



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:23 AM GMT

PDB ID : 3CBI
Title : Crystal structure of the ternary complex of phospholipase A2 with ajmaline and anisic acid at 3.1 Å resolution
Authors : Kumar, S.; Vikram, G.; Singh, N.; Sharma, S.; Kaur, P.; Singh, T.P.
Deposited on : 2008-02-22
Resolution : 3.15 Å(reported)

This is a full wwPDB validation 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/ValidationPDFNotes.html>

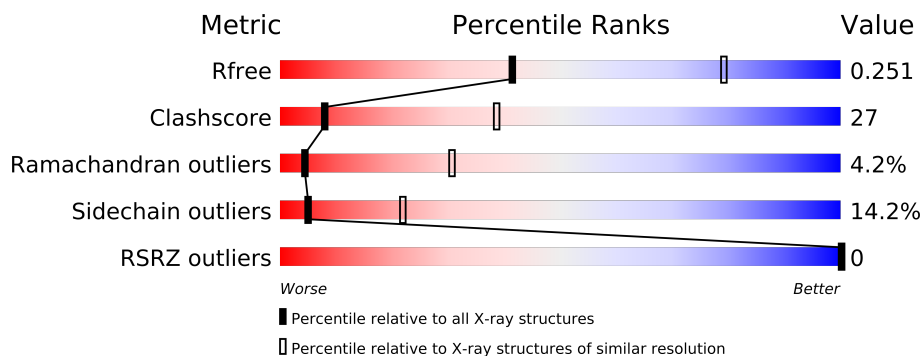
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.15 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}	66092	1360 (3.22-3.10)
Clashscore	79885	1681 (3.22-3.10)
Ramachandran outliers	78287	1639 (3.22-3.10)
Sidechain outliers	78261	1638 (3.22-3.10)
RSRZ outliers	66119	1361 (3.22-3.10)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	121	<div><div></div><div></div><div></div><div></div><div></div></div>
1	B	121	<div><div></div><div></div><div></div><div></div><div></div></div>
1	C	121	<div><div></div><div></div><div></div><div></div><div></div></div>
1	D	121	<div><div></div><div></div><div></div><div></div><div></div></div>

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	ANN	A	349	-	X
3	ANN	B	348	-	X
3	ANN	C	347	-	X
3	ANN	D	350	-	X

2 Entry composition i

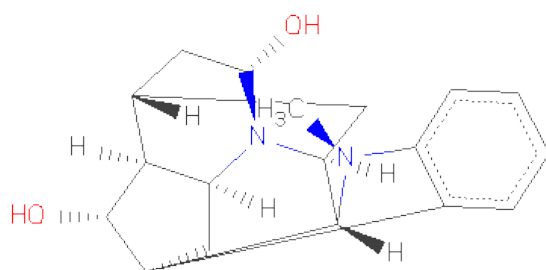
There are 3 unique types of molecules in this entry. The entry contains 3904 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Phospholipase A2 VRV-PL-VIIIa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	121	Total	C	N	O	S	0	0	0
			943	591	159	177	16			
1	B	121	Total	C	N	O	S	0	0	0
			943	591	159	177	16			
1	C	121	Total	C	N	O	S	0	0	0
			943	591	159	177	16			
1	D	121	Total	C	N	O	S	0	0	0
			943	591	159	177	16			

- Molecule 2 is AJMALINE (three-letter code: AJM) (formula: C₁₈H₂₂N₂O₂).



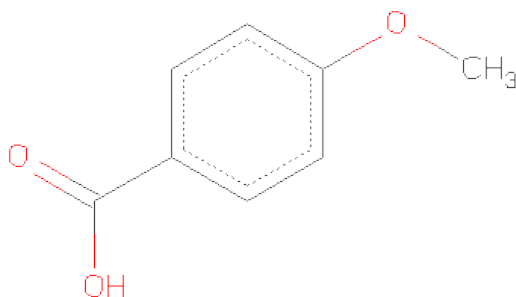
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			22	18	2	2		
2	B	1	Total	C	N	O	0	0
			22	18	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			22	18	2	2		
2	D	1	Total	C	N	O	0	0
			22	18	2	2		

- Molecule 3 is 4-METHOXYBENZOIC ACID (three-letter code: ANN) (formula: C₈H₈O₃).

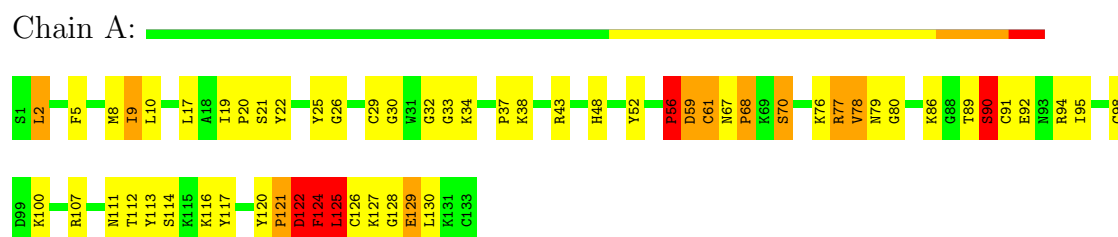


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			11	8	3		
3	B	1	Total	C	O	0	0
			11	8	3		
3	A	1	Total	C	O	0	0
			11	8	3		
3	D	1	Total	C	O	0	0
			11	8	3		

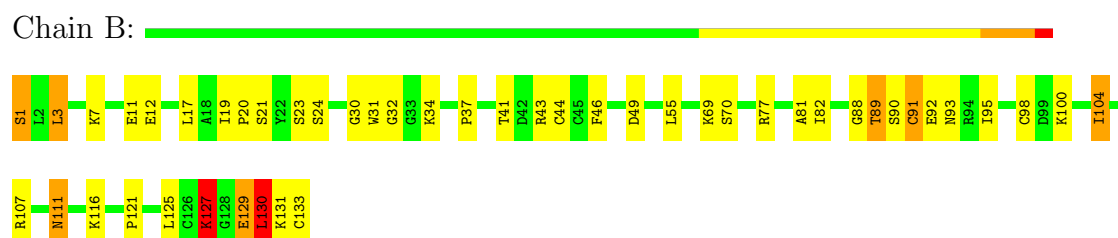
3 Residue-property plots

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.

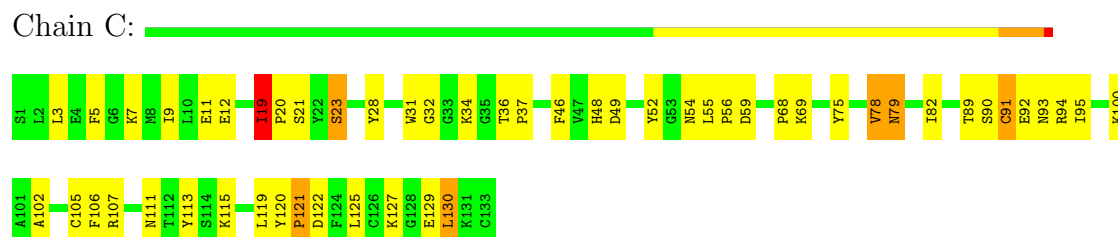
• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



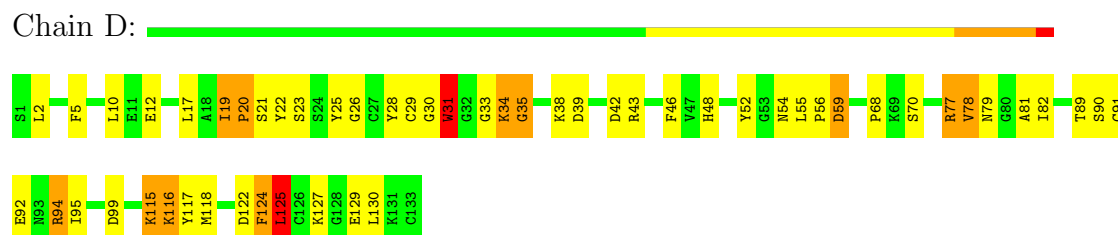
• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	108.80Å 108.80Å 43.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 3.15 23.04 – 3.15	Depositor EDS
% Data completeness (in resolution range)	92.2 (50.00-3.15) 97.5 (23.04-3.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 3.16Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.228 , 0.250 0.228 , 0.251	Depositor DCC
R_{free} test set	513 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.542	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 52.8	EDS
Estimated twinning fraction	0.420 for -h,-k,l 0.024 for h,-h-k,-l 0.017 for -k,-h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 9673 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3904	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.60	0/963	0.98	1/1292 (0.1%)
1	B	0.63	0/963	0.92	2/1292 (0.2%)
1	C	0.59	0/963	0.94	1/1292 (0.1%)
1	D	0.67	0/963	1.04	8/1292 (0.6%)
All	All	0.62	0/3852	0.97	12/5168 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	59	ASP	CB-CG-OD1	7.80	125.32	118.30
1	B	130	LEU	N-CA-C	-6.96	92.21	111.00
1	A	128	GLY	N-CA-C	6.57	129.51	113.10
1	C	19	ILE	C-N-CD	6.55	142.17	128.40
1	D	35	GLY	N-CA-C	6.25	128.72	113.10
1	D	19	ILE	N-CA-C	-6.13	94.45	111.00
1	D	59	ASP	CA-C-N	-5.98	104.05	117.20
1	D	59	ASP	O-C-N	5.88	132.11	122.70
1	D	59	ASP	CB-CG-OD2	-5.45	113.40	118.30
1	D	31	TRP	CA-CB-CG	-5.18	103.87	113.70
1	D	34	LYS	N-CA-C	5.13	124.85	111.00
1	B	88	GLY	N-CA-C	-5.12	100.29	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	943	0	907	64	0
1	B	943	0	907	44	0
1	C	943	0	907	61	0
1	D	943	0	907	42	0
2	A	22	0	22	6	0
2	B	22	0	22	3	0
2	C	22	0	22	7	0
2	D	22	0	22	0	0
3	A	11	0	7	2	0
3	B	11	0	7	2	0
3	C	11	0	7	4	0
3	D	11	0	7	2	0
All	All	3904	0	3744	206	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 27.

All (206) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:76:LYS:HE3	1:A:78:VAL:HG21	1.44	0.96
1:A:90:SER:O	1:A:94:ARG:HG3	1.68	0.93
1:B:31:TRP:O	2:B:202:AJM:H13	1.71	0.91
1:A:76:LYS:HE3	1:A:78:VAL:CG2	2.01	0.90
1:C:49:ASP:HB3	2:C:201:AJM:C16	2.04	0.88
1:C:37:PRO:HD3	1:C:130:LEU:CD1	2.04	0.88
1:C:49:ASP:HB3	2:C:201:AJM:H162	1.57	0.86
1:A:32:GLY:HA3	2:A:203:AJM:H161	1.60	0.83
1:C:37:PRO:HD3	1:C:130:LEU:HD13	1.61	0.83
1:C:121:PRO:HG2	1:C:125:LEU:HD22	1.58	0.82
1:B:91:CYS:O	1:B:95:ILE:HG12	1.79	0.82
1:C:32:GLY:HA3	2:C:201:AJM:H122	1.60	0.81
1:B:100:LYS:O	1:B:104:ILE:HD13	1.79	0.81
1:C:68:PRO:HB3	1:C:95:ILE:HD12	1.64	0.79
1:D:17:LEU:O	1:D:21:SER:HB2	1.81	0.79

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:10:LEU:HD13	1:A:17:LEU:HD23	1.66	0.78
1:B:90:SER:HA	1:B:93:ASN:HB2	1.65	0.77
1:A:30:GLY:HA2	3:A:349:ANN:C2	2.15	0.77
1:A:17:LEU:O	1:A:21:SER:HB2	1.83	0.77
1:C:7:LYS:O	1:C:11:GLU:HG3	1.86	0.76
1:A:77:ARG:O	1:A:78:VAL:HB	1.86	0.75
1:C:32:GLY:CA	2:C:201:AJM:H122	2.16	0.74
1:D:48:HIS:CE1	1:D:52:TYR:CZ	2.76	0.74
1:C:90:SER:HA	1:C:93:ASN:ND2	2.03	0.73
1:B:89:THR:OG1	1:B:92:GLU:HG3	1.88	0.73
1:A:120:TYR:OH	1:A:125:LEU:HD12	1.88	0.73
1:A:89:THR:H	1:A:92:GLU:HB2	1.54	0.72
1:B:82:ILE:HD12	1:B:104:ILE:HD11	1.71	0.72
1:C:78:VAL:O	1:C:78:VAL:HG13	1.90	0.72
1:A:89:THR:O	1:A:91:CYS:N	2.23	0.71
1:B:37:PRO:HD3	1:B:130:LEU:HD12	1.71	0.71
1:C:19:ILE:HA	1:C:23:SER:HB3	1.73	0.70
1:C:46:PHE:CE2	1:D:78:VAL:HG11	2.27	0.70
1:D:55:LEU:HD11	1:D:94:ARG:HB3	1.74	0.70
1:C:12:GLU:OE1	1:C:107:ARG:HD3	1.91	0.69
1:D:55:LEU:CD1	1:D:94:ARG:HB3	2.22	0.69
1:C:36:THR:HA	1:C:130:LEU:HD12	1.73	0.69
1:C:36:THR:HA	1:C:130:LEU:CD1	2.23	0.69
1:B:24:SER:CB	1:B:31:TRP:HE1	2.05	0.68
1:C:54:ASN:O	1:C:56:PRO:HD3	1.93	0.68
1:C:9:ILE:HD12	1:C:106:PHE:HD2	1.58	0.68
1:A:78:VAL:HG13	1:B:130:LEU:HD21	1.77	0.66
1:C:48:HIS:CD2	1:C:102:ALA:HB2	2.30	0.66
1:D:68:PRO:HB3	1:D:95:ILE:HD12	1.76	0.66
1:A:121:PRO:CB	1:A:124:PHE:CZ	2.79	0.66
1:D:19:ILE:HA	1:D:23:SER:OG	1.95	0.66
1:A:121:PRO:HB2	1:A:124:PHE:CZ	2.31	0.66
1:C:23:SER:HB2	3:C:347:ANN:H6	1.77	0.66
1:C:12:GLU:O	1:C:107:ARG:HG3	1.95	0.66
1:A:5:PHE:CE1	1:A:9:ILE:HD11	2.31	0.65
1:B:31:TRP:O	2:B:202:AJM:C13	2.46	0.64
1:B:24:SER:HB2	1:B:31:TRP:HE1	1.62	0.64
1:A:5:PHE:O	1:A:9:ILE:HD13	1.98	0.64
1:C:130:LEU:HD22	1:D:78:VAL:HG13	1.79	0.64
1:C:9:ILE:HD12	1:C:106:PHE:CD2	2.32	0.64
1:B:55:LEU:CD1	1:B:95:ILE:HD13	2.28	0.63
1:A:33:GLY:H	2:A:203:AJM:H122	1.64	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:7:LYS:O	1:B:11:GLU:HG3	1.98	0.62
1:C:75:TYR:CD1	1:C:82:ILE:HG23	2.33	0.62
1:D:19:ILE:HD13	1:D:23:SER:CB	2.28	0.62
1:D:116:LYS:HE2	1:D:117:TYR:CE2	2.34	0.62
1:B:55:LEU:HD13	1:B:95:ILE:HD13	1.82	0.61
1:B:11:GLU:HB2	1:B:77:ARG:NH2	2.15	0.61
1:B:111:ASN:H	1:B:111:ASN:ND2	1.98	0.61
1:C:90:SER:HA	1:C:93:ASN:HD22	1.66	0.61
1:C:23:SER:HA	3:C:347:ANN:H6	1.83	0.61
1:A:76:LYS:CE	1:A:78:VAL:HG21	2.27	0.60
1:A:124:PHE:O	1:A:125:LEU:HB2	2.01	0.60
1:C:19:ILE:HG22	1:C:20:PRO:HD3	1.84	0.60
1:B:17:LEU:O	1:B:21:SER:HB2	2.02	0.60
1:C:55:LEU:CD1	1:C:95:ILE:HD13	2.33	0.59
1:C:48:HIS:HD2	1:C:102:ALA:HB2	1.69	0.58
1:C:68:PRO:HB3	1:C:95:ILE:CD1	2.34	0.58
1:A:121:PRO:HB3	1:A:124:PHE:CZ	2.38	0.57
1:A:25:TYR:CE2	1:A:113:TYR:CE2	2.91	0.57
1:C:49:ASP:HB3	2:C:201:AJM:H161	1.85	0.57
1:C:19:ILE:O	1:C:23:SER:HB3	2.05	0.57
1:D:20:PRO:HA	1:D:23:SER:OG	2.03	0.57
1:D:89:THR:OG1	1:D:92:GLU:HG3	2.04	0.57
1:C:78:VAL:O	1:C:79:ASN:HB2	2.05	0.57
1:A:9:ILE:HD13	1:A:9:ILE:H	1.69	0.57
1:C:32:GLY:C	2:C:201:AJM:H122	2.24	0.57
1:D:19:ILE:HD13	1:D:23:SER:HB2	1.87	0.56
1:C:75:TYR:HD1	1:C:82:ILE:HG23	1.69	0.56
1:C:5:PHE:CE1	1:C:9:ILE:HD11	2.41	0.55
1:B:41:THR:O	1:B:44:CYS:HB2	2.06	0.55
1:D:68:PRO:HB3	1:D:95:ILE:CD1	2.36	0.55
1:B:11:GLU:CB	1:B:77:ARG:HH21	2.20	0.55
1:A:121:PRO:CB	1:A:124:PHE:CE2	2.89	0.55
1:D:125:LEU:HD23	1:D:127:LYS:HE2	1.89	0.54
1:D:115:LYS:HA	1:D:118:MET:HB2	1.90	0.54
1:C:55:LEU:HD21	1:C:94:ARG:CB	2.38	0.54
1:A:122:ASP:O	1:A:125:LEU:N	2.41	0.54
1:D:124:PHE:O	1:D:125:LEU:HB2	2.08	0.54
1:B:1:SER:OG	1:B:3:LEU:HB2	2.07	0.54
1:C:32:GLY:HA3	2:C:201:AJM:C12	2.34	0.53
1:B:90:SER:O	1:B:91:CYS:C	2.46	0.53
1:A:90:SER:C	1:A:94:ARG:HG3	2.29	0.53
1:B:127:LYS:H	1:B:127:LYS:HD2	1.72	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:78:VAL:O	1:A:80:GLY:N	2.42	0.53
1:C:52:TYR:CE2	1:C:68:PRO:HB2	2.44	0.53
1:A:37:PRO:HD3	1:A:130:LEU:HD12	1.90	0.53
1:B:130:LEU:HD22	1:B:131:LYS:H	1.74	0.53
1:C:130:LEU:CD2	1:D:78:VAL:HG13	2.39	0.52
1:A:56:PRO:CD	1:A:59:ASP:N	2.69	0.52
1:B:11:GLU:HB2	1:B:77:ARG:HH21	1.74	0.52
1:B:23:SER:HB3	3:B:348:ANN:H6	1.91	0.52
1:A:76:LYS:HE3	1:A:78:VAL:HG23	1.88	0.51
1:D:48:HIS:HE1	1:D:52:TYR:CZ	2.24	0.51
1:D:48:HIS:CE1	1:D:52:TYR:CE1	2.98	0.51
1:C:5:PHE:O	1:C:9:ILE:HG12	2.11	0.51
1:B:23:SER:CB	3:B:348:ANN:H6	2.41	0.51
1:A:121:PRO:HB2	1:A:124:PHE:CE2	2.44	0.51
1:C:91:CYS:O	1:C:95:ILE:HG12	2.11	0.51
1:A:90:SER:O	1:A:94:ARG:CG	2.50	0.51
1:C:121:PRO:HB2	1:C:125:LEU:HD13	1.94	0.50
1:A:9:ILE:N	1:A:9:ILE:CD1	2.75	0.49
1:A:120:TYR:HH	1:A:125:LEU:HD12	1.73	0.49
1:D:54:ASN:O	1:D:56:PRO:HD3	2.13	0.48
1:B:90:SER:O	1:B:93:ASN:HB2	2.13	0.48
1:A:122:ASP:O	1:A:126:CYS:N	2.40	0.48
1:A:32:GLY:CA	2:A:203:AJM:H161	2.40	0.48
1:C:68:PRO:CB	1:C:95:ILE:HD12	2.38	0.48
1:C:23:SER:CA	3:C:347:ANN:H6	2.42	0.48
1:C:19:ILE:HG22	1:C:20:PRO:CD	2.43	0.48
1:B:129:GLU:HG3	1:B:130:LEU:O	2.14	0.48
1:D:52:TYR:CE2	1:D:68:PRO:HB2	2.48	0.48
1:A:22:TYR:O	1:A:29:CYS:HB3	2.14	0.48
1:A:77:ARG:O	1:A:78:VAL:CB	2.60	0.48
1:A:37:PRO:HD3	1:A:130:LEU:CD1	2.44	0.47
1:A:33:GLY:N	2:A:203:AJM:H122	2.28	0.47
1:D:77:ARG:O	1:D:81:ALA:O	2.32	0.47
1:C:119:LEU:HD23	1:C:119:LEU:N	2.29	0.47
1:C:56:PRO:HD2	1:C:59:ASP:H	1.79	0.47
1:B:11:GLU:CB	1:B:77:ARG:NH2	2.76	0.47
1:A:30:GLY:HA2	3:A:349:ANN:C7	2.43	0.47
1:B:12:GLU:OE1	1:B:107:ARG:HD3	2.14	0.47
1:C:55:LEU:HD21	1:C:94:ARG:HB2	1.95	0.47
1:B:90:SER:O	1:B:93:ASN:N	2.49	0.46
1:A:89:THR:OG1	1:A:92:GLU:HG3	2.15	0.46
1:A:2:LEU:HA	1:A:2:LEU:HD12	1.75	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:120:TYR:HA	1:C:121:PRO:HD2	1.83	0.46
1:A:76:LYS:CE	1:A:78:VAL:CG2	2.83	0.46
1:D:43:ARG:O	1:D:46:PHE:HB3	2.16	0.46
1:C:19:ILE:CG2	1:C:20:PRO:HD3	2.44	0.46
1:C:102:ALA:O	1:C:105:CYS:HB3	2.16	0.46
1:D:25:TYR:CD1	1:D:26:GLY:N	2.84	0.46
1:B:89:THR:HG23	1:B:92:GLU:OE1	2.16	0.45
1:B:19:ILE:HG12	1:B:23:SER:HB3	1.99	0.45
1:D:55:LEU:HD13	1:D:94:ARG:HB3	1.95	0.45
1:B:95:ILE:O	1:B:98:CYS:HB2	2.16	0.45
1:A:33:GLY:H	2:A:203:AJM:H161	1.81	0.45
1:B:90:SER:CA	1:B:93:ASN:HB2	2.40	0.45
1:D:125:LEU:HD23	1:D:127:LYS:CE	2.47	0.45
1:A:129:GLU:O	1:A:130:LEU:HG	2.17	0.45
1:A:130:LEU:HA	1:A:130:LEU:HD23	1.76	0.45
1:A:48:HIS:CE1	1:A:52:TYR:CZ	3.04	0.45
1:A:107:ARG:HB3	1:A:107:ARG:HH11	1.81	0.45
1:C:55:LEU:HD13	1:C:95:ILE:HD13	1.97	0.44
1:B:77:ARG:HA	1:B:81:ALA:O	2.17	0.44
1:D:22:TYR:O	1:D:29:CYS:HB3	2.17	0.44
1:A:61:CYS:O	1:A:68:PRO:HD2	2.18	0.44
1:B:30:GLY:O	2:B:202:AJM:H161	2.17	0.44
1:D:116:LYS:HE2	1:D:117:TYR:CZ	2.53	0.44
1:D:38:LYS:O	1:D:39:ASP:HB3	2.17	0.44
1:A:19:ILE:HG22	1:A:20:PRO:HD3	2.00	0.44
1:C:121:PRO:HG2	1:C:125:LEU:CD2	2.40	0.44
1:C:23:SER:CB	3:C:347:ANN:H6	2.44	0.44
1:C:129:GLU:HG2	1:C:130:LEU:H	1.81	0.44
1:C:89:THR:OG1	1:C:92:GLU:HG3	2.18	0.43
1:A:8:MET:SD	1:A:100:LYS:HA	2.58	0.43
1:D:30:GLY:HA2	3:D:350:ANN:C3	2.47	0.43
1:B:121:PRO:HG2	1:B:125:LEU:CD1	2.48	0.43
1:C:113:TYR:C	1:C:113:TYR:CD1	2.91	0.43
1:A:125:LEU:HD22	1:A:125:LEU:HA	1.66	0.43
1:A:19:ILE:HG22	1:A:20:PRO:N	2.35	0.42
1:A:67:ASN:O	1:A:70:SER:N	2.47	0.42
1:D:129:GLU:HA	1:D:129:GLU:OE2	2.19	0.42
1:D:91:CYS:HA	1:D:94:ARG:HD3	2.01	0.42
1:A:25:TYR:HE2	1:A:113:TYR:CE2	2.34	0.42
1:A:76:LYS:HE2	1:B:133:CYS:C	2.40	0.42
1:A:95:ILE:O	1:A:98:CYS:HB2	2.20	0.42
1:D:28:TYR:HB2	1:D:42:ASP:OD1	2.19	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:31:TRP:N	1:D:31:TRP:CD2	2.82	0.42
1:D:19:ILE:HD13	1:D:23:SER:HB3	1.99	0.42
1:D:10:LEU:HD13	1:D:17:LEU:HD23	2.02	0.41
1:A:114:SER:HG	1:A:117:TYR:HD2	1.68	0.41
1:A:117:TYR:HD1	1:A:120:TYR:CD1	2.38	0.41
1:D:30:GLY:HA2	3:D:350:ANN:C2	2.50	0.41
1:B:46:PHE:CE1	1:B:130:LEU:HD13	2.55	0.41
1:B:32:GLY:HA2	1:B:49:ASP:OD2	2.20	0.41
1:B:43:ARG:O	1:B:46:PHE:HB3	2.21	0.41
1:A:111:ASN:C	1:A:113:TYR:H	2.23	0.41
1:A:90:SER:OG	1:A:91:CYS:N	2.54	0.41
1:C:28:TYR:HB3	1:C:49:ASP:OD2	2.20	0.41
1:D:5:PHE:CD1	1:D:99:ASP:HB3	2.56	0.41
1:A:33:GLY:HA2	2:A:203:AJM:H122	2.01	0.41
1:C:55:LEU:HD12	1:C:95:ILE:HD13	2.01	0.41
1:A:9:ILE:H	1:A:9:ILE:CD1	2.30	0.41
1:C:82:ILE:O	1:C:100:LYS:HE3	2.21	0.41
1:D:35:GLY:HA2	1:D:127:LYS:O	2.21	0.41
1:B:127:LYS:N	1:B:127:LYS:HD2	2.36	0.41
1:D:77:ARG:HA	1:D:77:ARG:HD3	1.90	0.41
1:A:67:ASN:HA	1:A:68:PRO:HD2	1.76	0.40
1:D:12:GLU:HG3	1:D:82:ILE:HD11	2.03	0.40
1:B:130:LEU:CD2	1:B:131:LYS:H	2.34	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	119/121 (98%)	95 (80%)	14 (12%)	10 (8%)	1	10
1	B	119/121 (98%)	107 (90%)	10 (8%)	2 (2%)	14	59
1	C	119/121 (98%)	102 (86%)	14 (12%)	3 (2%)	9	47
1	D	119/121 (98%)	103 (87%)	11 (9%)	5 (4%)	4	31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	476/484 (98%)	407 (86%)	49 (10%)	20 (4%)	4	31

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	90	SER
1	A	121	PRO
1	A	124	PHE
1	A	125	LEU
1	B	127	LYS
1	D	78	VAL
1	D	79	ASN
1	D	125	LEU
1	A	78	VAL
1	A	86	LYS
1	A	122	ASP
1	B	89	THR
1	D	31	TRP
1	D	33	GLY
1	A	56	PRO
1	C	31	TRP
1	C	79	ASN
1	C	121	PRO
1	A	26	GLY

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	104/104 (100%)	85 (82%)	19 (18%)	2	11
1	B	104/104 (100%)	91 (88%)	13 (12%)	7	29
1	C	104/104 (100%)	91 (88%)	13 (12%)	7	29
1	D	104/104 (100%)	90 (86%)	14 (14%)	6	25
All	All	416/416 (100%)	357 (86%)	59 (14%)	5	22

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	9	ILE
1	A	34	LYS
1	A	38	LYS
1	A	43	ARG
1	A	56	PRO
1	A	59	ASP
1	A	61	CYS
1	A	68	PRO
1	A	70	SER
1	A	77	ARG
1	A	90	SER
1	A	112	THR
1	A	116	LYS
1	A	122	ASP
1	A	124	PHE
1	A	125	LEU
1	A	127	LYS
1	A	129	GLU
1	B	1	SER
1	B	3	LEU
1	B	20	PRO
1	B	34	LYS
1	B	69	LYS
1	B	70	SER
1	B	91	CYS
1	B	104	ILE
1	B	111	ASN
1	B	116	LYS
1	B	127	LYS
1	B	129	GLU
1	B	130	LEU
1	C	3	LEU
1	C	19	ILE
1	C	21	SER
1	C	23	SER
1	C	34	LYS
1	C	69	LYS
1	C	78	VAL
1	C	91	CYS
1	C	111	ASN
1	C	115	LYS

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Mol	Chain	Res	Type
1	C	122	ASP
1	C	127	LYS
1	C	130	LEU
1	D	2	LEU
1	D	20	PRO
1	D	34	LYS
1	D	59	ASP
1	D	70	SER
1	D	77	ARG
1	D	90	SER
1	D	94	ARG
1	D	115	LYS
1	D	116	LYS
1	D	122	ASP
1	D	124	PHE
1	D	125	LEU
1	D	130	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	B	111	ASN
1	C	67	ASN
1	C	108	GLN
1	C	111	ASN
1	D	93	ASN
1	D	108	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

8 ligands 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	AJM	A	203	-	27,27,27	2.19	6 (22%)	46,46,46	1.95	11 (23%)
3	ANN	A	349	-	11,11,11	2.56	4 (36%)	14,14,14	1.04	1 (7%)
2	AJM	B	202	-	27,27,27	2.02	7 (25%)	46,46,46	2.01	13 (28%)
3	ANN	B	348	-	11,11,11	2.37	3 (27%)	14,14,14	1.03	0
2	AJM	C	201	-	27,27,27	2.22	8 (29%)	46,46,46	1.73	9 (19%)
3	ANN	C	347	-	11,11,11	2.46	3 (27%)	14,14,14	1.02	0
2	AJM	D	204	-	27,27,27	2.36	8 (29%)	46,46,46	1.80	8 (17%)
3	ANN	D	350	-	11,11,11	2.59	4 (36%)	14,14,14	1.07	1 (7%)

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	AJM	A	203	-	-	0/0/67/67	0/0/6/6
3	ANN	A	349	-	-	0/6/6/6	0/1/1/1
2	AJM	B	202	-	-	0/0/67/67	0/0/6/6
3	ANN	B	348	-	-	0/6/6/6	0/1/1/1
2	AJM	C	201	-	-	0/0/67/67	0/0/6/6
3	ANN	C	347	-	-	0/6/6/6	0/1/1/1
2	AJM	D	204	-	-	0/0/67/67	0/0/6/6
3	ANN	D	350	-	-	0/6/6/6	0/1/1/1

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	350	ANN	C6-C5	6.94	1.53	1.38
3	C	347	ANN	C6-C5	6.78	1.53	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	349	ANN	C6-C5	6.58	1.52	1.38
3	B	348	ANN	C6-C5	6.57	1.52	1.38
2	D	204	AJM	C16-C17	6.12	1.58	1.51
2	A	203	AJM	C16-C17	5.52	1.57	1.51
2	D	204	AJM	C11-C6	5.30	1.46	1.39
2	C	201	AJM	C11-C6	5.09	1.46	1.39
2	A	203	AJM	C11-C6	4.78	1.45	1.39
2	C	201	AJM	C16-C17	4.73	1.56	1.51
2	B	202	AJM	C11-C6	4.53	1.45	1.39
2	D	204	AJM	C3-N2	4.35	1.55	1.49
2	B	202	AJM	C16-C17	4.09	1.56	1.51
2	D	204	AJM	C14-C3	3.82	1.59	1.54
2	A	203	AJM	C3-N2	3.77	1.54	1.49
2	A	203	AJM	C14-C3	3.68	1.59	1.54
2	B	202	AJM	C2-N2	3.55	1.54	1.49
2	C	201	AJM	C14-C3	3.52	1.58	1.54
2	C	201	AJM	C3-N2	3.42	1.54	1.49
2	B	202	AJM	C14-C3	3.30	1.58	1.54
2	C	201	AJM	C1-C2	-3.16	1.50	1.53
3	A	349	ANN	C4-C5	3.13	1.45	1.38
2	C	201	AJM	C4-C3	3.05	1.58	1.53
3	D	350	ANN	C4-C5	2.92	1.44	1.38
3	A	349	ANN	C7-C2	2.71	1.44	1.39
3	D	350	ANN	C3-C2	2.71	1.44	1.39
3	A	349	ANN	C3-C2	2.61	1.43	1.39
3	C	347	ANN	C3-C2	2.59	1.43	1.39
2	B	202	AJM	C17-N2	2.58	1.52	1.45
2	A	203	AJM	C2-N2	2.58	1.52	1.49
3	D	350	ANN	C7-C2	2.53	1.43	1.39
2	B	202	AJM	C13-C14	2.48	1.58	1.53
3	C	347	ANN	C4-C5	2.44	1.43	1.38
2	B	202	AJM	C3-N2	2.40	1.52	1.49
3	B	348	ANN	C4-C5	2.40	1.43	1.38
2	A	203	AJM	C5-C6	-2.37	1.47	1.51
3	B	348	ANN	C3-C2	2.31	1.43	1.39
2	C	201	AJM	C2-N2	2.22	1.52	1.49
2	C	201	AJM	C5-C15	2.18	1.55	1.53
2	D	204	AJM	C2-N2	2.13	1.52	1.49
2	D	204	AJM	C5-C15	2.09	1.55	1.53
2	D	204	AJM	C16-C13	2.03	1.58	1.53
2	D	204	AJM	C4-C3	2.01	1.56	1.53

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	202	AJM	C18-N1-C1	-8.08	107.01	116.24
2	A	203	AJM	C18-N1-C1	-7.52	107.64	116.24
2	D	204	AJM	C18-N1-C1	-6.80	108.47	116.24
2	C	201	AJM	C18-N1-C1	-6.17	109.19	116.24
2	A	203	AJM	C4-C3-C14	-5.17	101.61	105.18
2	D	204	AJM	C4-C3-C14	-5.12	101.65	105.18
2	C	201	AJM	C4-C3-C14	-4.01	102.42	105.18
2	B	202	AJM	C4-C3-C14	-3.74	102.60	105.18
2	B	202	AJM	C12-C2-C1	-3.58	101.53	110.27
2	B	202	AJM	C18-N1-C11	-3.21	109.07	120.84
2	A	203	AJM	C4-C3-N2	-3.12	106.40	109.53
2	D	204	AJM	C18-N1-C11	-3.04	109.69	120.84
2	A	203	AJM	C18-N1-C11	-2.96	110.00	120.84
2	D	204	AJM	C4-C3-N2	-2.91	106.61	109.53
2	D	204	AJM	C5-C6-C11	-2.90	105.44	110.26
2	A	203	AJM	C15-C14-C3	2.87	106.95	104.34
2	B	202	AJM	C1-C2-N2	2.86	117.93	110.72
2	C	201	AJM	C5-C1-N1	-2.85	99.57	102.11
2	C	201	AJM	C5-C6-C11	-2.84	105.55	110.26
2	C	201	AJM	C18-N1-C11	-2.81	110.54	120.84
2	C	201	AJM	C12-C2-C1	-2.80	103.44	110.27
2	A	203	AJM	C12-C2-C1	-2.73	103.60	110.27
2	B	202	AJM	C4-C3-N2	-2.58	106.94	109.53
2	A	203	AJM	C5-C6-C11	-2.57	105.99	110.26
2	A	203	AJM	C4-C5-C1	2.57	112.50	108.44
2	B	202	AJM	C12-C13-C14	2.53	114.00	110.64
2	B	202	AJM	C11-N1-C1	2.53	106.00	103.88
2	B	202	AJM	C15-C14-C3	2.51	106.63	104.34
2	D	204	AJM	C15-C14-C3	2.50	106.62	104.34
2	C	201	AJM	C15-C5-C1	2.39	113.49	110.42
2	B	202	AJM	C13-C12-C2	-2.38	105.75	108.50
2	B	202	AJM	C4-C5-C1	2.36	112.17	108.44
2	A	203	AJM	C1-C2-N2	2.35	116.64	110.72
2	B	202	AJM	C5-C6-C11	-2.33	106.39	110.26
2	D	204	AJM	C7-C6-C5	2.27	134.19	130.09
2	A	203	AJM	C16-C13-C14	-2.21	107.70	110.64
2	C	201	AJM	C7-C6-C5	2.18	134.02	130.09
3	D	350	ANN	C8-O3-C5	-2.11	112.40	117.54
2	B	202	AJM	C12-C13-C16	-2.11	104.00	109.17
3	A	349	ANN	C8-O3-C5	-2.10	112.43	117.54
2	D	204	AJM	C4-C5-C1	2.05	111.69	108.44
2	A	203	AJM	C11-N1-C1	2.03	105.58	103.88
2	C	201	AJM	C1-C2-N2	2.03	115.83	110.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

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 ⓘ

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	121/121 (100%)	-0.11	0 100 100	11, 38, 66, 75	0
1	B	121/121 (100%)	-0.17	0 100 100	6, 25, 52, 64	0
1	C	121/121 (100%)	-0.14	0 100 100	8, 26, 52, 61	0
1	D	121/121 (100%)	-0.10	0 100 100	7, 36, 66, 74	0
All	All	484/484 (100%)	-0.13	0 100 100	6, 32, 61, 75	0

There are no RSRZ outliers to report.

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

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ANN	B	348	11/11	0.42	8.88	39,40,42,42	11
3	ANN	C	347	11/11	0.44	8.82	36,39,40,41	11

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ANN	D	350	11/11	0.40	2.93	51,52,53,53	11
3	ANN	A	349	11/11	0.36	2.46	36,38,39,39	11
2	AJM	B	202	22/22	0.46	1.84	80,81,82,82	22
2	AJM	C	201	22/22	0.39	1.05	54,56,56,56	22
2	AJM	D	204	22/22	0.31	0.54	53,55,55,56	22
2	AJM	A	203	22/22	0.35	0.00	67,69,69,69	22

6.5 Other polymers ⓘ

There are no such residues in this entry.