



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:19 PM UTC

PDB ID : 3CBI / pdb_00003cbi
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 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

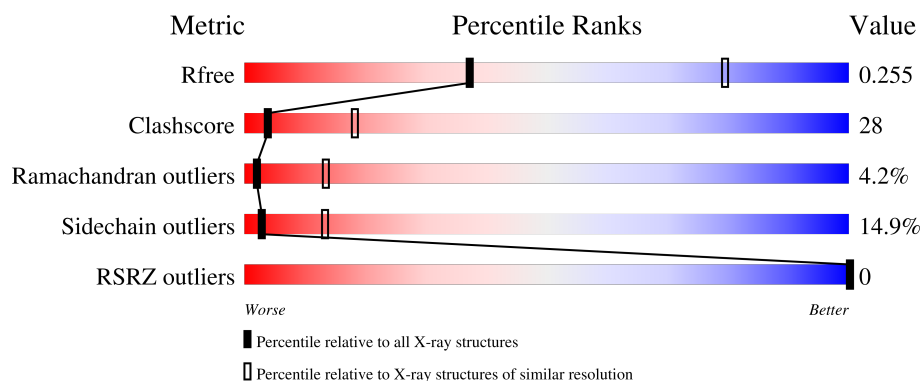
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	121	
1	B	121	
1	C	121	
1	D	121	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ANN	C	347	-	-	X	-

2 Entry composition [i](#)

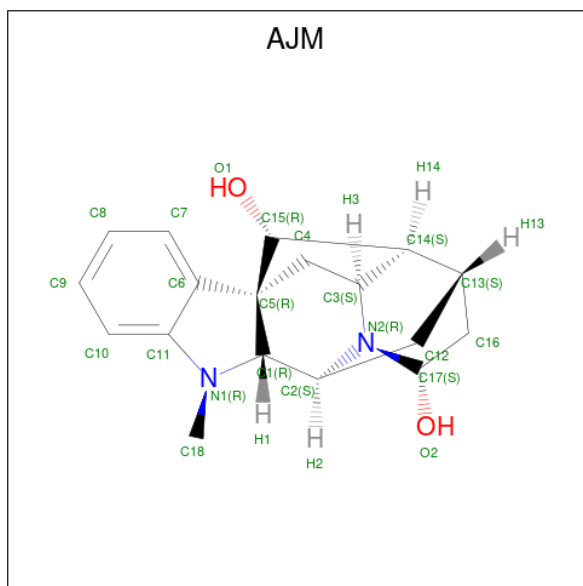
There are 3 unique types of molecules in this entry. The entry contains 3904 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 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 (CCD ID: AJM) (formula: $C_{18}H_{22}N_2O_2$).



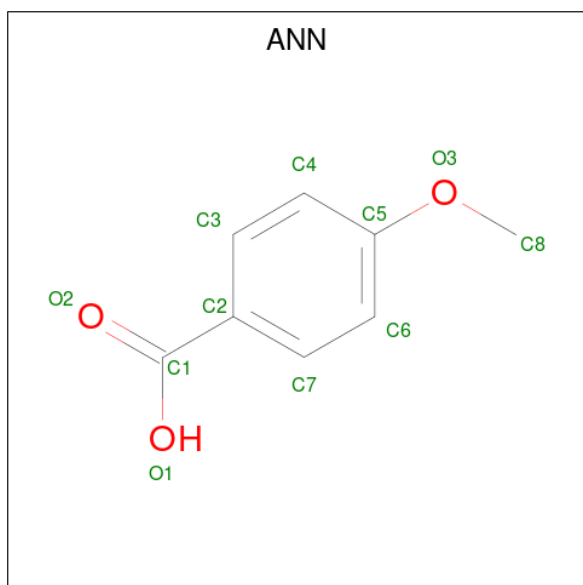
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			22	18	2	2		
2	B	1	Total	C	N	O	0	0
			22	18	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	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 (CCD ID: ANN) (formula: C₈H₈O₃).

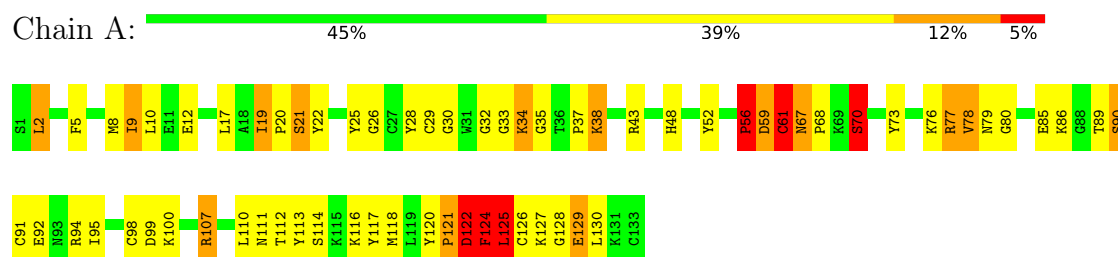


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	8	3		
3	B	1	Total	C	O	0	0
			11	8	3		
3	C	1	Total	C	O	0	0
			11	8	3		
3	D	1	Total	C	O	0	0
			11	8	3		

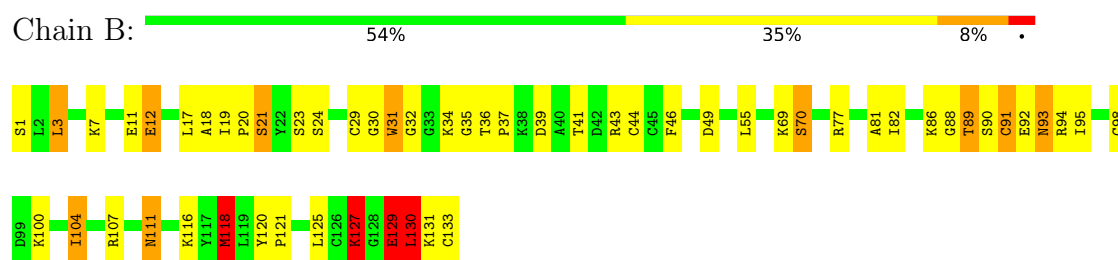
3 Residue-property plots

These plots are drawn for all protein, RNA, 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.

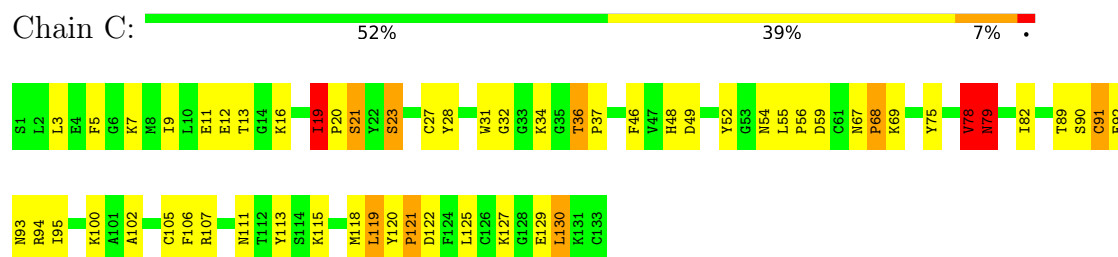
• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



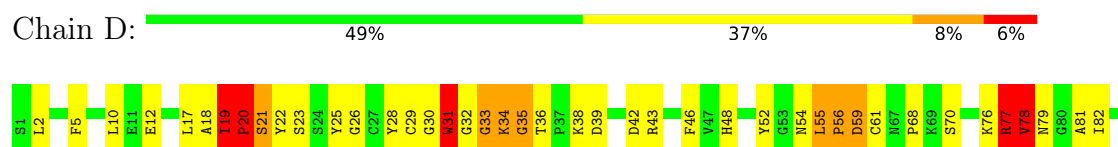
• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



• Molecule 1: Phospholipase A2 VRV-PL-VIIIa



T89	S90	C91	E92	N93	R94	I95	D99	R107	Q108	N109	L110	N111	K115	K116	Y117	M118	D122	F124	L125	G126	K127	G128	E129	L130	K131	C133
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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 50.00 – 3.15	Depositor EDS
% Data completeness (in resolution range)	92.2 (50.00-3.15) 97.3 (50.00-3.15)	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 , 0.255	Depositor DCC
R_{free} test set	513 reflections (5.30%)	wwPDB-VP
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 , 60.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.420 for -h,-k,l 0.024 for h,-h-k,-l 0.017 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.93	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.

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: 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	0.74	0/963	1.56	23/1292 (1.8%)
1	B	0.78	1/963 (0.1%)	1.36	16/1292 (1.2%)
1	C	0.73	1/963 (0.1%)	1.45	10/1292 (0.8%)
1	D	0.91	5/963 (0.5%)	1.59	21/1292 (1.6%)
All	All	0.79	7/3852 (0.2%)	1.49	70/5168 (1.4%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	19	ILE	N-CA	-10.06	1.37	1.46
1	C	122	ASP	CA-C	7.10	1.62	1.52
1	D	59	ASP	CA-C	-6.38	1.44	1.52
1	D	19	ILE	CA-C	-5.48	1.48	1.53
1	D	61	CYS	CA-C	5.28	1.58	1.52
1	D	20	PRO	CA-CB	-5.27	1.44	1.53
1	B	86	LYS	CA-C	5.12	1.59	1.52

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	128	GLY	N-CA-C	17.13	129.51	112.04
1	C	19	ILE	CA-C-N	-16.65	102.06	120.04
1	C	19	ILE	C-N-CA	-16.65	102.06	120.04
1	D	35	GLY	N-CA-C	16.48	128.72	112.08
1	D	19	ILE	N-CA-C	-11.90	94.45	107.77
1	D	34	LYS	N-CA-C	10.59	124.85	110.55
1	B	130	LEU	N-CA-C	-10.35	92.21	109.80
1	C	78	VAL	N-CA-C	-9.91	101.37	111.88
1	A	19	ILE	CA-C-N	-9.28	109.35	119.19
1	A	19	ILE	C-N-CA	-9.28	109.35	119.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	77	ARG	N-CA-C	9.04	121.21	111.36
1	D	21	SER	N-CA-C	8.92	122.12	111.33
1	A	34	LYS	N-CA-C	8.81	122.28	109.14
1	B	129	GLU	N-CA-C	8.81	123.42	109.50
1	D	61	CYS	CA-C-N	-8.64	112.19	123.34
1	D	61	CYS	C-N-CA	-8.64	112.19	123.34
1	A	35	GLY	N-CA-C	8.37	120.58	112.04
1	D	18	ALA	N-CA-C	8.29	121.13	111.02
1	C	118	MET	N-CA-C	-8.15	99.76	110.43
1	A	61	CYS	N-CA-C	8.13	120.72	109.18
1	D	19	ILE	CA-C-O	7.85	124.93	119.20
1	B	94	ARG	N-CA-C	7.76	119.82	111.36
1	A	56	PRO	CA-C-N	-7.72	110.53	122.49
1	A	56	PRO	C-N-CA	-7.72	110.53	122.49
1	B	21	SER	N-CA-C	7.71	119.69	111.28
1	C	36	THR	N-CA-C	-7.48	100.66	109.93
1	B	118	MET	N-CA-C	-7.35	99.40	110.28
1	B	70	SER	N-CA-C	7.34	121.95	112.41
1	D	59	ASP	O-C-N	7.16	132.11	122.59
1	B	93	ASN	N-CA-C	-6.99	102.49	111.02
1	A	61	CYS	CA-C-N	-6.86	114.19	123.46
1	A	61	CYS	C-N-CA	-6.86	114.19	123.46
1	D	116	LYS	N-CA-C	-6.68	104.61	112.89
1	A	19	ILE	N-CA-C	6.63	123.20	108.88
1	A	85	GLU	N-CA-C	6.62	119.94	109.81
1	D	109	ASN	N-CA-C	6.49	120.83	112.26
1	C	19	ILE	N-CA-C	6.47	122.86	108.88
1	A	56	PRO	N-CA-C	-6.47	99.15	112.47
1	A	70	SER	N-CA-C	6.43	121.65	113.55
1	B	29	CYS	N-CA-C	6.37	118.32	109.15
1	B	36	THR	N-CA-C	-6.27	102.17	110.07
1	B	31	TRP	N-CA-C	6.25	117.89	111.14
1	D	61	CYS	O-C-N	-6.12	117.40	123.46
1	A	38	LYS	N-CA-C	5.91	117.72	111.28
1	A	122	ASP	N-CA-C	-5.90	98.23	110.80
1	D	56	PRO	CB-CA-C	-5.81	103.61	111.85
1	A	21	SER	N-CA-C	5.70	118.26	111.71
1	A	67	ASN	CA-C-N	5.66	126.91	119.84
1	A	67	ASN	C-N-CA	5.66	126.91	119.84
1	C	68	PRO	N-CA-C	-5.65	106.79	113.86
1	D	36	THR	N-CA-C	-5.62	98.37	109.10
1	C	21	SER	N-CA-C	5.51	119.97	112.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	12	GLU	N-CA-C	5.49	120.85	113.72
1	B	88	GLY	N-CA-C	-5.44	100.29	113.18
1	D	55	LEU	CA-C-N	5.42	125.09	119.56
1	D	55	LEU	C-N-CA	5.42	125.09	119.56
1	D	111	ASN	N-CA-C	5.39	119.11	112.54
1	D	19	ILE	O-C-N	-5.30	117.32	121.46
1	B	35	GLY	N-CA-C	5.21	121.68	112.83
1	A	28	TYR	N-CA-C	5.20	120.62	113.97
1	A	12	GLU	N-CA-C	5.16	116.59	111.07
1	C	79	ASN	N-CA-C	-5.15	99.84	110.80
1	B	39	ASP	N-CA-C	5.12	114.60	107.73
1	D	31	TRP	CA-CB-CG	-5.12	103.87	113.60
1	B	120	TYR	CA-C-N	5.12	125.92	119.98
1	B	120	TYR	C-N-CA	5.12	125.92	119.98
1	D	78	VAL	N-CA-C	5.06	119.87	109.34
1	A	73	TYR	N-CA-C	5.05	117.35	109.52
1	C	27	CYS	N-CA-C	5.03	119.56	113.38
1	A	118	MET	N-CA-C	-5.01	102.64	110.10

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	943	0	907	69	0
1	B	943	0	907	47	0
1	C	943	0	907	62	0
1	D	943	0	907	44	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	11	0	7	2	0
All	All	3904	0	3744	217	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 28.

All (217) 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: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
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:HG13	1:C:78:VAL:O	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:9:ILE:HD12	1:C:106:PHE:HD2	1.58	0.68
1:C:54:ASN:O	1:C:56:PRO:HD3	1.93	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
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:C:32:GLY:C	2:C:201:AJM:H122	2.24	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:SER:O	1:B:91:CYS:C	2.46	0.58
1:A:90:SER:C	1:A:94:ARG:HG3	2.29	0.58
1:C:68:PRO:HB3	1:C:95:ILE:CD1	2.34	0.58
1:A:25:TYR:CE2	1:A:113:TYR:CE2	2.91	0.57
1:A:121:PRO:HB3	1:A:124:PHE:CZ	2.38	0.57
1:C:19:ILE:O	1:C:23:SER:HB3	2.05	0.57
1:C:49:ASP:HB3	2:C:201:AJM:H161	1.85	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:A:9:ILE:HD13	1:A:9:ILE:H	1.69	0.57
1:C:78:VAL:O	1:C:79:ASN:HB2	2.05	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:A:122:ASP:O	1:A:125:LEU:N	2.41	0.54
1:B:1:SER:OG	1:B:3:LEU:HB2	2.07	0.54
1:C:55:LEU:HD21	1:C:94:ARG:CB	2.38	0.54
1:D:124:PHE:O	1:D:125:LEU:HB2	2.08	0.54
1:C:32:GLY:HA3	2:C:201:AJM:C12	2.34	0.53
1:B:127:LYS:H	1:B:127:LYS:HD2	1.72	0.53
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:A:56:PRO:CD	1:A:59:ASP:N	2.69	0.52
1:C:130:LEU:CD2	1:D:78:VAL:HG13	2.39	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:B:23:SER:CB	3:B:348:ANN:H6	2.41	0.51
1:C:5:PHE:O	1:C:9:ILE:HG12	2.11	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:CD1	1:A:9:ILE:N	2.75	0.49
1:D:54:ASN:O	1:D:56:PRO:HD3	2.13	0.48
1:A:122:ASP:O	1:A:126:CYS:N	2.40	0.48
1:B:90:SER:O	1:B:93:ASN:HB2	2.13	0.48
1:A:32:GLY:CA	2:A:203:AJM:H161	2.40	0.48
1:C:23:SER:CA	3:C:347:ANN:H6	2.42	0.48
1:C:68:PRO:CB	1:C:95:ILE:HD12	2.38	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:C:113:TYR:CD1	1:C:113:TYR:C	2.91	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:C:119:LEU:N	1:C:119:LEU:HD23	2.29	0.47
1:D:77:ARG:O	1:D:81:ALA:O	2.32	0.47
1:B:11:GLU:CB	1:B:77:ARG:NH2	2.76	0.47
1:C:56:PRO:HD2	1:C:59:ASP:H	1.79	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:2:LEU:HD12	1:A:2:LEU:HA	1.75	0.46
1:A:89:THR:OG1	1:A:92:GLU:HG3	2.15	0.46
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:A:111:ASN:C	1:A:113:TYR:H	2.23	0.46
1:A:76:LYS:HE2	1:B:133:CYS:C	2.40	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:19:ILE:HG12	1:B:23:SER:HB3	1.99	0.45
1:B:89:THR:HG23	1:B:92:GLU:OE1	2.16	0.45
1:A:122:ASP:CG	1:A:124:PHE:N	2.73	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:SER:CA	1:B:93:ASN:HB2	2.40	0.45
1:A:48:HIS:CE1	1:A:52:TYR:CZ	3.04	0.45
1:A:107:ARG:HH11	1:A:107:ARG:HB3	1.81	0.45
1:A:129:GLU:O	1:A:130:LEU:HG	2.17	0.45
1:A:130:LEU:HD23	1:A:130:LEU:HA	1.76	0.45
1:D:125:LEU:HD23	1:D:127:LYS:CE	2.47	0.45
1:A:120:TYR:HH	1:A:125:LEU:HD12	1.78	0.44
1:B:77:ARG:HA	1:B:81:ALA:O	2.17	0.44
1:C:55:LEU:HD13	1:C:95:ILE:HD13	1.97	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:D:90:SER:C	1:D:94:ARG:HD2	2.43	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:A:19:ILE:HG22	1:A:20:PRO:HD3	2.00	0.44
1:C:23:SER:CB	3:C:347:ANN:H6	2.44	0.44
1:C:121:PRO:HG2	1:C:125:LEU:CD2	2.40	0.44
1:C:129:GLU:HG2	1:C:130:LEU:H	1.81	0.44
1:D:38:LYS:O	1:D:39:ASP:HB3	2.17	0.44
1:A:8:MET:SD	1:A:100:LYS:HA	2.58	0.43
1:C:89:THR:OG1	1:C:92:GLU:HG3	2.18	0.43
1:D:30:GLY:HA2	3:D:350:ANN:C3	2.47	0.43
1:D:32:GLY:O	1:D:33:GLY:C	2.60	0.43
1:A:78:VAL:C	1:A:80:GLY:H	2.26	0.43
1:B:121:PRO:HG2	1:B:125:LEU:CD1	2.48	0.43
1:A:125:LEU:HA	1:A:125:LEU:HD22	1.66	0.43
1:B:92:GLU:O	1:B:93:ASN:C	2.63	0.42
1:A:67:ASN:O	1:A:70:SER:N	2.47	0.42
1:A:99:ASP:O	1:A:100:LYS:C	2.61	0.42
1:D:129:GLU:HA	1:D:129:GLU:OE2	2.19	0.42
1:A:8:MET:O	1:A:9:ILE:C	2.63	0.42
1:A:19:ILE:HG22	1:A:20:PRO:N	2.35	0.42
1:A:25:TYR:HE2	1:A:113:TYR:CE2	2.34	0.42
1:D:91:CYS:HA	1:D:94:ARG:HD3	2.01	0.42
1:A:95:ILE:O	1:A:98:CYS:HB2	2.20	0.42
1:A:90:SER:O	1:A:91:CYS:C	2.63	0.42
1:D:28:TYR:HB2	1:D:42:ASP:OD1	2.19	0.42
1:D:31:TRP:CD2	1:D:31:TRP:N	2.82	0.42
1:D:19:ILE:HD13	1:D:23:SER:HB3	1.99	0.42
1:A:114:SER:HG	1:A:117:TYR:HD2	1.68	0.41
1:D:10:LEU:HD13	1:D:17:LEU:HD23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:32:GLY:HA2	1:B:49:ASP:OD2	2.20	0.41
1:B:46:PHE:CE1	1:B:130:LEU:HD13	2.55	0.41
1:C:13:THR:C	1:C:16:LYS:N	2.79	0.41
1:B:43:ARG:O	1:B:46:PHE:HB3	2.21	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:9:ILE:H	1:A:9:ILE:CD1	2.30	0.41
1:A:33:GLY:HA2	2:A:203:AJM:H122	2.01	0.41
1:B:127:LYS:HD2	1:B:127:LYS:N	2.36	0.41
1:C:55:LEU:HD12	1:C:95:ILE:HD13	2.01	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:17:LEU:O	1:B:18:ALA:C	2.64	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:118:MET:HB3	1:B:118:MET:HE3	1.83	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 [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	119/121 (98%)	95 (80%)	14 (12%)	10 (8%)	0	3
1	B	119/121 (98%)	107 (90%)	10 (8%)	2 (2%)	7	30
1	C	119/121 (98%)	102 (86%)	14 (12%)	3 (2%)	4	22
1	D	119/121 (98%)	103 (87%)	11 (9%)	5 (4%)	2	13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	476/484 (98%)	407 (86%)	49 (10%)	20 (4%)	2	13

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	8
1	B	104/104 (100%)	91 (88%)	13 (12%)	4	19
1	C	104/104 (100%)	90 (86%)	14 (14%)	4	16
1	D	104/104 (100%)	88 (85%)	16 (15%)	2	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	416/416 (100%)	354 (85%)	62 (15%)	3 13

All (62) 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	70	SER
1	A	77	ARG
1	A	107	ARG
1	A	110	LEU
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	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	118	MET
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

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Mol	Chain	Res	Type
1	C	67	ASN
1	C	69	LYS
1	C	78	VAL
1	C	91	CYS
1	C	111	ASN
1	C	115	LYS
1	C	119	LEU
1	C	127	LYS
1	C	130	LEU
1	D	2	LEU
1	D	19	ILE
1	D	20	PRO
1	D	34	LYS
1	D	59	ASP
1	D	70	SER
1	D	76	LYS
1	D	77	ARG
1	D	90	SER
1	D	107	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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

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

5.3.3 RNA ⓘ

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](#)

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	D	204	-	26,27,27	2.46	10 (38%)	23,46,46	1.67	6 (26%)
2	AJM	A	203	-	26,27,27	2.31	9 (34%)	23,46,46	1.74	6 (26%)
3	ANN	C	347	-	11,11,11	2.76	3 (27%)	14,14,14	1.12	0
3	ANN	B	348	-	11,11,11	2.65	3 (27%)	14,14,14	1.12	0
2	AJM	C	201	-	26,27,27	2.34	9 (34%)	23,46,46	1.63	5 (21%)
2	AJM	B	202	-	26,27,27	2.15	10 (38%)	23,46,46	1.89	7 (30%)
3	ANN	D	350	-	11,11,11	2.89	4 (36%)	14,14,14	1.16	1 (7%)
3	ANN	A	349	-	11,11,11	2.85	4 (36%)	14,14,14	1.14	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
3	ANN	C	347	-	-	0/6/6/6	0/1/1/1
3	ANN	B	348	-	-	0/6/6/6	0/1/1/1
3	ANN	A	349	-	-	0/6/6/6	0/1/1/1
3	ANN	D	350	-	-	0/6/6/6	0/1/1/1

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	350	ANN	C6-C5	7.81	1.53	1.38
3	C	347	ANN	C6-C5	7.63	1.53	1.38
3	A	349	ANN	C6-C5	7.41	1.52	1.38
3	B	348	ANN	C6-C5	7.39	1.52	1.38
2	D	204	AJM	C11-C6	6.15	1.46	1.39
2	C	201	AJM	C11-C6	5.90	1.46	1.39
2	A	203	AJM	C11-C6	5.54	1.45	1.39
2	D	204	AJM	C16-C17	5.37	1.58	1.51
2	B	202	AJM	C11-C6	5.24	1.45	1.39
2	A	203	AJM	C16-C17	4.83	1.57	1.51
2	D	204	AJM	C3-N2	4.55	1.55	1.48
2	C	201	AJM	C16-C17	4.14	1.56	1.51
2	A	203	AJM	C3-N2	3.96	1.54	1.48
2	D	204	AJM	C14-C3	3.83	1.59	1.54
2	B	202	AJM	C2-N2	3.74	1.54	1.48
2	A	203	AJM	C14-C3	3.69	1.59	1.54
2	C	201	AJM	C3-N2	3.60	1.54	1.48
2	B	202	AJM	C16-C17	3.58	1.56	1.51
2	C	201	AJM	C14-C3	3.52	1.58	1.54
3	A	349	ANN	C4-C5	3.46	1.45	1.38
2	C	201	AJM	C4-C3	3.43	1.58	1.53
2	B	202	AJM	C14-C3	3.29	1.58	1.54
3	D	350	ANN	C4-C5	3.23	1.44	1.38
2	C	201	AJM	C1-C2	-3.14	1.50	1.53
2	B	202	AJM	C13-C14	3.04	1.58	1.54
3	A	349	ANN	C7-C2	3.01	1.44	1.39
3	D	350	ANN	C3-C2	3.01	1.44	1.39
3	A	349	ANN	C3-C2	2.89	1.43	1.39
3	C	347	ANN	C3-C2	2.87	1.43	1.39
3	D	350	ANN	C7-C2	2.81	1.43	1.39
2	A	203	AJM	C2-N2	2.74	1.52	1.48
3	C	347	ANN	C4-C5	2.68	1.43	1.38
3	B	348	ANN	C4-C5	2.64	1.43	1.38
2	A	203	AJM	C5-C6	-2.60	1.47	1.51
2	B	202	AJM	C3-N2	2.56	1.52	1.48
3	B	348	ANN	C3-C2	2.55	1.43	1.39
2	C	201	AJM	C2-N2	2.37	1.52	1.48
2	B	202	AJM	C10-C11	2.35	1.43	1.39
2	D	204	AJM	C2-N2	2.28	1.52	1.48
2	A	203	AJM	C10-C11	2.25	1.43	1.39
2	C	201	AJM	C10-C11	2.23	1.43	1.39
2	D	204	AJM	C4-C3	2.19	1.56	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	204	AJM	C10-C11	2.18	1.43	1.39
2	A	203	AJM	C13-C14	2.17	1.57	1.54
2	C	201	AJM	C11-N1	2.15	1.43	1.39
2	D	204	AJM	C12-C2	2.15	1.58	1.53
2	D	204	AJM	C16-C13	2.14	1.58	1.53
2	B	202	AJM	C5-C6	-2.13	1.48	1.51
2	B	202	AJM	C4-C3	2.13	1.56	1.53
2	A	203	AJM	C11-N1	2.11	1.43	1.39
2	D	204	AJM	C13-C14	2.09	1.57	1.54
2	B	202	AJM	C7-C6	2.02	1.41	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	202	AJM	C18-N1-C1	-4.56	107.01	117.38
2	A	203	AJM	C18-N1-C1	-4.28	107.64	117.38
2	D	204	AJM	C18-N1-C1	-3.91	108.47	117.38
2	B	202	AJM	C18-N1-C11	-3.83	109.07	120.94
2	D	204	AJM	C18-N1-C11	-3.63	109.69	120.94
2	C	201	AJM	C18-N1-C1	-3.60	109.19	117.38
2	A	203	AJM	C18-N1-C11	-3.53	110.00	120.94
2	C	201	AJM	C18-N1-C11	-3.35	110.54	120.94
2	B	202	AJM	C12-C13-C14	3.19	114.00	110.40
2	B	202	AJM	C12-C13-C16	-3.02	104.00	109.06
2	D	204	AJM	C5-C6-C11	-2.78	105.44	110.20
2	C	201	AJM	C5-C6-C11	-2.72	105.55	110.20
2	A	203	AJM	C12-C13-C16	-2.69	104.56	109.06
2	C	201	AJM	C12-C13-C16	-2.62	104.66	109.06
2	A	203	AJM	C5-C6-C11	-2.46	105.99	110.20
2	D	204	AJM	C12-C13-C16	-2.45	104.95	109.06
2	A	203	AJM	C16-C13-C14	-2.38	107.70	110.40
3	D	350	ANN	C8-O3-C5	-2.38	112.40	117.50
3	A	349	ANN	C8-O3-C5	-2.36	112.43	117.50
2	B	202	AJM	C13-C12-C2	-2.31	105.75	108.50
2	D	204	AJM	C7-C6-C5	2.24	134.19	130.21
2	A	203	AJM	C10-C11-C6	-2.23	119.60	121.76
2	B	202	AJM	C5-C6-C11	-2.22	106.39	110.20
2	C	201	AJM	C7-C6-C5	2.14	134.02	130.21
2	D	204	AJM	C16-C13-C14	-2.08	108.05	110.40
2	B	202	AJM	C10-C11-C6	-2.05	119.78	121.76

There are no chirality outliers.

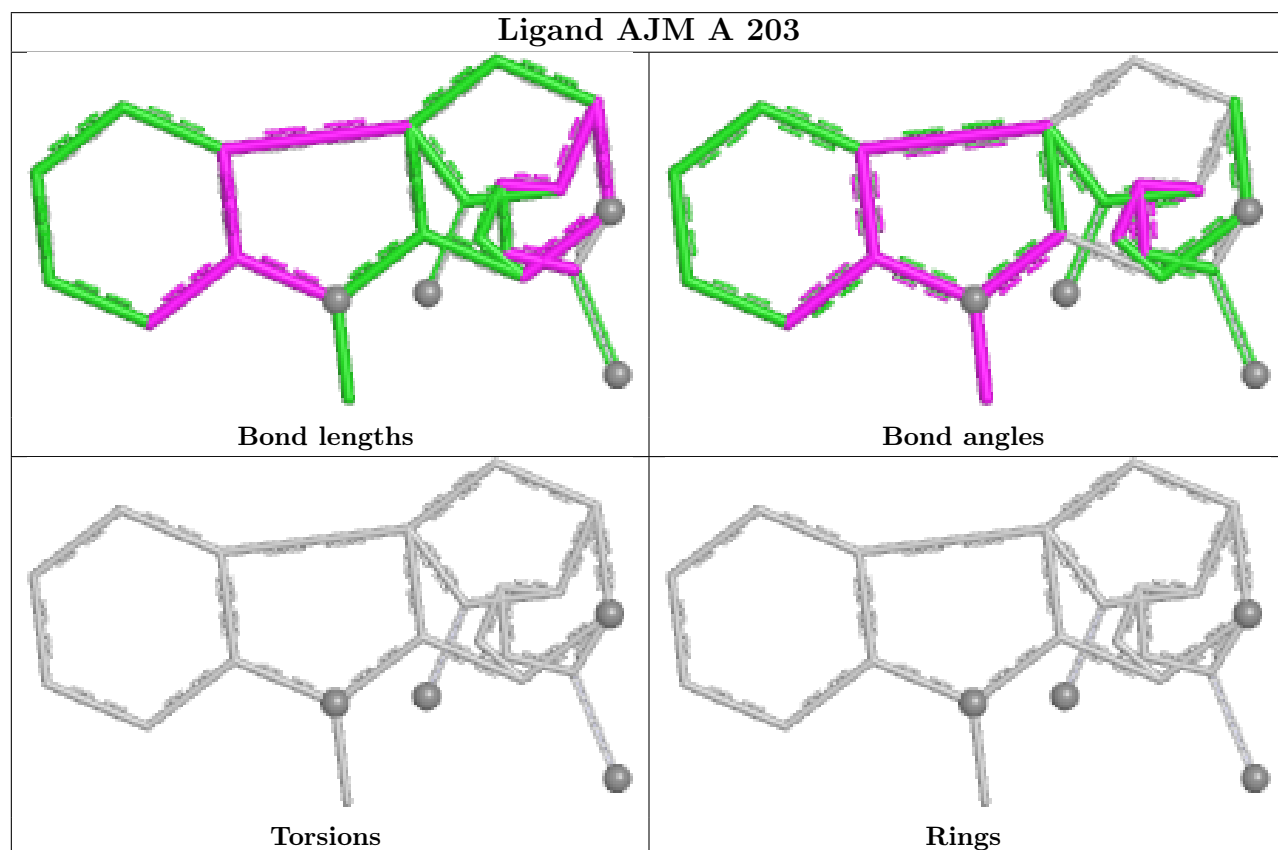
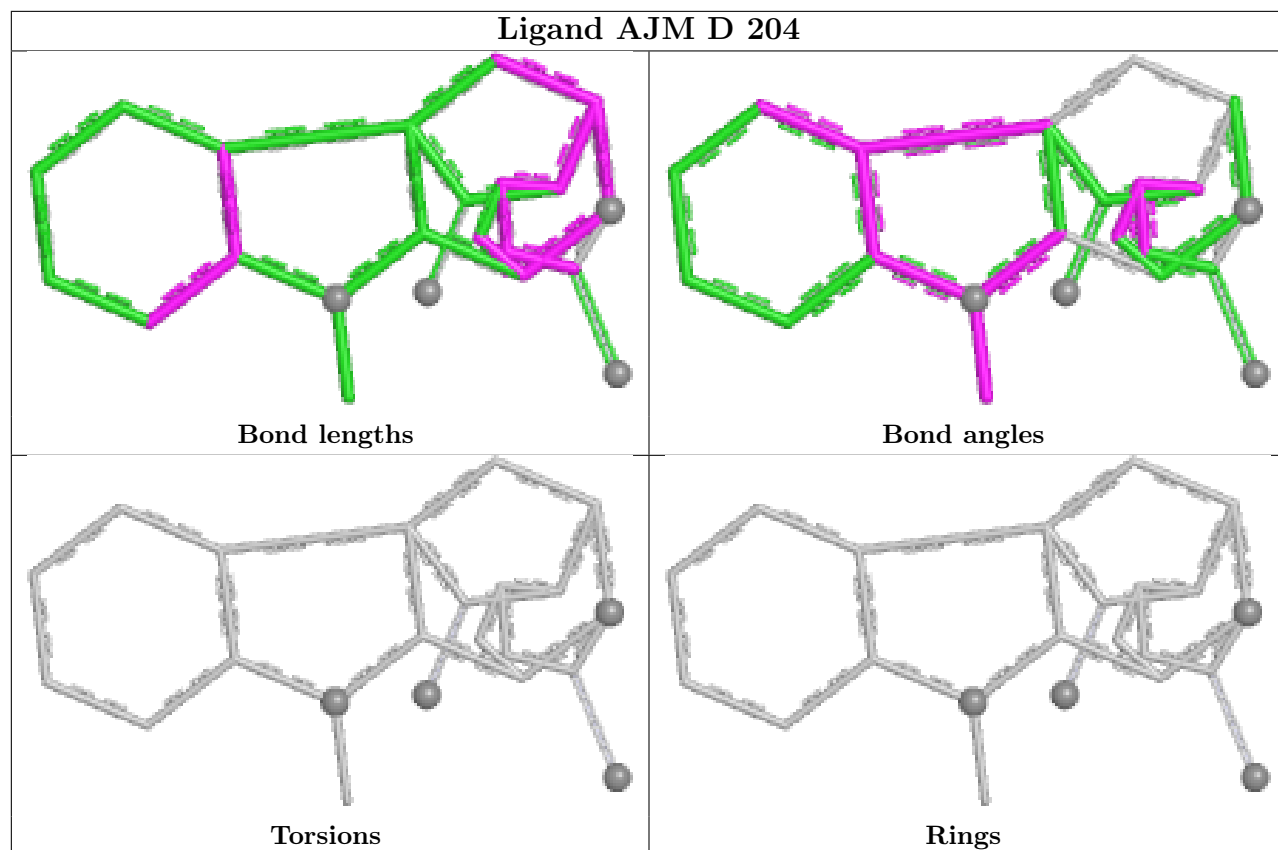
There are no torsion outliers.

There are no ring outliers.

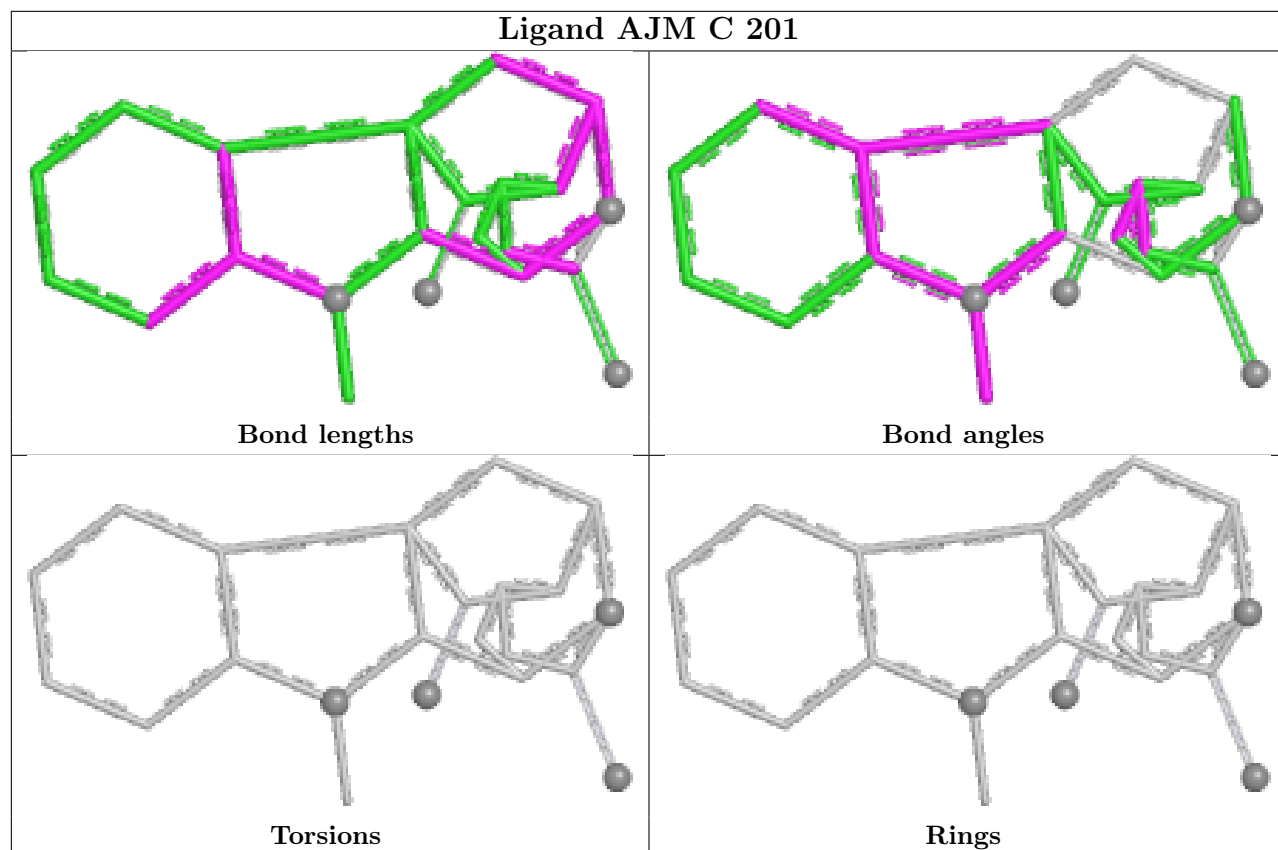
7 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	203	AJM	6	0
3	C	347	ANN	4	0
3	B	348	ANN	2	0
2	C	201	AJM	7	0
2	B	202	AJM	3	0
3	D	350	ANN	2	0
3	A	349	ANN	2	0

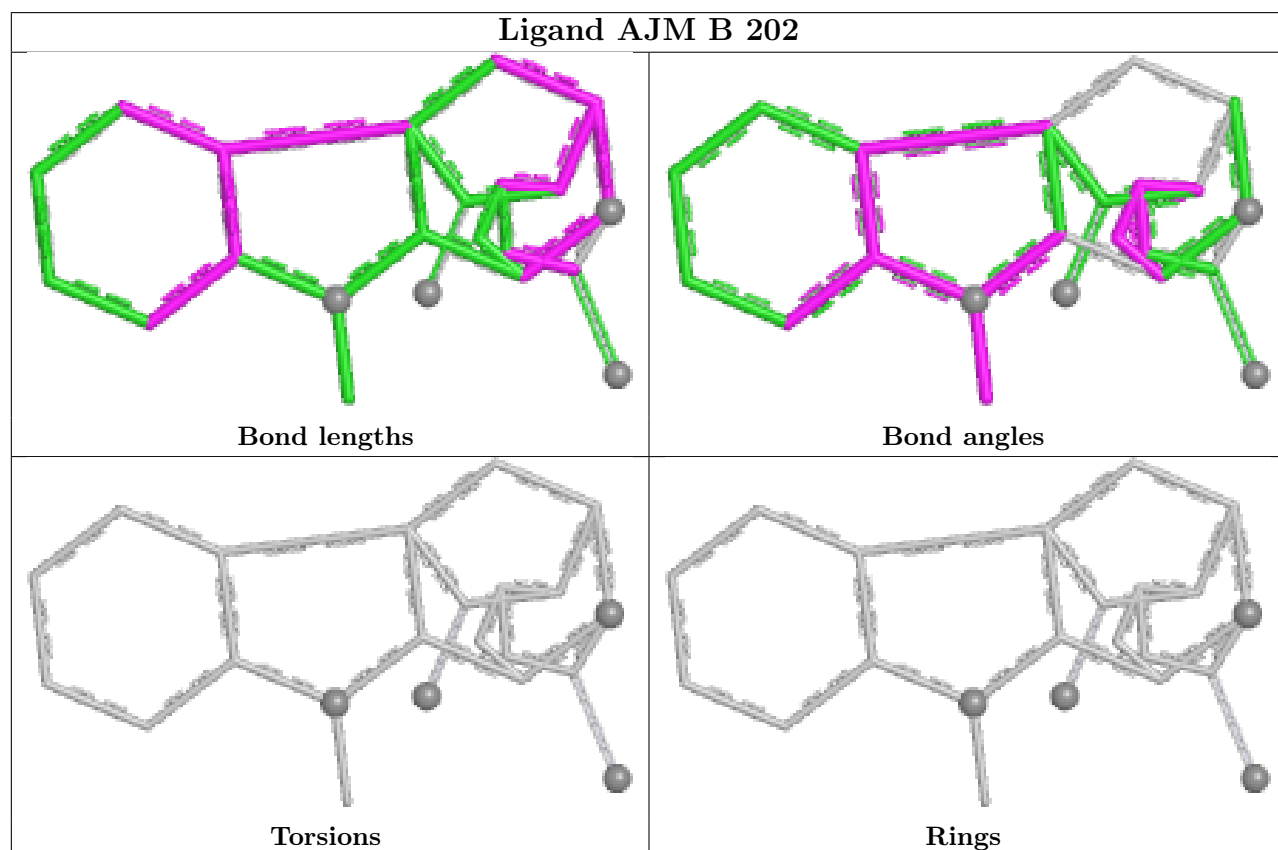
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand AJM C 201



Ligand AJM B 202



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	121/121 (100%)	-1.29	0 100 100	11, 38, 66, 75	0
1	B	121/121 (100%)	-1.42	0 100 100	6, 25, 52, 64	0
1	C	121/121 (100%)	-1.37	0 100 100	8, 26, 52, 61	0
1	D	121/121 (100%)	-1.35	0 100 100	7, 36, 66, 74	0
All	All	484/484 (100%)	-1.36	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 [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(Å ²)	Q<0.9
3	ANN	D	350	11/11	0.94	0.19	51,52,53,53	11
3	ANN	C	347	11/11	0.95	0.18	36,39,40,41	11
3	ANN	A	349	11/11	0.95	0.15	36,38,39,39	11
2	AJM	B	202	22/22	0.96	0.10	80,81,82,82	22

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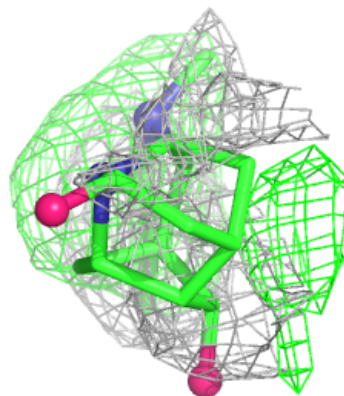
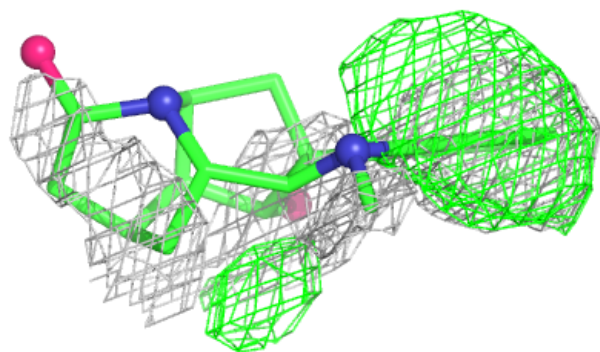
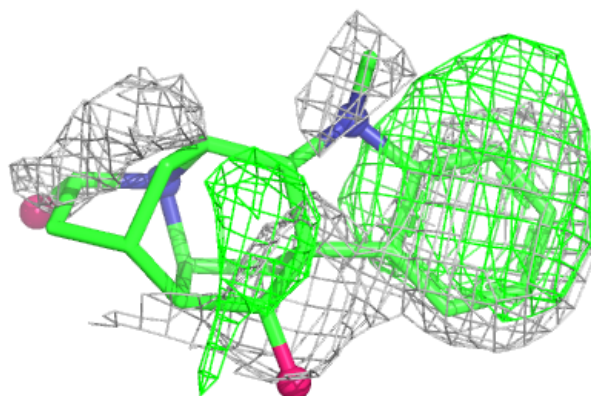
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ANN	B	348	11/11	0.96	0.17	39,40,42,42	11
2	AJM	A	203	22/22	0.97	0.08	67,69,69,69	22
2	AJM	D	204	22/22	0.97	0.07	53,55,55,56	22
2	AJM	C	201	22/22	0.98	0.07	54,56,56,56	22

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

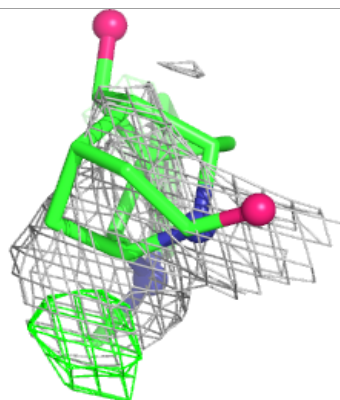
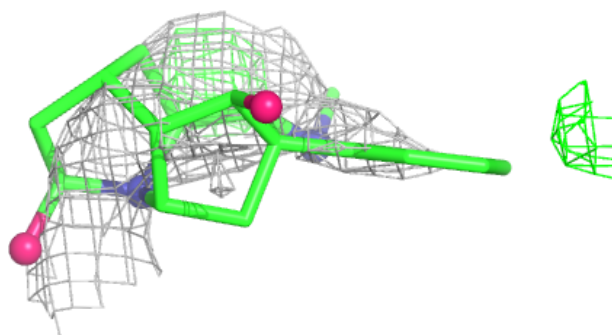
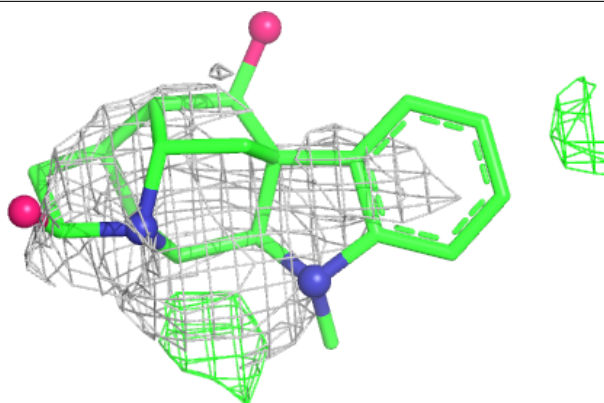
Electron density around AJM B 202:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

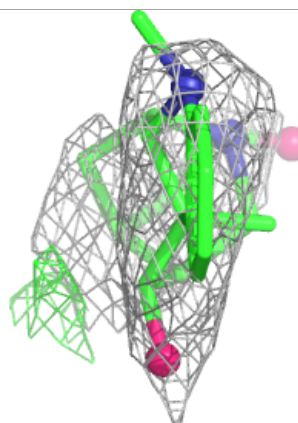
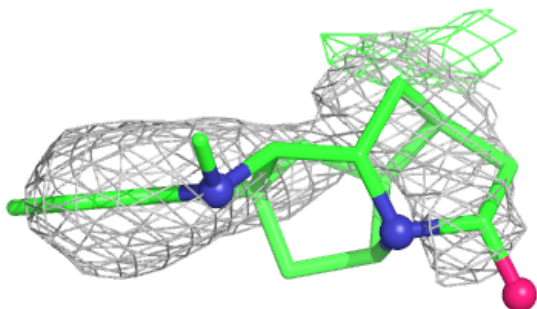
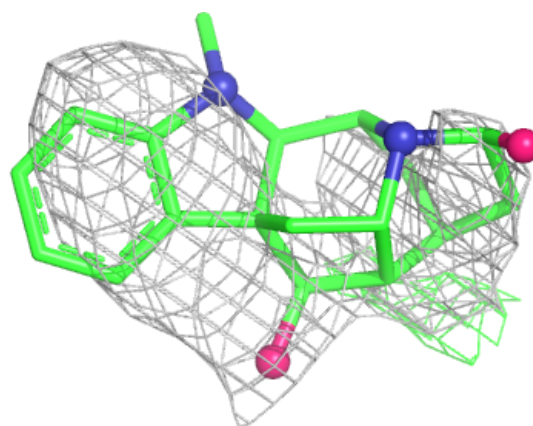


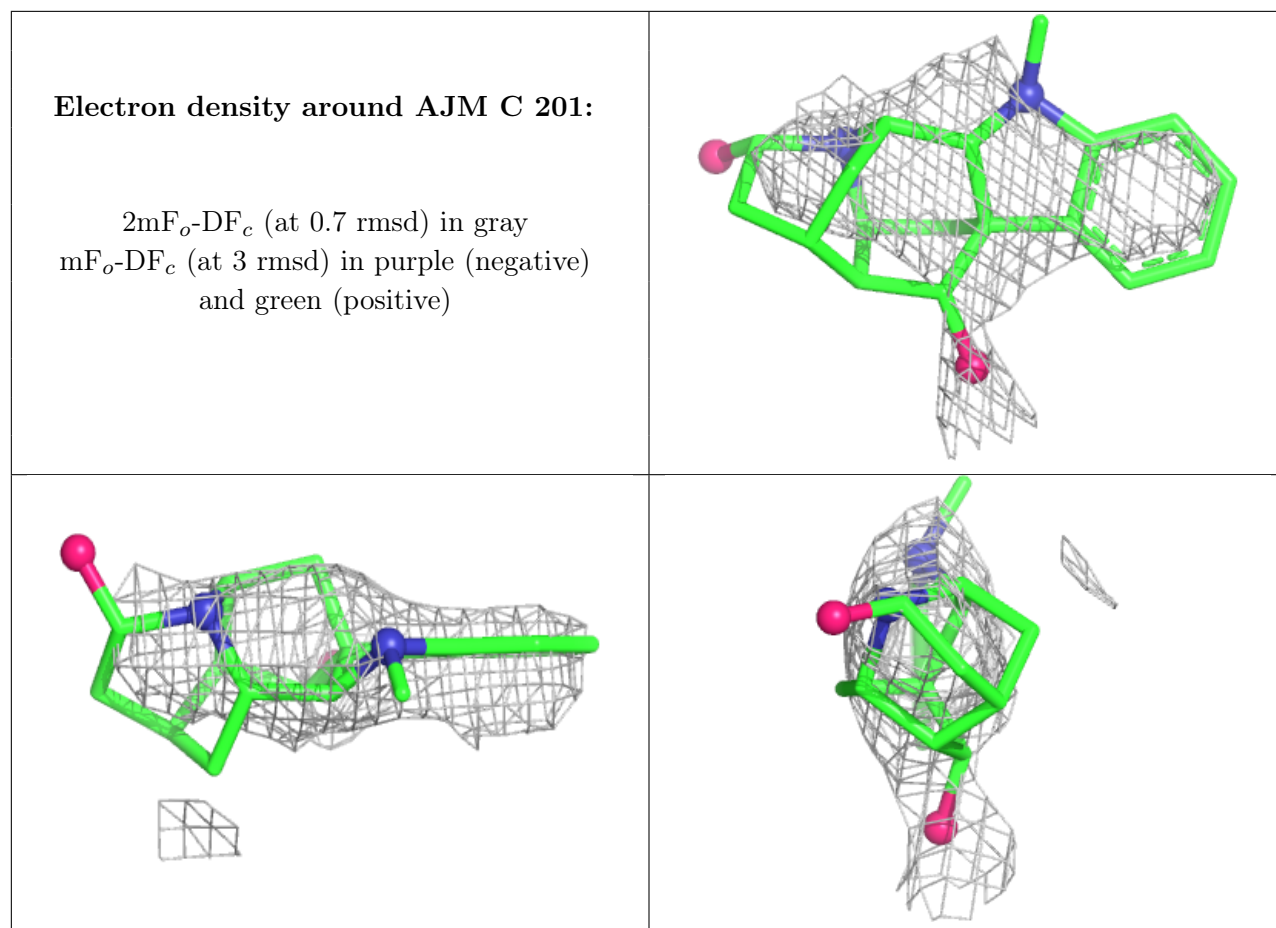
Electron density around AJM A 203:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around AJM D 204:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers ⓘ

There are no such residues in this entry.