



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2024 – 09:45 pm BST

PDB ID : 4ZGV
Title : The Crystal Structure of the Ferredoxin Receptor FusA from *Pectobacterium atrosepticum* SCRI1043
Authors : Grinter, R.; Josts, I.; Roszak, A.W.; Cogdell, R.J.; Walker, D.
Deposited on : 2015-04-24
Resolution : 3.20 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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

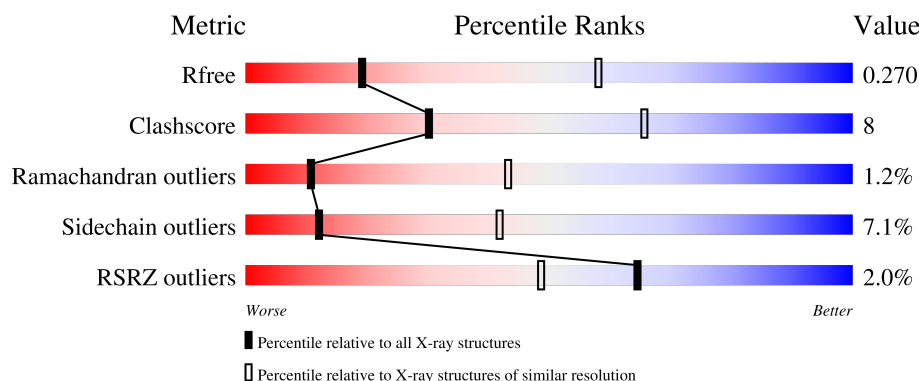
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	164625	1370 (3.20-3.20)
Clashscore	180529	1497 (3.20-3.20)
Ramachandran outliers	177936	1479 (3.20-3.20)
Sidechain outliers	177891	1478 (3.20-3.20)
RSRZ outliers	164620	1371 (3.20-3.20)

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	868	
1	B	868	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13030 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 Ferredoxin receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	809	Total	C	N	O	S	0	0	0
			6450	4062	1090	1279	19			
1	B	809	Total	C	N	O	S	0	0	0
			6450	4062	1090	1279	19			

There are 42 discrepancies between the modelled and reference sequences:

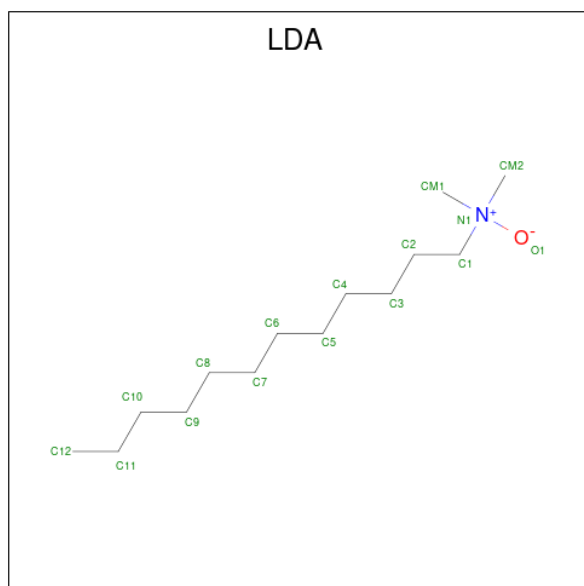
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q6D8U4
A	1	GLY	-	expression tag	UNP Q6D8U4
A	2	SER	-	expression tag	UNP Q6D8U4
A	3	SER	-	expression tag	UNP Q6D8U4
A	4	HIS	-	expression tag	UNP Q6D8U4
A	5	HIS	-	expression tag	UNP Q6D8U4
A	6	HIS	-	expression tag	UNP Q6D8U4
A	7	HIS	-	expression tag	UNP Q6D8U4
A	8	HIS	-	expression tag	UNP Q6D8U4
A	9	HIS	-	expression tag	UNP Q6D8U4
A	10	SER	-	expression tag	UNP Q6D8U4
A	11	SER	-	expression tag	UNP Q6D8U4
A	12	GLY	-	expression tag	UNP Q6D8U4
A	13	LEU	-	expression tag	UNP Q6D8U4
A	14	VAL	-	expression tag	UNP Q6D8U4
A	15	PRO	-	expression tag	UNP Q6D8U4
A	16	ARG	-	expression tag	UNP Q6D8U4
A	17	GLY	-	expression tag	UNP Q6D8U4
A	18	SER	-	expression tag	UNP Q6D8U4
A	19	HIS	-	expression tag	UNP Q6D8U4
A	20	MET	-	expression tag	UNP Q6D8U4
B	0	MET	-	initiating methionine	UNP Q6D8U4
B	1	GLY	-	expression tag	UNP Q6D8U4
B	2	SER	-	expression tag	UNP Q6D8U4
B	3	SER	-	expression tag	UNP Q6D8U4

Continued on next page...

Continued from previous page...

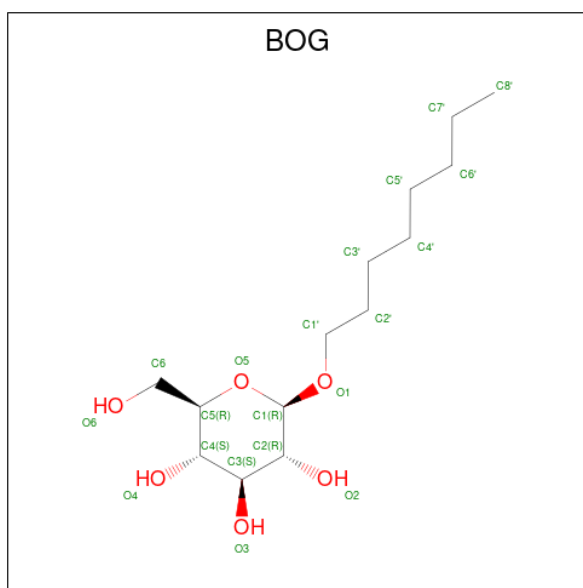
Chain	Residue	Modelled	Actual	Comment	Reference
B	4	HIS	-	expression tag	UNP Q6D8U4
B	5	HIS	-	expression tag	UNP Q6D8U4
B	6	HIS	-	expression tag	UNP Q6D8U4
B	7	HIS	-	expression tag	UNP Q6D8U4
B	8	HIS	-	expression tag	UNP Q6D8U4
B	9	HIS	-	expression tag	UNP Q6D8U4
B	10	SER	-	expression tag	UNP Q6D8U4
B	11	SER	-	expression tag	UNP Q6D8U4
B	12	GLY	-	expression tag	UNP Q6D8U4
B	13	LEU	-	expression tag	UNP Q6D8U4
B	14	VAL	-	expression tag	UNP Q6D8U4
B	15	PRO	-	expression tag	UNP Q6D8U4
B	16	ARG	-	expression tag	UNP Q6D8U4
B	17	GLY	-	expression tag	UNP Q6D8U4
B	18	SER	-	expression tag	UNP Q6D8U4
B	19	HIS	-	expression tag	UNP Q6D8U4
B	20	MET	-	expression tag	UNP Q6D8U4

- Molecule 2 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: $C_{14}H_{31}NO$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			32	28	2	2		
2	A	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 3 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			20	14	6		
3	A	1	Total	C	O	0	0
			20	14	6		
3	A	1	Total	C	O	0	0
			20	14	6		
3	B	1	Total	C	O	0	0
			20	14	6		

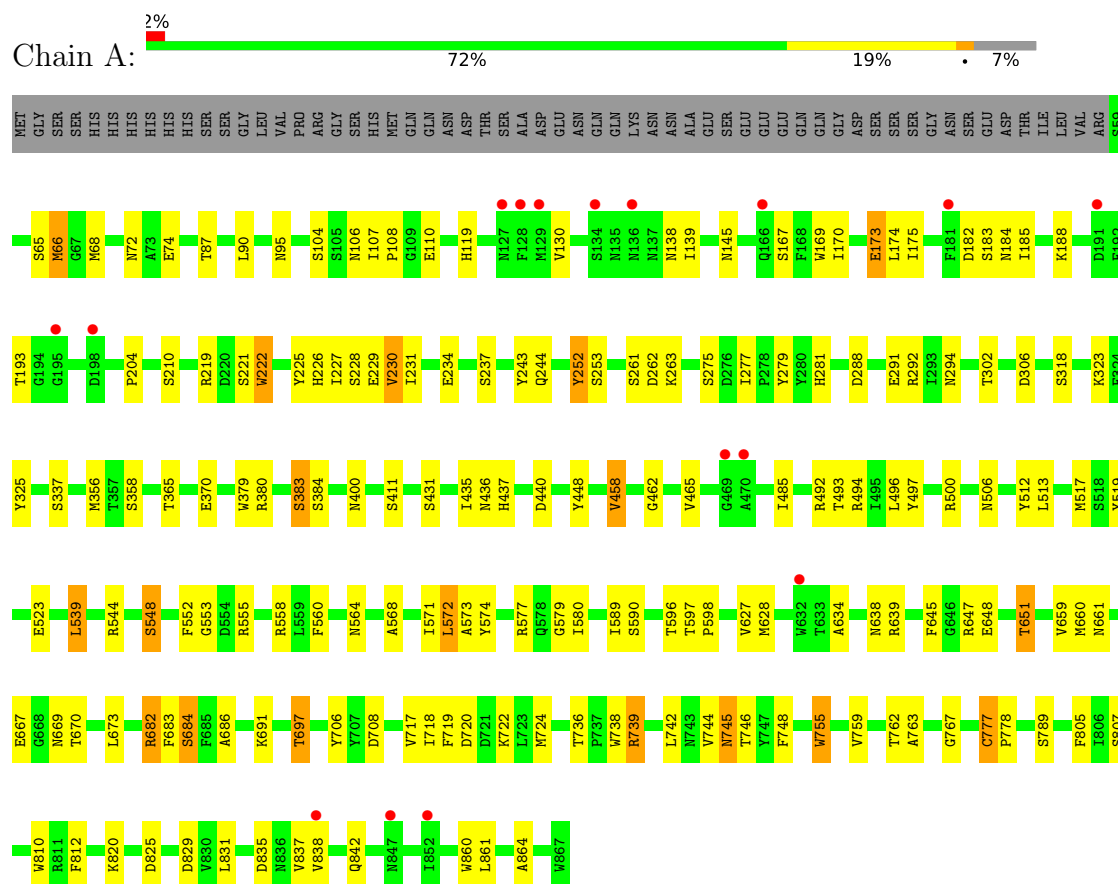
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		
4	B	1	Total	O	0	0
			1	1		

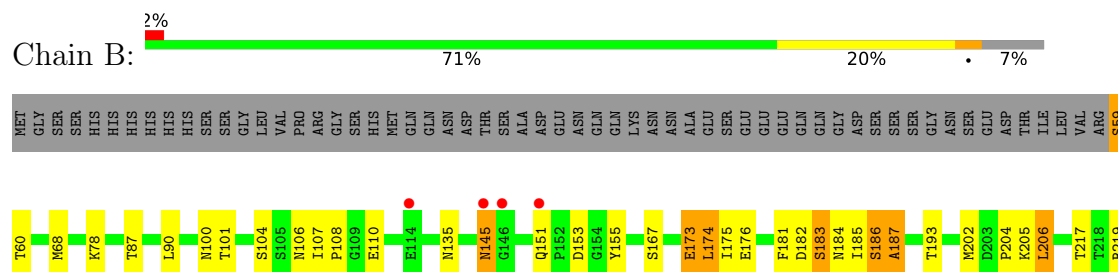
3 Residue-property plots

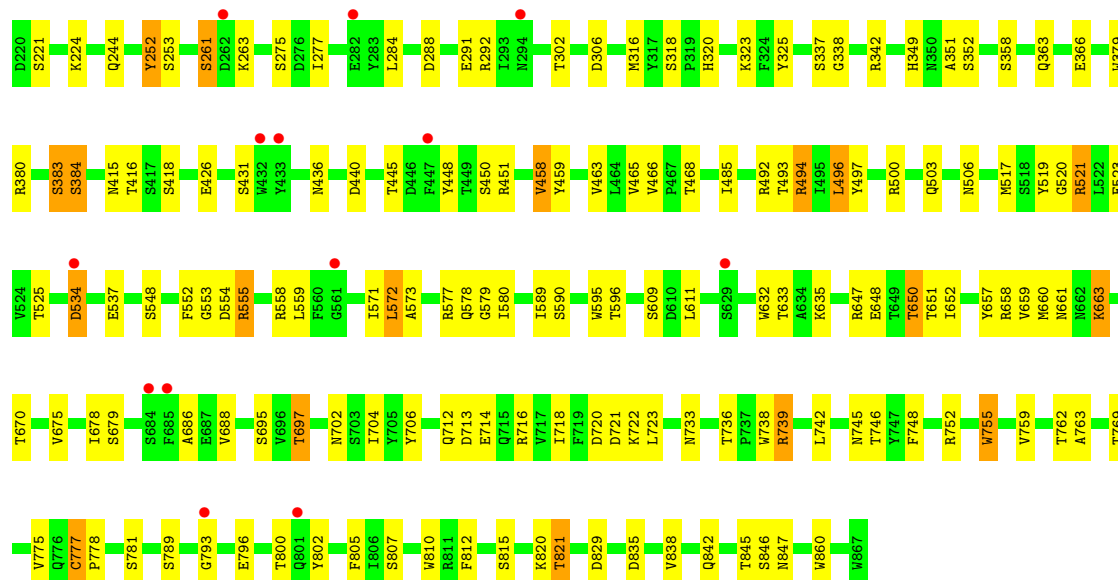
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Ferredoxin receptor



• Molecule 1: Ferredoxin receptor





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	137.27Å 79.89Å 137.91Å 90.00° 90.71° 90.00°	Depositor
Resolution (Å)	48.95 – 3.20 48.95 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.95-3.20) 99.9 (48.95-3.20)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.218 , 0.271 0.222 , 0.270	Depositor DCC
R_{free} test set	2511 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	74.4	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 38.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.006 for l,k,-h 0.026 for h,-k,-l 0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	13030	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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: BOG, LDA

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	0.59	0/6625	0.76	1/9003 (0.0%)
1	B	0.58	0/6625	0.78	3/9003 (0.0%)
All	All	0.59	0/13250	0.77	4/18006 (0.0%)

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	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	723	LEU	CA-CB-CG	6.57	130.41	115.30
1	B	572	LEU	CA-CB-CG	6.16	129.47	115.30
1	B	284	LEU	N-CA-C	-5.51	96.13	111.00
1	A	684	SER	N-CA-C	5.47	125.77	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	683	PHE	Peptide
1	A	777	CYS	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	B	552	PHE	Peptide
1	B	777	CYS	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	6450	0	6024	109	0
1	B	6450	0	6024	102	0
2	A	48	0	93	0	0
3	A	60	0	84	0	0
3	B	20	0	28	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	13030	0	12253	210	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 8.

All (210) 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:110:GLU:OE2	1:A:292:ARG:NH1	1.82	1.12
1:A:261:SER:HB2	1:A:263:LYS:H	1.51	0.75
1:B:521:ARG:HD2	1:B:553:GLY:HA2	1.69	0.73
1:B:523:GLU:OE1	1:B:548:SER:HB3	1.89	0.73
1:A:523:GLU:OE1	1:A:548:SER:HB3	1.89	0.72
1:A:574:TYR:CD1	1:A:660:MET:HE3	2.26	0.70
1:A:574:TYR:CE1	1:A:660:MET:HE2	2.28	0.68
1:A:219:ARG:HB2	1:A:221:SER:O	1.93	0.68
1:A:682:ARG:O	1:A:684:SER:HA	1.93	0.68
1:A:572:LEU:C	1:A:572:LEU:HD23	2.15	0.67
1:A:647:ARG:NH1	1:A:708:ASP:O	2.29	0.66
1:B:182:ASP:OD2	1:B:183:SER:OG	2.11	0.66
1:A:226:HIS:CD2	1:A:837:VAL:HG22	2.31	0.65
1:A:574:TYR:CD1	1:A:660:MET:CE	2.79	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:670:THR:HG22	1:B:697:THR:HG22	1.77	0.64
1:A:66:MET:HG3	1:A:560:PHE:CZ	2.33	0.64
1:B:78:LYS:HB3	1:B:745:ASN:HD22	1.61	0.63
1:A:670:THR:HG22	1:A:697:THR:HG22	1.80	0.63
1:B:739:ARG:HB2	1:B:762:THR:HG23	1.82	0.62
1:B:87:THR:HA	1:B:90:LEU:HD13	1.82	0.62
1:B:173:GLU:O	1:B:175:ILE:HG22	1.98	0.62
1:A:87:THR:HA	1:A:90:LEU:HD13	1.82	0.62
1:A:139:ILE:HA	1:A:572:LEU:CD2	2.30	0.61
1:B:648:GLU:HG3	1:B:661:ASN:HB3	1.83	0.61
1:A:745:ASN:N	1:A:745:ASN:HD22	1.99	0.60
1:B:261:SER:HB2	1:B:263:LYS:H	1.66	0.60
1:B:186:SER:O	1:B:187:ALA:CB	2.47	0.60
1:B:571:ILE:O	1:B:572:LEU:HB3	2.00	0.60
1:A:138:ASN:O	1:A:572:LEU:HD22	2.02	0.59
1:B:500:ARG:NH1	1:B:579:GLY:HA3	2.17	0.59
1:A:500:ARG:NH1	1:A:579:GLY:HA3	2.18	0.59
1:B:647:ARG:HD2	1:B:658:ARG:HD3	1.85	0.59
1:B:204:PRO:HG2	1:B:302:THR:HG23	1.83	0.58
1:B:217:THR:HG21	1:B:219:ARG:HH12	1.68	0.58
1:B:186:SER:O	1:B:187:ALA:HB3	2.02	0.58
1:B:555:ARG:HB2	1:B:555:ARG:CZ	2.33	0.58
1:A:574:TYR:CE1	1:A:660:MET:CE	2.86	0.58
1:A:648:GLU:HG3	1:A:661:ASN:HB3	1.85	0.58
1:A:108:PRO:O	1:A:244:GLN:OE1	2.22	0.58
1:A:718:ILE:O	1:A:722:LYS:O	2.22	0.58
1:A:229:GLU:O	1:A:231:ILE:N	2.37	0.57
1:B:755:TRP:HH2	1:B:810:TRP:CE2	2.22	0.57
1:A:755:TRP:HH2	1:A:810:TRP:CE2	2.22	0.57
1:A:500:ARG:HH12	1:A:579:GLY:HA3	1.70	0.57
1:A:173:GLU:O	1:A:175:ILE:HG22	2.03	0.57
1:A:572:LEU:HD23	1:A:572:LEU:O	2.04	0.57
1:A:596:THR:O	1:A:598:PRO:HD3	2.05	0.56
1:B:500:ARG:HH12	1:B:579:GLY:HA3	1.71	0.56
1:A:139:ILE:O	1:A:573:ALA:HB2	2.06	0.56
1:A:193:THR:HG23	1:A:568:ALA:HB2	1.88	0.55
1:A:458:VAL:HG22	1:A:497:TYR:HB2	1.88	0.55
1:A:227:ILE:HG22	1:A:227:ILE:O	2.07	0.55
1:B:244:GLN:HG2	1:B:277:ILE:HG21	1.88	0.55
1:B:217:THR:HG21	1:B:219:ARG:NH1	2.22	0.55
1:A:138:ASN:O	1:A:571:ILE:O	2.25	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:720:ASP:O	1:B:722:LYS:N	2.38	0.55
1:B:829:ASP:HB2	1:B:860:TRP:HB2	1.88	0.54
1:A:539:LEU:C	1:A:539:LEU:HD12	2.27	0.54
1:A:139:ILE:HA	1:A:572:LEU:HD22	1.89	0.54
1:B:185:ILE:HG22	1:B:186:SER:H	1.72	0.54
1:B:176:GLU:HG2	1:B:202:MET:HA	1.90	0.54
1:B:342:ARG:HE	1:B:363:GLN:HE21	1.57	0.53
1:A:65:SER:HA	1:A:68:MET:HG2	1.91	0.53
1:A:261:SER:HB2	1:A:263:LYS:N	2.24	0.53
1:A:400:ASN:O	1:A:400:ASN:ND2	2.41	0.53
1:B:688:VAL:HG12	1:B:746:THR:HG23	1.90	0.52
1:A:244:GLN:HG2	1:A:277:ILE:HG21	1.91	0.52
1:A:106:ASN:HB2	1:A:107:ILE:HG13	1.92	0.52
1:B:173:GLU:O	1:B:175:ILE:N	2.42	0.52
1:A:744:VAL:C	1:A:745:ASN:HD22	2.12	0.52
1:B:648:GLU:OE2	1:B:663:LYS:NZ	2.37	0.52
1:B:573:ALA:O	1:B:577:ARG:HD2	2.09	0.51
1:B:650:THR:HG23	1:B:657:TYR:HB2	1.91	0.51
1:A:221:SER:O	1:A:222:TRP:CB	2.59	0.51
1:A:589:ILE:HG13	1:A:590:SER:N	2.25	0.50
1:A:221:SER:O	1:A:222:TRP:CD1	2.64	0.50
1:A:755:TRP:CZ3	1:A:812:PHE:CZ	2.99	0.50
1:B:755:TRP:CZ3	1:B:812:PHE:CZ	2.99	0.50
1:A:436:ASN:O	1:A:517:MET:HA	2.11	0.50
1:B:78:LYS:HB3	1:B:745:ASN:ND2	2.27	0.50
1:A:173:GLU:O	1:A:175:ILE:N	2.45	0.50
1:B:100:ASN:OD1	1:B:733:ASN:OD1	2.30	0.50
1:A:767:GLY:HA3	1:A:842:GLN:OE1	2.11	0.50
1:A:288:ASP:HB3	1:A:325:TYR:CZ	2.46	0.49
1:A:448:TYR:CZ	1:A:506:ASN:HB2	2.47	0.49
1:A:596:THR:OG1	1:A:597:THR:N	2.44	0.49
1:B:151:GLN:HG2	1:B:155:TYR:CE2	2.47	0.49
1:A:739:ARG:HB2	1:A:762:THR:HG23	1.94	0.49
1:A:571:ILE:O	1:A:572:LEU:HD22	2.11	0.49
1:A:738:TRP:CE2	1:A:763:ALA:HB2	2.46	0.49
1:A:204:PRO:HG2	1:A:302:THR:HG23	1.93	0.49
1:A:745:ASN:N	1:A:745:ASN:ND2	2.60	0.49
1:B:416:THR:HG22	1:B:450:SER:OG	2.12	0.49
1:B:555:ARG:HB2	1:B:555:ARG:NH1	2.26	0.49
1:B:738:TRP:CE2	1:B:763:ALA:HB2	2.47	0.49
1:B:534:ASP:O	1:B:537:GLU:HG3	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:752:ARG:NH1	1:B:815:SER:O	2.46	0.49
1:A:188:LYS:HG3	1:A:512:TYR:CZ	2.48	0.49
1:A:130:VAL:HG11	1:A:170:ILE:HD12	1.95	0.48
1:A:107:ILE:CG2	1:A:110:GLU:HG3	2.44	0.48
1:B:458:VAL:HG22	1:B:497:TYR:HB2	1.96	0.48
1:A:370:GLU:HG3	1:A:411:SER:HB3	1.95	0.47
1:B:291:GLU:HB2	1:B:323:LYS:HB3	1.96	0.47
1:A:252:TYR:N	1:A:252:TYR:CD1	2.80	0.47
1:B:712:GLN:C	1:B:713:ASP:O	2.51	0.47
1:B:320:HIS:NE2	1:B:338:GLY:O	2.48	0.47
1:B:686:ALA:HB2	1:B:748:PHE:CD1	2.49	0.47
1:B:145:ASN:OD1	1:B:706:TYR:HB3	2.15	0.47
1:B:835:ASP:CG	1:B:835:ASP:O	2.53	0.47
1:A:742:LEU:HB3	1:A:759:VAL:HG12	1.97	0.47
1:B:577:ARG:O	1:B:580:ILE:HG22	2.15	0.47
1:B:802:TYR:OH	1:B:842:GLN:NE2	2.48	0.47
1:A:686:ALA:HB2	1:A:748:PHE:CD1	2.49	0.46
1:A:119:HIS:HD2	1:A:670:THR:CG2	2.28	0.46
1:B:106:ASN:HB2	1:B:107:ILE:HG13	1.97	0.46
1:B:100:ASN:O	1:B:101:THR:HG23	2.16	0.46
1:A:169:TRP:HE3	1:A:294:ASN:HD21	1.62	0.46
1:A:435:ILE:HG13	1:A:437:HIS:CE1	2.51	0.46
1:B:845:THR:OG1	1:B:847:ASN:HB3	2.16	0.46
1:B:100:ASN:OD1	1:B:733:ASN:HA	2.16	0.46
1:B:755:TRP:CE3	1:B:812:PHE:CE1	3.04	0.46
1:A:651:THR:HG22	1:A:651:THR:O	2.14	0.46
1:A:829:ASP:HB2	1:A:860:TRP:HB2	1.96	0.46
1:A:72:ASN:ND2	1:A:74:GLU:HG2	2.30	0.46
1:B:135:ASN:CG	1:B:135:ASN:O	2.54	0.46
1:B:181:PHE:CD1	1:B:181:PHE:N	2.83	0.46
1:A:95:ASN:ND2	1:A:182:ASP:OD2	2.48	0.45
1:A:227:ILE:O	1:A:229:GLU:N	2.48	0.45
1:A:221:SER:O	1:A:222:TRP:CG	2.70	0.45
1:B:820:LYS:O	1:B:821:THR:HG22	2.17	0.45
1:A:107:ILE:HA	1:A:244:GLN:OE1	2.17	0.45
1:B:805:PHE:CD2	1:B:838:VAL:HG11	2.52	0.45
1:A:234:GLU:O	1:A:237:SER:OG	2.31	0.44
1:B:275:SER:HB3	1:B:292:ARG:HB2	1.98	0.44
1:A:279:TYR:HE2	1:A:281:HIS:CE1	2.35	0.44
1:A:719:PHE:CD2	1:A:724:MET:HE3	2.51	0.44
1:B:451:ARG:HG2	1:B:503:GLN:HG3	1.98	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:775:VAL:O	1:B:793:GLY:HA2	2.17	0.44
1:A:755:TRP:HH2	1:A:810:TRP:CD2	2.35	0.44
1:B:695:SER:OG	1:B:739:ARG:NH1	2.50	0.44
1:B:553:GLY:O	1:B:555:ARG:N	2.41	0.44
1:A:139:ILE:HG12	1:A:572:LEU:HD21	1.98	0.44
1:A:519:TYR:O	1:A:519:TYR:CG	2.70	0.44
1:A:231:ILE:HD13	1:A:243:TYR:OH	2.17	0.44
1:A:825:ASP:HB2	1:A:864:ALA:HB3	1.98	0.44
1:A:230:VAL:HG22	1:A:231:ILE:N	2.33	0.43
1:B:108:PRO:O	1:B:244:GLN:OE1	2.36	0.43
1:B:379:TRP:CE2	1:B:485:ILE:HD11	2.53	0.43
1:A:462:GLY:HA3	1:A:493:THR:OG1	2.18	0.43
1:B:494:ARG:HB3	1:B:595:TRP:CZ3	2.53	0.43
1:A:107:ILE:HG22	1:A:110:GLU:CG	2.49	0.43
1:B:718:ILE:HB	1:B:796:GLU:HA	2.00	0.43
1:A:719:PHE:HD2	1:A:724:MET:HE3	1.84	0.43
1:B:205:LYS:O	1:B:206:LEU:CB	2.66	0.43
1:A:379:TRP:CE2	1:A:485:ILE:HD11	2.54	0.43
1:B:366:GLU:HG3	1:B:415:ASN:HB3	2.01	0.43
1:A:183:SER:O	1:A:184:ASN:C	2.56	0.43
1:B:383:SER:O	1:B:384:SER:C	2.57	0.43
1:B:652:ILE:O	1:B:652:ILE:HG22	2.18	0.43
1:A:638:ASN:HD22	1:A:669:ASN:ND2	2.16	0.43
1:A:746:THR:HB	1:A:755:TRP:HB3	2.01	0.43
1:B:379:TRP:CD2	1:B:485:ILE:HD11	2.54	0.43
1:B:448:TYR:CZ	1:B:506:ASN:HB2	2.54	0.43
1:A:275:SER:HB3	1:A:292:ARG:HB2	2.00	0.43
1:A:383:SER:O	1:A:384:SER:C	2.56	0.43
1:A:755:TRP:CE3	1:A:812:PHE:CE1	3.07	0.42
1:B:632:TRP:CG	1:B:675:VAL:HG12	2.55	0.42
1:B:805:PHE:CG	1:B:838:VAL:HG11	2.54	0.42
1:A:661:ASN:OD1	1:A:661:ASN:C	2.57	0.42
1:B:745:ASN:OD1	1:B:745:ASN:N	2.53	0.42
1:B:205:LYS:O	1:B:206:LEU:HG	2.20	0.42
1:B:523:GLU:OE2	1:B:525:THR:CG2	2.67	0.42
1:B:769:THR:HG21	1:B:846:SER:HA	2.01	0.42
1:B:288:ASP:HB3	1:B:325:TYR:CZ	2.54	0.42
1:B:519:TYR:O	1:B:520:GLY:C	2.58	0.42
1:B:589:ILE:HG13	1:B:590:SER:N	2.34	0.42
1:B:742:LEU:HB3	1:B:759:VAL:HG12	2.00	0.42
1:B:755:TRP:HH2	1:B:810:TRP:CD2	2.37	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:634:ALA:HB2	1:A:673:LEU:CD1	2.50	0.42
1:A:645:PHE:HA	1:A:661:ASN:O	2.19	0.42
1:A:831:LEU:HD23	1:A:831:LEU:HA	1.90	0.42
1:A:835:ASP:CG	1:A:835:ASP:O	2.58	0.42
1:B:746:THR:HB	1:B:755:TRP:HB3	2.02	0.42
1:A:435:ILE:HG22	1:A:519:TYR:HD1	1.85	0.42
1:B:380:ARG:HD3	1:B:383:SER:CB	2.50	0.42
1:B:349:HIS:CE1	1:B:351:ALA:HB2	2.55	0.41
1:A:262:ASP:HA	1:B:652:ILE:CD1	2.50	0.41
1:A:553:GLY:O	1:A:555:ARG:N	2.42	0.41
1:A:805:PHE:CD2	1:A:838:VAL:HG21	2.55	0.41
1:B:261:SER:HB2	1:B:263:LYS:N	2.34	0.41
1:B:611:LEU:HD21	1:B:660:MET:HB3	2.01	0.41
1:B:702:ASN:HB3	1:B:704:ILE:H	1.85	0.41
1:B:718:ILE:O	1:B:722:LYS:O	2.39	0.41
1:A:706:TYR:CE2	1:A:708:ASP:HB2	2.56	0.41
1:B:463:VAL:HG22	1:B:493:THR:HG23	2.03	0.41
1:A:627:VAL:HG13	1:A:628:MET:N	2.36	0.41
1:B:252:TYR:CD1	1:B:252:TYR:N	2.88	0.41
1:A:544:ARG:HD3	1:A:564:ASN:HB3	2.02	0.41
1:B:107:ILE:CG2	1:B:110:GLU:HG3	2.51	0.41
1:A:291:GLU:HB2	1:A:323:LYS:HB3	2.03	0.41
1:A:577:ARG:O	1:A:580:ILE:HG22	2.21	0.41
1:B:59:SER:OG	1:B:60:THR:N	2.51	0.41
1:A:225:TYR:O	1:A:226:HIS:ND1	2.54	0.40
1:B:459:TYR:CE1	1:B:496:LEU:HG	2.55	0.40
1:A:193:THR:O	1:A:639:ARG:NH1	2.54	0.40
1:B:436:ASN:O	1:B:517:MET:HA	2.22	0.40
1:A:380:ARG:HD3	1:A:383:SER:HB3	2.04	0.40
1:B:183:SER:O	1:B:184:ASN:C	2.59	0.40
1:B:466:VAL:O	1:B:468:THR:HG23	2.21	0.40
1:B:534:ASP:O	1:B:537:GLU:CG	2.69	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	807/868 (93%)	743 (92%)	55 (7%)	9 (1%)	12	44
1	B	807/868 (93%)	740 (92%)	56 (7%)	11 (1%)	9	40
All	All	1614/1736 (93%)	1483 (92%)	111 (7%)	20 (1%)	11	43

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	720	ASP
1	A	778	PRO
1	B	554	ASP
1	B	778	PRO
1	A	145	ASN
1	A	174	LEU
1	A	228	SER
1	A	820	LYS
1	B	145	ASN
1	B	174	LEU
1	B	187	ALA
1	B	206	LEU
1	B	384	SER
1	B	721	ASP
1	A	173	GLU
1	A	222	TRP
1	A	230	VAL
1	B	173	GLU
1	B	578	GLN
1	B	821	THR

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	702/754 (93%)	660 (94%)	42 (6%)	16	48
1	B	702/754 (93%)	645 (92%)	57 (8%)	9	36
All	All	1404/1508 (93%)	1305 (93%)	99 (7%)	12	42

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	MET
1	A	104	SER
1	A	167	SER
1	A	185	ILE
1	A	210	SER
1	A	252	TYR
1	A	253	SER
1	A	306	ASP
1	A	318	SER
1	A	337	SER
1	A	356	MET
1	A	358	SER
1	A	365	THR
1	A	383	SER
1	A	431	SER
1	A	440	ASP
1	A	458	VAL
1	A	465	VAL
1	A	492	ARG
1	A	494	ARG
1	A	496	LEU
1	A	513	LEU
1	A	539	LEU
1	A	548	SER
1	A	552	PHE
1	A	558	ARG
1	A	572	LEU
1	A	651	THR
1	A	659	VAL
1	A	667	GLU
1	A	682	ARG
1	A	691	LYS
1	A	697	THR
1	A	717	VAL
1	A	736	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	739	ARG
1	A	745	ASN
1	A	755	TRP
1	A	777	CYS
1	A	789	SER
1	A	807	SER
1	A	861	LEU
1	B	59	SER
1	B	68	MET
1	B	104	SER
1	B	153	ASP
1	B	167	SER
1	B	174	LEU
1	B	183	SER
1	B	186	SER
1	B	193	THR
1	B	221	SER
1	B	224	LYS
1	B	252	TYR
1	B	253	SER
1	B	261	SER
1	B	306	ASP
1	B	316	MET
1	B	318	SER
1	B	337	SER
1	B	352	SER
1	B	358	SER
1	B	383	SER
1	B	418	SER
1	B	426	GLU
1	B	431	SER
1	B	440	ASP
1	B	445	THR
1	B	458	VAL
1	B	465	VAL
1	B	492	ARG
1	B	494	ARG
1	B	496	LEU
1	B	521	ARG
1	B	534	ASP
1	B	555	ARG
1	B	558	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	559	LEU
1	B	596	THR
1	B	609	SER
1	B	633	THR
1	B	635	LYS
1	B	650	THR
1	B	651	THR
1	B	659	VAL
1	B	663	LYS
1	B	678	ILE
1	B	679	SER
1	B	697	THR
1	B	714	GLU
1	B	716	ARG
1	B	736	THR
1	B	739	ARG
1	B	755	TRP
1	B	777	CYS
1	B	781	SER
1	B	789	SER
1	B	800	THR
1	B	807	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	HIS
1	A	400	ASN
1	A	437	HIS
1	A	514	GLN
1	A	665	HIS
1	A	669	ASN
1	A	745	ASN
1	A	801	GLN
1	B	95	ASN
1	B	100	ASN
1	B	363	GLN
1	B	400	ASN
1	B	437	HIS
1	B	514	GLN
1	B	662	ASN
1	B	745	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	757	GLN
1	B	842	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

7 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	LDA	A	901[B]	-	12,15,15	2.15	1 (8%)	14,17,17	0.37	0
3	BOG	A	905	-	20,20,20	0.86	1 (5%)	25,25,25	1.86	3 (12%)
3	BOG	A	903	-	20,20,20	1.13	2 (10%)	25,25,25	1.62	6 (24%)
3	BOG	A	904	-	20,20,20	0.86	1 (5%)	25,25,25	1.80	2 (8%)
2	LDA	A	902	-	12,15,15	2.18	1 (8%)	14,17,17	0.82	1 (7%)
2	LDA	A	901[A]	-	12,15,15	2.11	1 (8%)	14,17,17	0.53	0
3	BOG	B	901	-	20,20,20	1.21	1 (5%)	25,25,25	2.62	6 (24%)

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	LDA	A	901[B]	-	-	8/13/13/13	-
3	BOG	A	905	-	-	5/11/31/31	0/1/1/1
3	BOG	A	903	-	-	6/11/31/31	0/1/1/1
3	BOG	A	904	-	-	4/11/31/31	0/1/1/1
2	LDA	A	902	-	-	3/13/13/13	-
2	LDA	A	901[A]	-	-	8/13/13/13	-
3	BOG	B	901	-	-	9/11/31/31	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	902	LDA	O1-N1	-7.47	1.24	1.42
2	A	901[B]	LDA	O1-N1	-7.39	1.24	1.42
2	A	901[A]	LDA	O1-N1	-7.27	1.25	1.42
3	B	901	BOG	O1-C1	4.94	1.48	1.40
3	A	903	BOG	O1-C1	3.86	1.46	1.40
3	A	903	BOG	O5-C1	3.10	1.49	1.41
3	A	905	BOG	O1-C1	2.97	1.45	1.40
3	A	904	BOG	O1-C1	2.19	1.43	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	901	BOG	O1-C1-C2	7.99	120.78	108.30
3	A	904	BOG	O5-C1-C2	7.81	126.88	110.35
3	A	905	BOG	O1-C1-C2	7.14	119.45	108.30
3	B	901	BOG	O5-C1-C2	-6.98	95.58	110.35
3	B	901	BOG	O5-C1-O1	-4.74	98.76	109.97
3	A	903	BOG	O1-C1-C2	3.95	114.46	108.30
3	A	905	BOG	C1'-O1-C1	3.87	120.25	113.84
3	B	901	BOG	C4-C3-C2	3.61	117.12	110.82
3	A	903	BOG	C4-C3-C2	-3.54	104.64	110.82
3	A	903	BOG	O5-C1-C2	2.73	116.12	110.35
3	A	903	BOG	C1'-O1-C1	-2.66	109.43	113.84
3	A	904	BOG	C1-O5-C5	-2.54	108.71	113.69
3	A	903	BOG	O5-C5-C4	2.23	113.74	109.69
2	A	902	LDA	CM2-N1-C1	2.21	114.88	110.23
3	B	901	BOG	C1'-O1-C1	-2.11	110.35	113.84
3	A	903	BOG	C1-C2-C3	-2.09	105.63	110.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	905	BOG	C3-C4-C5	2.09	113.96	110.24
3	B	901	BOG	O3-C3-C2	-2.05	105.60	110.35

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901[A]	LDA	C2-C1-N1-O1
2	A	901[A]	LDA	C2-C1-N1-CM2
2	A	902	LDA	N1-C1-C2-C3
3	B	901	BOG	O5-C1-O1-C1'
3	A	905	BOG	O5-C5-C6-O6
3	A	905	BOG	C4-C5-C6-O6
3	B	901	BOG	O5-C5-C6-O6
3	A	903	BOG	O5-C5-C6-O6
3	A	903	BOG	C4-C5-C6-O6
3	B	901	BOG	C4-C5-C6-O6
3	B	901	BOG	O1-C1'-C2'-C3'
2	A	901[A]	LDA	C3-C4-C5-C6
2	A	901[B]	LDA	C4-C5-C6-C7
2	A	902	LDA	C11-C10-C9-C8
2	A	901[B]	LDA	C6-C7-C8-C9
3	B	901	BOG	C2-C1-O1-C1'
2	A	901[A]	LDA	C5-C6-C7-C8
3	A	903	BOG	C3'-C4'-C5'-C6'
2	A	901[B]	LDA	C11-C10-C9-C8
3	A	904	BOG	C2'-C1'-O1-C1
3	A	905	BOG	C2'-C3'-C4'-C5'
3	B	901	BOG	C1'-C2'-C3'-C4'
2	A	901[B]	LDA	C1-C2-C3-C4
3	A	905	BOG	C4'-C5'-C6'-C7'
2	A	902	LDA	C1-C2-C3-C4
3	B	901	BOG	C2'-C3'-C4'-C5'
2	A	901[B]	LDA	C5-C6-C7-C8
2	A	901[A]	LDA	C9-C10-C11-C12
3	B	901	BOG	C5'-C6'-C7'-C8'
2	A	901[A]	LDA	C7-C8-C9-C10
2	A	901[B]	LDA	C3-C4-C5-C6
3	A	904	BOG	O1-C1'-C2'-C3'
3	A	904	BOG	C4-C5-C6-O6
2	A	901[A]	LDA	C2-C1-N1-CM1
3	A	903	BOG	C5'-C6'-C7'-C8'

Continued on next page...

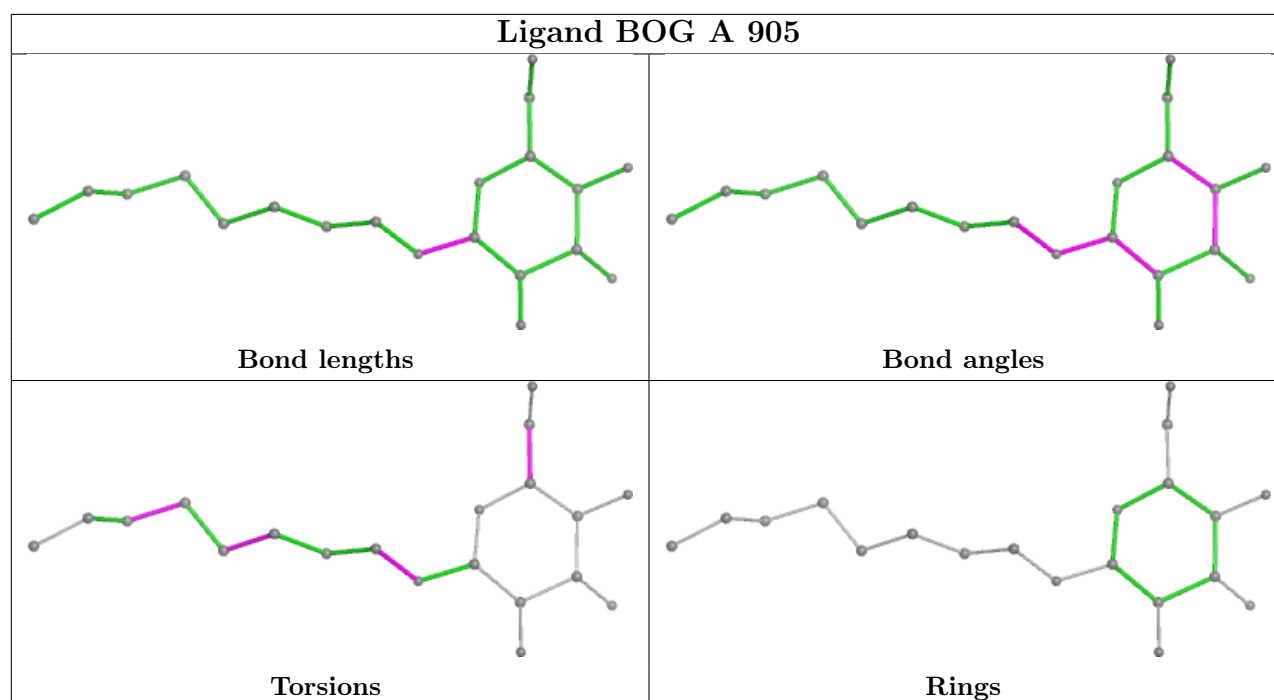
Continued from previous page...

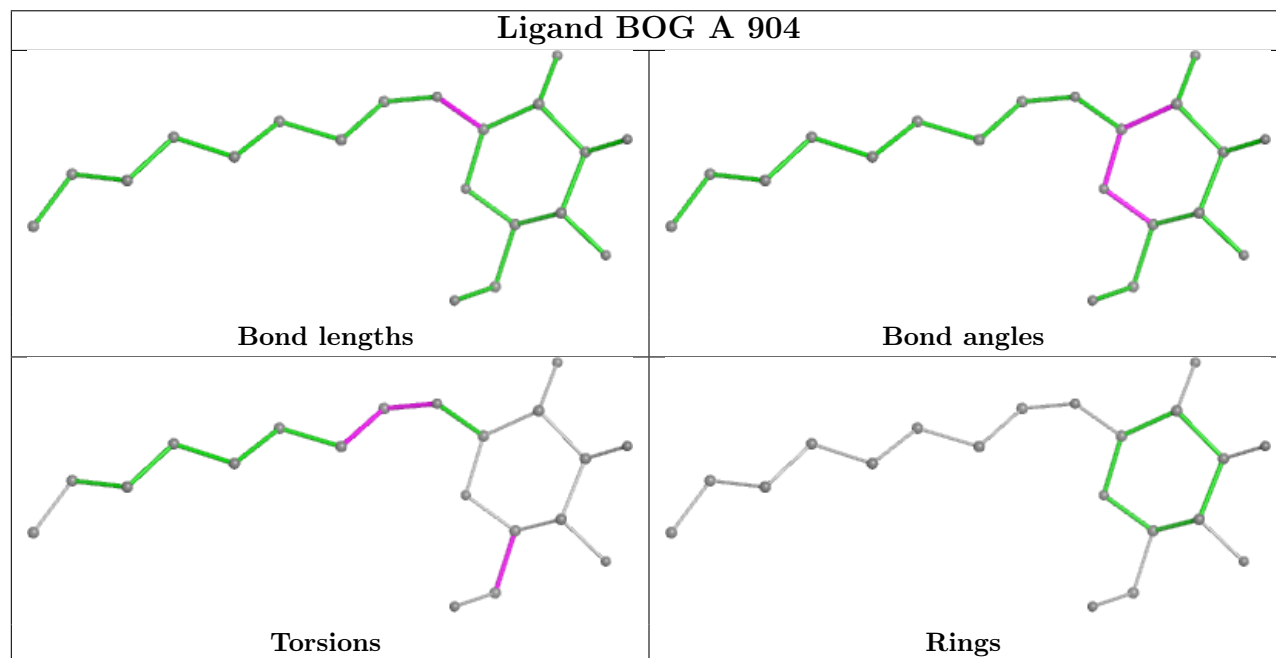
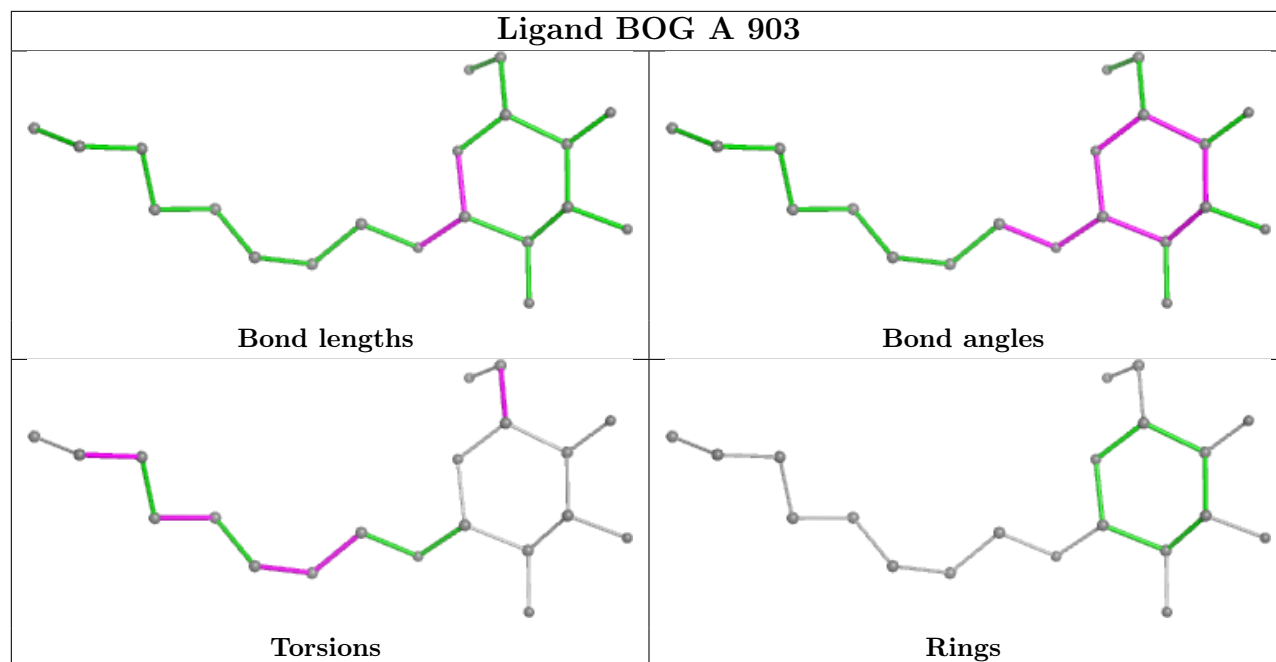
Mol	Chain	Res	Type	Atoms
3	A	903	BOG	C1'-C2'-C3'-C4'
2	A	901[B]	LDA	C9-C10-C11-C12
2	A	901[A]	LDA	C4-C5-C6-C7
3	A	903	BOG	O1-C1'-C2'-C3'
3	B	901	BOG	C3'-C4'-C5'-C6'
3	A	904	BOG	O5-C5-C6-O6
2	A	901[B]	LDA	C7-C8-C9-C10
3	A	905	BOG	C2'-C1'-O1-C1

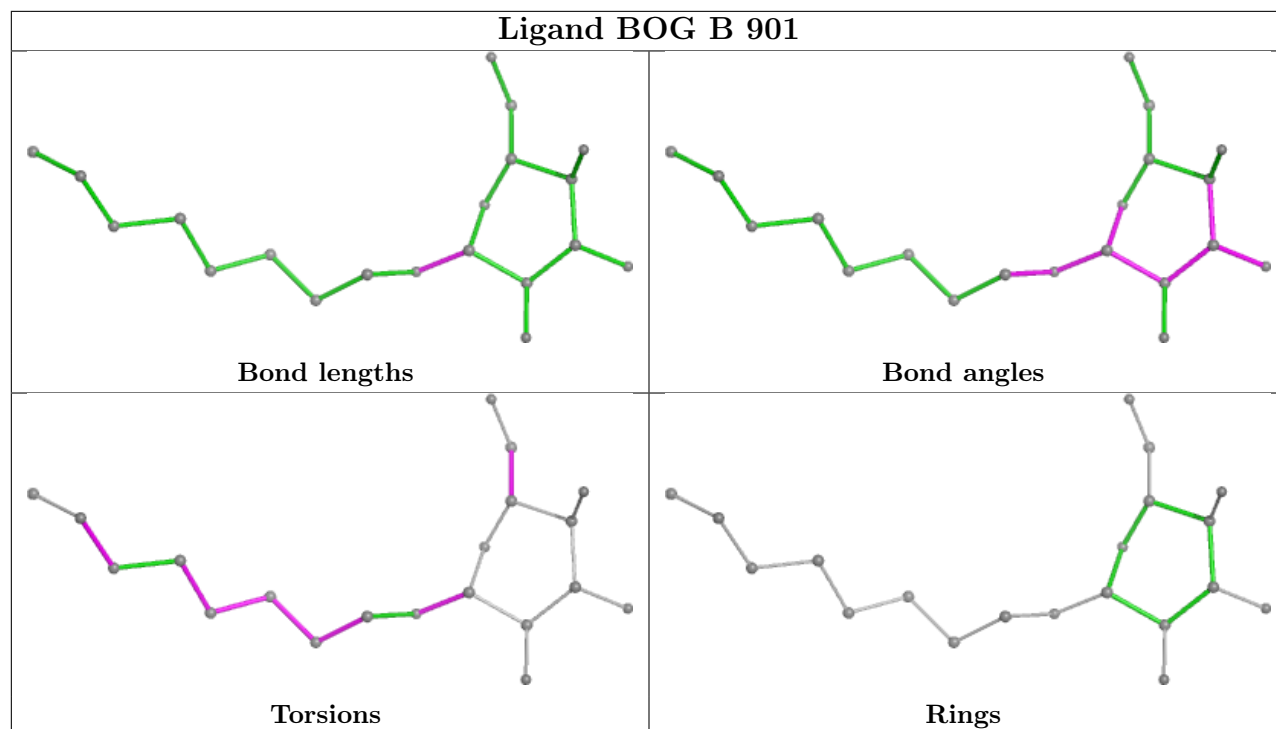
There are no ring outliers.

No monomer is involved in short contacts.

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.







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

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	809/868 (93%)	-0.10	16 (1%)	64 49	42, 75, 114, 151	0
1	B	809/868 (93%)	-0.04	17 (2%)	63 47	46, 85, 120, 165	0
All	All	1618/1736 (93%)	-0.07	33 (2%)	64 49	42, 80, 118, 165	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	847	ASN	4.9
1	A	134	SER	4.4
1	B	114	GLU	4.2
1	B	262	ASP	4.1
1	A	195	GLY	3.8
1	B	433	TYR	3.8
1	A	166	GLN	3.5
1	A	127	ASN	3.5
1	B	629	SER	3.4
1	A	129	MET	3.3
1	B	432	TRP	3.3
1	A	191	ASP	3.0
1	B	801	GLN	3.0
1	A	470	ALA	3.0
1	B	146	GLY	2.8
1	B	684	SER	2.8
1	B	282	GLU	2.8
1	A	469	GLY	2.8
1	A	852	ILE	2.6
1	B	793	GLY	2.5
1	B	447	PHE	2.5
1	A	128	PHE	2.4
1	B	145	ASN	2.4
1	B	294	ASN	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	685	PHE	2.3
1	A	198	ASP	2.3
1	B	534	ASP	2.2
1	A	181	PHE	2.2
1	B	151	GLN	2.2
1	A	632	TRP	2.2
1	B	561	GLY	2.1
1	A	838	VAL	2.1
1	A	136	ASN	2.1

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](#)

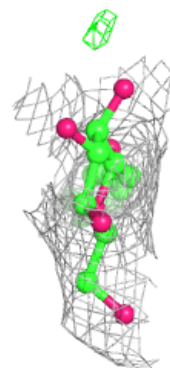
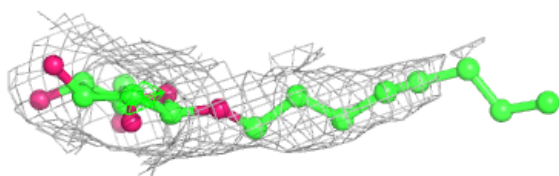
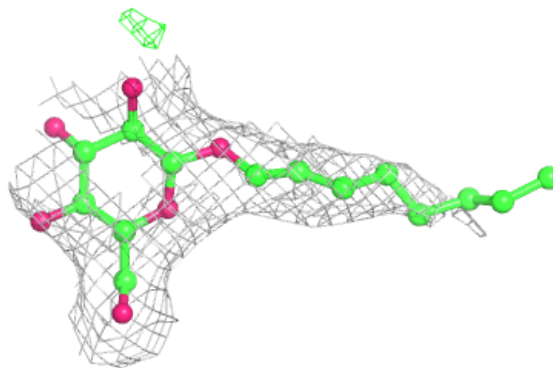
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	LDA	A	902	16/16	0.80	0.35	92,118,155,159	0
2	LDA	A	901[B]	16/16	0.82	0.50	68,87,91,92	16
2	LDA	A	901[A]	16/16	0.82	0.50	70,94,103,105	16
3	BOG	A	905	20/20	0.87	0.15	117,134,152,157	0
3	BOG	A	904	20/20	0.89	0.15	82,127,142,142	0
3	BOG	B	901	20/20	0.94	0.11	81,88,96,96	0
3	BOG	A	903	20/20	0.95	0.11	63,77,86,95	0

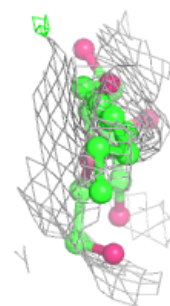
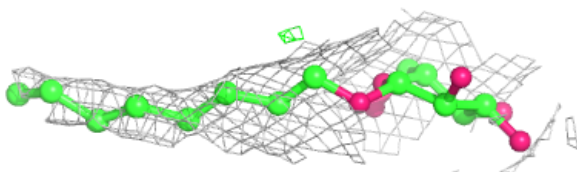
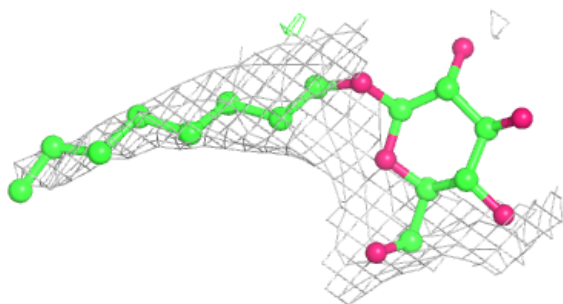
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BOG A 905:

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

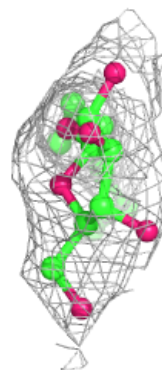
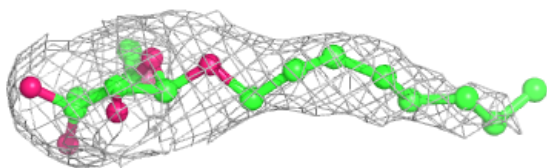
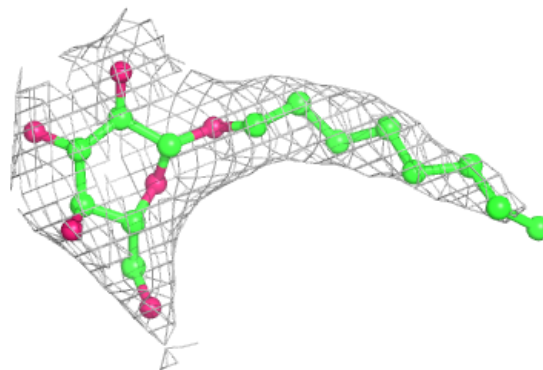
**Electron density around BOG A 904:**

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

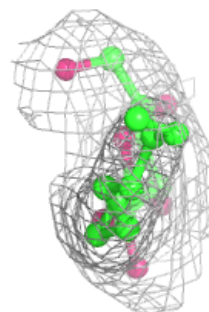
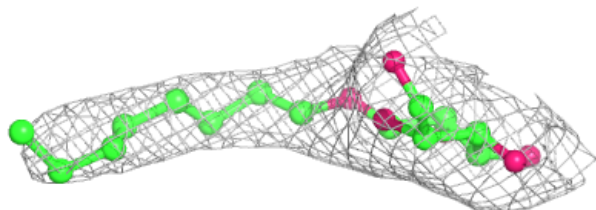
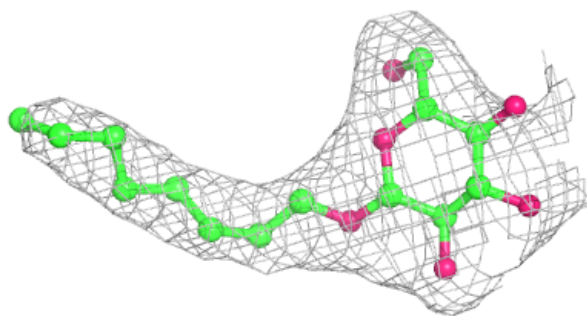


Electron density around BOG B 901:

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

**Electron density around BOG A 903:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.