



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:55 AM GMT

PDB ID : 3CNV
Title : Crystal structure of the ligand-binding domain of a putative GntR-family transcriptional regulator from Bordetella bronchiseptica
Authors : Zimmerman, M.D.; Xu, X.; Cui, H.; Filippova, E.V.; Savchenko, A.; Edwards, A.M.; Joachimiak, A.; Minor, W.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2008-03-26
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

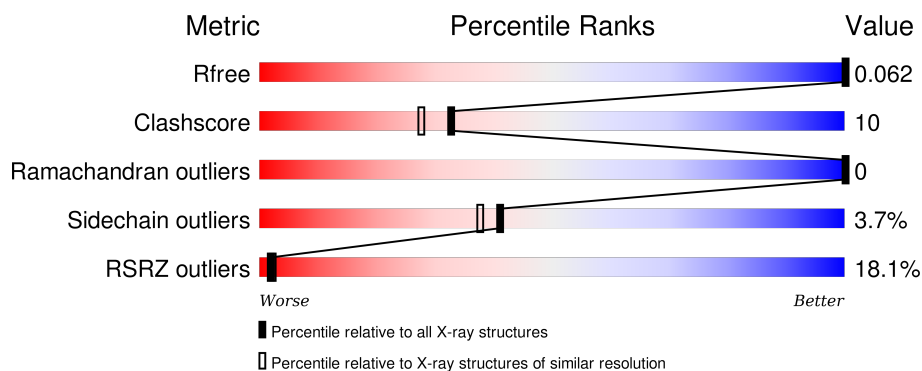
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	162	<div> <div>17%</div> <div>83%</div> <div>13%</div> <div>• •</div> </div>
1	B	162	<div> <div>17%</div> <div>75%</div> <div>20%</div> <div>• •</div> </div>
1	C	162	<div> <div>20%</div> <div>84%</div> <div>14%</div> <div>•</div> </div>
1	D	162	<div> <div>15%</div> <div>83%</div> <div>9%</div> <div>• • •</div> </div>

2 Entry composition [i](#)

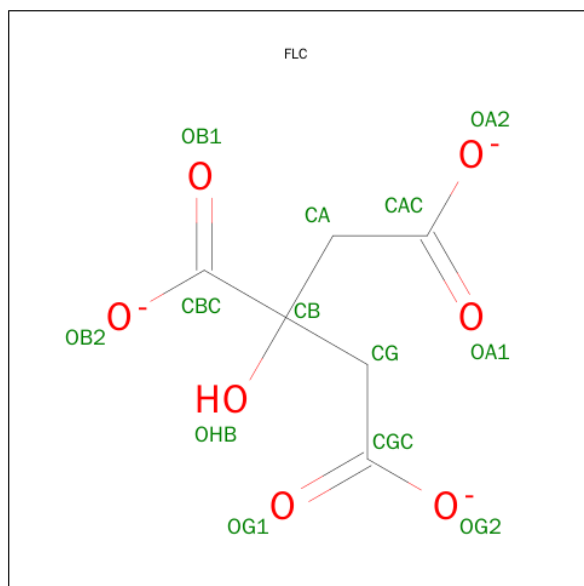
There are 4 unique types of molecules in this entry. The entry contains 5485 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Putative GntR-family transcriptional regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	Se	0	1	0
			1230	782	225	219	1	3			
1	B	159	Total	C	N	O	S	Se	1	1	0
			1268	802	235	226	1	4			
1	C	162	Total	C	N	O	S	Se	0	2	0
			1289	814	243	228	1	3			
1	D	157	Total	C	N	O	S	Se	0	1	0
			1246	790	235	216	1	4			

- Molecule 2 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		
3	A	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		

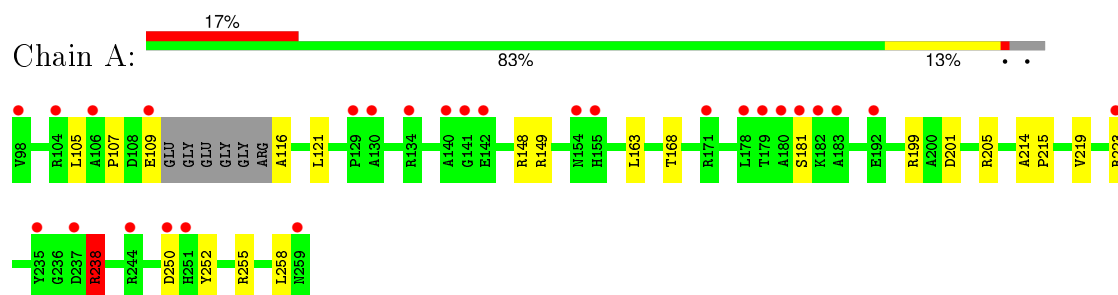
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	98	Total	O	0	0
			98	98		
4	B	105	Total	O	0	0
			105	105		
4	C	98	Total	O	0	0
			98	98		
4	D	94	Total	O	0	0
			94	94		

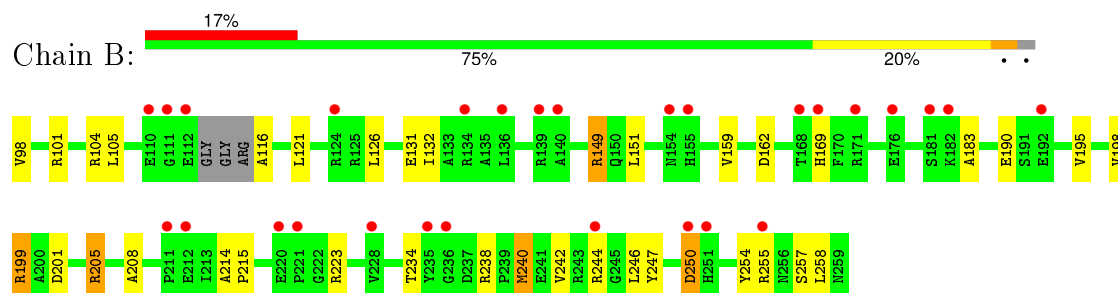
3 Residue-property plots

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.

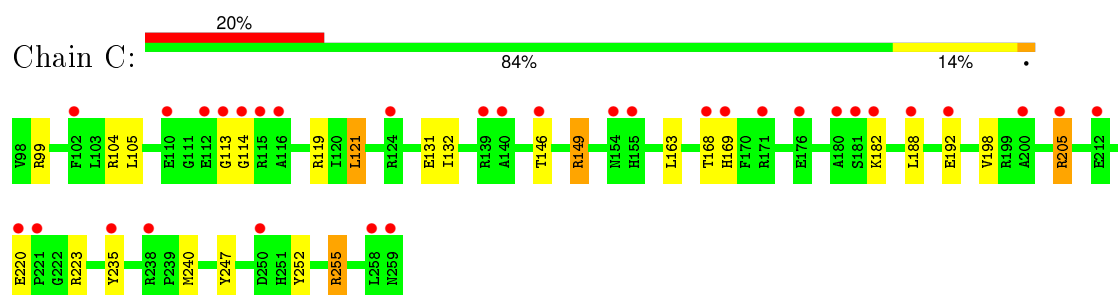
- Molecule 1: Putative GntR-family transcriptional regulator



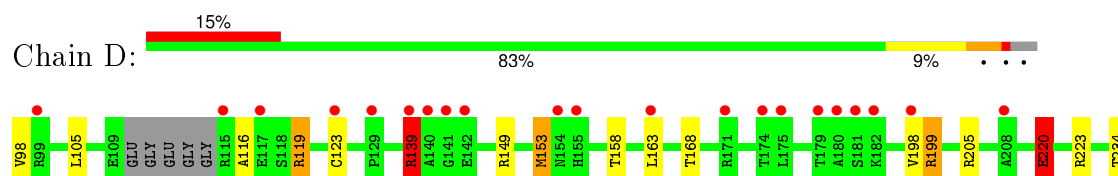
- Molecule 1: Putative GntR-family transcriptional regulator

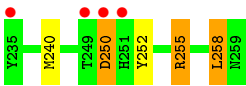


- Molecule 1: Putative GntR-family transcriptional regulator



- Molecule 1: Putative GntR-family transcriptional regulator





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.20Å 58.07Å 65.58Å 82.36° 86.52° 68.31°	Depositor
Resolution (Å)	29.91 – 2.00 29.91 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (29.91-2.00) 95.9 (29.91-2.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3149.24 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.4.0069	Depositor
R, R_{free}	0.164 , 0.217 0.307 , 0.062	Depositor DCC
R_{free} test set	2710 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.61$, $\langle L^2 \rangle = 0.47$	Xtriage
Outliers	0 of 53378 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5485	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.98% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	1.05	0/1254	1.08	3/1696 (0.2%)
1	B	1.13	4/1292 (0.3%)	1.14	11/1742 (0.6%)
1	C	1.11	2/1317 (0.2%)	1.08	7/1776 (0.4%)
1	D	1.05	2/1270 (0.2%)	1.04	6/1715 (0.3%)
All	All	1.09	8/5133 (0.2%)	1.09	27/6929 (0.4%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	149	ARG	CB-CG	-7.73	1.31	1.52
1	D	123	CYS	CB-SG	6.56	1.93	1.82
1	B	208	ALA	CA-CB	6.42	1.66	1.52
1	B	190	GLU	CG-CD	5.49	1.60	1.51
1	C	220	GLU	CD-OE1	5.48	1.31	1.25
1	B	242	VAL	CB-CG1	5.48	1.64	1.52
1	D	220	GLU	CG-CD	5.32	1.59	1.51
1	C	149	ARG	CB-CG	-5.26	1.38	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	199	ARG	NE-CZ-NH1	-10.05	115.28	120.30
1	B	101	ARG	NE-CZ-NH1	8.59	124.59	120.30
1	A	163	LEU	CA-CB-CG	7.61	132.79	115.30
1	B	101	ARG	NE-CZ-NH2	-7.48	116.56	120.30
1	D	139	ARG	NE-CZ-NH1	-7.32	116.64	120.30
1	D	163	LEU	CA-CB-CG	7.18	131.81	115.30
1	D	199	ARG	NE-CZ-NH1	-6.16	117.22	120.30
1	D	255	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	D	250	ASP	CB-CG-OD1	5.92	123.63	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	250	ASP	CB-CG-OD1	5.78	123.50	118.30
1	C	121	LEU	CB-CG-CD1	5.76	120.80	111.00
1	A	238	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	A	148	ARG	NE-CZ-NH2	-5.73	117.43	120.30
1	C	255	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	B	240[A]	MSE	CG-SE-CE	-5.68	86.41	98.90
1	B	240[B]	MSE	CG-SE-CE	-5.68	86.41	98.90
1	C	163	LEU	CA-CB-CG	5.66	128.32	115.30
1	B	126	LEU	CB-CG-CD2	-5.58	101.51	111.00
1	B	199	ARG	NE-CZ-NH2	5.56	123.08	120.30
1	C	119	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	C	99	ARG	NE-CZ-NH1	-5.45	117.58	120.30
1	C	119	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	B	205	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	C	255	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	B	162	ASP	CB-CG-OD1	-5.20	113.62	118.30
1	B	149	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	D	255	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	1230	0	1233	17	0
1	B	1268	0	1279	52	0
1	C	1289	0	1301	25	0
1	D	1246	0	1260	28	0
2	A	13	0	5	0	0
2	B	13	0	5	0	0
2	C	13	0	5	0	0
2	D	13	0	5	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	1	0	0	0	0
4	A	98	0	0	3	0
4	B	105	0	0	6	0
4	C	98	0	0	11	0
4	D	94	0	0	6	0
All	All	5485	0	5093	106	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 10.

All (106) 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:C:205[B]:ARG:HG3	1:C:205[B]:ARG:HH11	1.02	1.09
1:A:258[A]:LEU:HD22	1:B:104:ARG:NH2	1.69	1.07
1:B:244:ARG:HH12	1:B:246:LEU:CD1	1.71	1.03
1:C:205[B]:ARG:CG	1:C:205[B]:ARG:HH11	1.73	1.01
1:D:223:ARG:NH1	1:D:250:ASP:OD2	1.96	0.98
1:A:105:LEU:HD22	1:A:252:TYR:HB3	1.45	0.97
1:B:198:VAL:HG23	1:D:255:ARG:NE	1.83	0.93
1:C:205[B]:ARG:NH1	1:C:205[B]:ARG:HG3	1.84	0.92
1:C:169:HIS:ND1	1:C:240:MSE:HE1	1.91	0.85
1:B:199:ARG:NH1	1:B:201:ASP:OD1	2.12	0.83
1:D:116:ALA:HB2	1:D:153:MSE:HE3	1.66	0.78
1:B:244:ARG:NH1	1:B:246:LEU:CD1	2.46	0.77
1:B:244:ARG:HH12	1:B:246:LEU:HD12	1.52	0.73
1:B:198:VAL:CG2	1:D:255:ARG:CD	2.67	0.72
1:B:244:ARG:NH1	1:B:246:LEU:HD11	2.04	0.72
1:C:205[B]:ARG:CD	4:C:343:HOH:O	2.38	0.71
1:B:199:ARG:HH12	1:B:201:ASP:CG	1.94	0.71
1:B:169:HIS:CG	1:B:240[B]:MSE:HE1	2.28	0.69
1:C:205[B]:ARG:NH1	1:C:205[B]:ARG:CG	2.39	0.68
1:A:199:ARG:NH1	1:A:201:ASP:OD2	2.27	0.67
1:B:198:VAL:HG23	1:D:255:ARG:CD	2.25	0.66
1:C:113:GLY:O	1:C:114:GLY:C	2.32	0.66
1:B:255:ARG:CD	1:D:198:VAL:CG2	2.74	0.65
1:A:168:THR:HG23	4:A:362:HOH:O	1.96	0.65
1:C:205[B]:ARG:HD2	4:C:343:HOH:O	1.95	0.64
1:D:119:ARG:HG3	4:D:371:HOH:O	1.98	0.64
1:B:223:ARG:NH1	1:B:250:ASP:OD2	2.30	0.64
1:D:105:LEU:HD22	1:D:252:TYR:HB3	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:ARG:HD2	1:D:198:VAL:HG22	1.79	0.64
1:B:234:THR:CG2	1:B:240[A]:MSE:HG3	2.29	0.62
1:C:235:TYR:CD2	4:C:338:HOH:O	2.51	0.62
1:B:116:ALA:N	4:B:363:HOH:O	2.31	0.62
1:B:238:ARG:HB3	1:B:240[A]:MSE:HE2	1.82	0.61
1:C:205[B]:ARG:HD3	4:C:343:HOH:O	2.01	0.60
1:C:235:TYR:HD2	4:C:338:HOH:O	1.84	0.59
1:B:244:ARG:HH12	1:B:246:LEU:HD11	1.54	0.58
1:A:223:ARG:NH1	1:A:250:ASP:OD2	2.36	0.58
1:B:199:ARG:NH1	1:B:201:ASP:CG	2.56	0.58
1:B:255:ARG:NE	1:D:198:VAL:CG2	2.66	0.58
1:C:168:THR:HG23	4:C:336:HOH:O	2.04	0.58
1:C:255:ARG:HG3	1:C:255:ARG:HH11	1.69	0.58
1:B:234:THR:HG23	1:B:240[A]:MSE:HG3	1.86	0.57
1:C:169:HIS:CG	1:C:240:MSE:HE1	2.40	0.57
1:C:188:LEU:HD12	1:C:192:GLU:HG2	1.85	0.57
1:B:255:ARG:CD	1:D:198:VAL:HG22	2.36	0.55
1:B:255:ARG:HD2	1:D:198:VAL:CG2	2.36	0.55
1:A:107:PRO:HB2	1:A:109:GLU:HG3	1.90	0.54
1:B:255:ARG:HH11	1:B:255:ARG:HG3	1.72	0.54
1:D:168:THR:HG23	4:D:357:HOH:O	2.07	0.53
1:B:198:VAL:CG2	1:D:255:ARG:HD2	2.39	0.53
1:A:105:LEU:HD22	1:A:252:TYR:CB	2.30	0.52
1:D:205:ARG:NE	4:D:382:HOH:O	2.28	0.52
1:A:255:ARG:HH11	1:A:255:ARG:HG3	1.75	0.52
1:B:244:ARG:NH1	1:B:246:LEU:HG	2.25	0.51
1:B:255:ARG:NE	1:D:198:VAL:HG23	2.24	0.51
1:A:258[A]:LEU:HD22	1:B:104:ARG:HH22	1.66	0.50
1:C:223:ARG:HD2	4:C:335:HOH:O	2.11	0.50
1:C:114:GLY:HA3	4:C:391:HOH:O	2.10	0.50
1:B:195:VAL:HG13	1:B:234:THR:HG21	1.92	0.49
1:C:168:THR:HB	4:C:378:HOH:O	2.11	0.49
1:C:104:ARG:HG3	1:C:104:ARG:NH1	2.26	0.49
1:B:240[A]:MSE:HE1	4:B:351:HOH:O	2.12	0.49
1:B:205:ARG:HG2	4:B:328:HOH:O	2.11	0.49
1:B:244:ARG:HH12	1:B:246:LEU:CG	2.24	0.48
1:C:205[B]:ARG:CB	1:C:205[B]:ARG:NH1	2.76	0.48
1:B:205:ARG:NH1	4:B:328:HOH:O	2.41	0.48
1:A:214:ALA:N	1:A:215:PRO:CD	2.77	0.48
1:B:214:ALA:HB3	1:B:215:PRO:HD3	1.96	0.48
1:C:146:THR:HG23	4:C:383:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:ARG:NH1	1:B:255:ARG:HG3	2.29	0.47
1:C:105:LEU:HD22	1:C:252:TYR:HB3	1.97	0.47
1:B:214:ALA:N	1:B:215:PRO:HD2	2.29	0.47
1:B:151:LEU:HB2	1:B:159:VAL:HB	1.96	0.47
1:B:244:ARG:NH1	1:B:246:LEU:CG	2.79	0.46
1:B:98:VAL:N	4:B:330:HOH:O	2.47	0.46
1:D:205:ARG:NH1	4:D:324:HOH:O	2.38	0.46
1:B:183:ALA:O	4:B:344:HOH:O	2.20	0.46
1:B:240[A]:MSE:CE	1:B:240[A]:MSE:CG	2.94	0.46
1:B:234:THR:HG21	1:B:240[A]:MSE:HG3	1.97	0.46
1:D:98:VAL:N	4:D:328:HOH:O	2.48	0.46
1:D:220:GLU:CD	1:D:220:GLU:H	2.20	0.45
1:B:234:THR:CG2	1:B:240[B]:MSE:HG3	2.46	0.45
1:D:258:LEU:HD23	1:D:258:LEU:N	2.30	0.45
1:B:214:ALA:N	1:B:215:PRO:CD	2.80	0.44
1:B:131:GLU:HG2	1:B:132:ILE:N	2.31	0.44
1:A:238:ARG:HG2	1:A:238:ARG:H	1.68	0.43
1:A:199:ARG:NH1	1:A:201:ASP:CG	2.71	0.43
1:A:199:ARG:HD2	1:A:199:ARG:HH11	1.67	0.43
1:A:205:ARG:NH1	4:A:330:HOH:O	2.43	0.43
1:B:247:TYR:CD1	1:B:247:TYR:N	2.87	0.43
1:B:257:SER:OG	1:D:199:ARG:HD3	2.19	0.42
1:C:131:GLU:HG2	1:C:132:ILE:H	1.84	0.42
1:D:139:ARG:HB3	1:D:139:ARG:HE	1.18	0.42
1:B:247:TYR:HD1	1:B:247:TYR:N	2.18	0.42
1:A:116:ALA:N	4:A:376:HOH:O	2.53	0.42
1:C:205[A]:ARG:NH1	4:C:343:HOH:O	2.51	0.41
1:B:198:VAL:HG23	1:D:255:ARG:HE	1.79	0.41
1:D:205:ARG:CD	4:D:382:HOH:O	2.68	0.41
1:B:198:VAL:HG22	1:D:255:ARG:HD2	2.02	0.41
1:A:214:ALA:HB1	1:A:219:VAL:O	2.21	0.41
1:B:255:ARG:NE	1:D:198:VAL:HG22	2.35	0.40
1:D:105:LEU:CD1	1:D:158:THR:CG2	2.99	0.40
1:D:234:THR:HG22	1:D:240:MSE:HG2	2.03	0.40
1:C:198:VAL:HG11	1:C:235:TYR:CZ	2.55	0.40
1:B:258:LEU:HD12	1:B:258:LEU:N	2.37	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	153/162 (94%)	152 (99%)	1 (1%)	0	100	100
1	B	156/162 (96%)	155 (99%)	1 (1%)	0	100	100
1	C	162/162 (100%)	158 (98%)	4 (2%)	0	100	100
1	D	154/162 (95%)	154 (100%)	0	0	100	100
All	All	625/648 (96%)	619 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	127/132 (96%)	123 (97%)	4 (3%)	47	46
1	B	132/132 (100%)	128 (97%)	4 (3%)	48	47
1	C	132/132 (100%)	126 (96%)	6 (4%)	34	29
1	D	128/132 (97%)	122 (95%)	6 (5%)	32	27
All	All	519/528 (98%)	499 (96%)	20 (4%)	41	35

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	LEU
1	A	149	ARG

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Mol	Chain	Res	Type
1	A	181	SER
1	A	238	ARG
1	B	105	LEU
1	B	121	LEU
1	B	149	ARG
1	B	254	TYR
1	C	121	LEU
1	C	149	ARG
1	C	182	LYS
1	C	205[A]	ARG
1	C	205[B]	ARG
1	C	247	TYR
1	D	119	ARG
1	D	139	ARG
1	D	149	ARG
1	D	153	MSE
1	D	220	GLU
1	D	258	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FLC	A	301	-	3,12,12	1.76	1 (33%)	3,17,17	1.63	1 (33%)
2	FLC	B	301	-	3,12,12	1.65	1 (33%)	3,17,17	2.47	2 (66%)
2	FLC	C	301	-	3,12,12	1.43	0	3,17,17	3.61	2 (66%)
2	FLC	D	301	-	3,12,12	1.31	0	3,17,17	1.50	0

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	FLC	A	301	-	-	0/6/16/16	0/0/0/0
2	FLC	B	301	-	-	0/6/16/16	0/0/0/0
2	FLC	C	301	-	-	0/6/16/16	0/0/0/0
2	FLC	D	301	-	-	0/6/16/16	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	FLC	CG-CB	-2.43	1.51	1.54
2	A	301	FLC	CG-CB	-2.15	1.51	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	FLC	CB-CG-CGC	-5.00	106.97	114.96
2	C	301	FLC	CB-CA-CAC	-3.75	108.96	114.96
2	B	301	FLC	CB-CG-CGC	-3.42	109.49	114.96
2	B	301	FLC	CB-CA-CAC	-2.56	110.86	114.96
2	A	301	FLC	CB-CA-CAC	-2.05	111.69	114.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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	152/162 (93%)	1.13	27 (17%) 2 2	11, 18, 29, 41	0
1	B	155/162 (95%)	1.06	28 (18%) 2 2	12, 19, 30, 39	0
1	C	158/162 (97%)	1.21	32 (20%) 1 1	12, 19, 31, 42	0
1	D	153/162 (94%)	1.16	25 (16%) 2 3	12, 19, 33, 40	0
All	All	618/648 (95%)	1.14	112 (18%) 2 2	11, 19, 33, 42	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	113	GLY	11.1
1	D	182	LYS	9.4
1	A	182	LYS	7.8
1	C	235	TYR	7.2
1	D	180	ALA	6.9
1	A	155	HIS	6.3
1	C	155	HIS	6.2
1	B	235	TYR	5.8
1	C	112	GLU	5.5
1	B	140	ALA	5.3
1	D	115	ARG	5.2
1	C	182	LYS	5.1
1	A	180	ALA	5.0
1	B	110	GLU	4.7
1	A	154	ASN	4.5
1	C	140	ALA	4.4
1	D	179	THR	4.3
1	A	181	SER	4.3
1	C	250	ASP	4.2
1	A	141	GLY	4.2
1	B	220	GLU	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	171	ARG	4.1
1	B	139	ARG	4.0
1	A	130	ALA	4.0
1	D	181	SER	4.0
1	B	155	HIS	4.0
1	A	179	THR	3.8
1	A	140	ALA	3.7
1	A	237	ASP	3.7
1	C	115	ARG	3.7
1	B	212	GLU	3.6
1	C	181	SER	3.6
1	D	117	GLU	3.6
1	B	182	LYS	3.5
1	B	154	ASN	3.5
1	B	168	THR	3.5
1	B	169	HIS	3.5
1	A	235	TYR	3.5
1	B	171	ARG	3.4
1	A	251	HIS	3.4
1	C	176	GLU	3.3
1	B	250	ASP	3.2
1	D	250	ASP	3.1
1	D	175	LEU	3.1
1	C	168	THR	3.0
1	C	259	ASN	3.0
1	D	154	ASN	3.0
1	B	112	GLU	2.9
1	D	140	ALA	2.9
1	A	142	GLU	2.9
1	B	244	ARG	2.9
1	D	174	THR	2.9
1	B	176	GLU	2.9
1	D	171	ARG	2.8
1	A	250	ASP	2.8
1	C	154	ASN	2.8
1	C	110	GLU	2.7
1	C	116	ALA	2.7
1	A	244	ARG	2.7
1	C	114	GLY	2.7
1	D	235	TYR	2.7
1	C	238	ARG	2.7
1	B	181	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	221	PRO	2.7
1	A	171	ARG	2.6
1	B	251	HIS	2.6
1	C	124	ARG	2.5
1	C	188	LEU	2.5
1	B	236	GLY	2.5
1	D	142	GLU	2.5
1	C	139	ARG	2.5
1	D	139	ARG	2.5
1	A	259	ASN	2.5
1	C	102	PHE	2.5
1	B	111	GLY	2.4
1	C	200	ALA	2.4
1	A	192	GLU	2.4
1	A	109	GLU	2.4
1	A	183	ALA	2.3
1	C	180	ALA	2.3
1	B	192	GLU	2.3
1	B	134	ARG	2.3
1	A	178	LEU	2.3
1	D	208	ALA	2.3
1	C	220	GLU	2.3
1	A	129	PRO	2.3
1	A	223	ARG	2.2
1	C	258	LEU	2.2
1	C	169	HIS	2.2
1	D	163	LEU	2.2
1	B	211	PRO	2.2
1	C	146	THR	2.2
1	D	249	THR	2.2
1	A	98	VAL	2.2
1	A	134	ARG	2.2
1	A	104	ARG	2.1
1	D	123	CYS	2.1
1	B	228	VAL	2.1
1	B	255	ARG	2.1
1	D	141	GLY	2.1
1	D	198	VAL	2.1
1	C	205[A]	ARG	2.1
1	D	251	HIS	2.1
1	A	106	ALA	2.1
1	C	212	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	155	HIS	2.1
1	C	221	PRO	2.0
1	D	129	PRO	2.0
1	B	124	ARG	2.0
1	D	99	ARG	2.0
1	B	136	LEU	2.0
1	C	192	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	FLC	B	301	13/13	0.89	0.19	0.96	12,15,22,24	0
2	FLC	D	301	13/13	0.94	0.17	-0.01	12,18,21,21	0
2	FLC	A	301	13/13	0.89	0.16	-0.19	13,15,18,19	0
2	FLC	C	301	13/13	0.93	0.15	-0.50	10,16,21,22	0
3	CL	C	2	1/1	0.97	0.07	-2.70	17,17,17,17	0
3	CL	B	3	1/1	0.97	0.04	-3.20	28,28,28,28	0
3	CL	A	4	1/1	0.97	0.04	-3.42	27,27,27,27	0
3	CL	B	1	1/1	0.97	0.09	-3.56	16,16,16,16	0
3	CL	D	5	1/1	0.71	0.16	-	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.