D	166	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	D	203	ASP	CB-CG-OD1	6.25	123.92	118.30
1	D	203	ASP	CB-CG-OD2	-5.85	113.03	118.30
1	C	575	ARG	NE-CZ-NH1	5.83	123.22	120.30
1	B	161	LEU	CA-CB-CG	5.74	128.50	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	81	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	B	320	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	B	424	PRO	N-CA-C	5.55	126.52	112.10
1	A	161	LEU	CA-CB-CG	5.51	127.98	115.30
1	D	161	LEU	CA-CB-CG	5.51	127.96	115.30
1	A	343	LYS	CD-CE-NZ	5.50	124.36	111.70
1	C	161	LEU	CA-CB-CG	5.50	127.94	115.30
1	B	360	ARG	NE-CZ-NH1	5.41	123.01	120.30
1	B	198	LEU	CA-CB-CG	5.39	127.70	115.30
1	B	396	ARG	NE-CZ-NH2	-5.29	117.66	120.30
1	A	688	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	D	688	ARG	NE-CZ-NH1	5.19	122.89	120.30
1	C	424	PRO	N-CA-C	5.19	125.59	112.10
1	C	360	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	D	320	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	A	391	ASP	CB-CG-OD1	-5.09	113.72	118.30
1	C	320	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	C	688	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	B	575	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	B	437	ASP	CB-CG-OD1	5.05	122.85	118.30

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 (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	5703	0	5543	19	0
1	B	5773	0	5615	23	0
1	C	5765	0	5611	18	0
1	D	5651	0	5490	17	0
2	W	48	0	45	1	0
2	X	39	0	39	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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
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	68	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.

All (68) 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: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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LEU:CD2	1:A:148:ILE:CD1	2.78	0.60
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: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: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: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:B:290:MET:HE1	1:B:328:LEU:HB3	1.94	0.49
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:A:424:PRO:O	1:A:425:THR:OG1	2.29	0.49
1:C:274:PRO:O	1:C:275:ASN:OD1	2.31	0.48
1:C:188:LYS:HB2	5:C:924:HOH:O	2.13	0.48
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: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
1:D:103:GLN:NE2	1:D:130:SER:O	2.51	0.42
1:A:108:MET:HE3	1:A:163:THR:HB	2.01	0.42
1:A:95:PHE:CE2	2:W:-3:ASN:HB2	2.55	0.42
1:C:90:ASN:O	1:C:90:ASN:CG	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:119:LEU:HD12	1:B:166[A]:ARG:CD	2.51	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 (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	704/724 (97%)	674 (96%)	29 (4%)	1 (0%)	51	49
1	B	715/724 (99%)	684 (96%)	29 (4%)	2 (0%)	41	37
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
2	Z	6/32 (19%)	6 (100%)	0	0	100	100
All	All	2845/3024 (94%)	2727 (96%)	115 (4%)	3 (0%)	51	49

All (3) Ramachandran outliers are listed below:

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

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Mol	Chain	Res	Type
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%)	59	63
1	B	625/628 (100%)	613 (98%)	12 (2%)	57	61
1	C	624/628 (99%)	614 (98%)	10 (2%)	62	67
1	D	615/628 (98%)	603 (98%)	12 (2%)	55	58
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%)	2458 (98%)	45 (2%)	59	63

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	69	PHE
1	A	77	VAL
1	A	171	ASP
1	A	426	GLN
1	A	437	ASP
1	A	440	ARG
1	A	527	PHE
1	A	606	ASP
1	A	635	PRO
1	A	678	ASP
1	B	10	LEU
1	B	69	PHE
1	B	77	VAL

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Mol	Chain	Res	Type
1	B	110	ASP
1	B	135	TYR
1	B	171	ASP
1	B	259	CYS
1	B	423	ASP
1	B	527	PHE
1	B	606	ASP
1	B	675	SER
1	B	678	ASP
1	C	69	PHE
1	C	77	VAL
1	C	110	ASP
1	C	135	TYR
1	C	171	ASP
1	C	185	HIS
1	C	259	CYS
1	C	440	ARG
1	C	527	PHE
1	C	606	ASP
1	D	69	PHE
1	D	77	VAL
1	D	152	LEU
1	D	166	ARG
1	D	171	ASP
1	D	204	HIS
1	D	277	LEU
1	D	437	ASP
1	D	440	ARG
1	D	527	PHE
1	D	606	ASP
1	D	678	ASP

Sometimes 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 monosaccharides 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	0.22	0	6,6,6	0.39	0
4	SO4	B	802	-	4,4,4	0.20	0	6,6,6	0.25	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 (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	709/724 (97%)	-0.31	9 (1%)	77	76	20, 36, 69, 128	0
1	B	717/724 (99%)	-0.38	11 (1%)	73	72	16, 33, 65, 118	0
1	C	716/724 (98%)	-0.37	5 (0%)	87	87	16, 33, 71, 130	0
1	D	702/724 (96%)	-0.33	6 (0%)	84	83	21, 36, 67, 107	0
2	W	7/32 (21%)	0.08	0	100	100	37, 41, 50, 52	0
2	X	6/32 (18%)	-0.21	0	100	100	31, 33, 40, 40	0
2	Y	6/32 (18%)	-0.43	0	100	100	29, 30, 37, 41	0
2	Z	8/32 (25%)	-0.06	0	100	100	35, 43, 47, 47	0
All	All	2871/3024 (94%)	-0.35	31 (1%)	80	79	16, 35, 68, 130	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	GLY	6.7
1	D	152	LEU	6.1
1	A	152	LEU	5.6
1	A	112	LEU	4.2
1	B	9	PRO	4.1
1	D	425	THR	3.9
1	C	286	SER	3.9
1	A	425	THR	3.7
1	B	421	LYS	3.4
1	B	112	LEU	3.3
1	A	284	GLU	3.0
1	B	426	GLN	3.0
1	B	113	ASP	2.9
1	C	426	GLN	2.9
1	A	150	SER	2.9
1	D	27	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	425	THR	2.7
1	C	113	ASP	2.6
1	D	554	ARG	2.6
1	D	260	ASP	2.6
1	A	426	GLN	2.6
1	A	27	GLY	2.4
1	B	115	LYS	2.4
1	A	283	ARG	2.3
1	B	259	CYS	2.3
1	B	696	TYR	2.2
1	C	425	THR	2.2
1	D	130	SER	2.2
1	B	285	ASP	2.2
1	C	287	ALA	2.1
1	B	428	LYS	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MG	C	801	1/1	0.98	0.05	35,35,35,35	0
4	SO4	B	802	5/5	0.98	0.06	37,37,38,38	0
4	SO4	C	802	5/5	0.98	0.07	36,36,37,38	0
3	MG	B	801	1/1	0.99	0.06	38,38,38,38	0

6.5 Other polymers (i)

There are no such residues in this entry.