Full wwPDB X-ray Structure Validation Report (i)

Nov 6, 2023 – 04:58 AM EST

PDB ID	:	5VVI
Title	:	Crystal Structure of the Ligand Binding Domain of LysR-type Transcriptional Regulator, OccR from Agrobacterium tumefaciens in the Complex with Octopine
Authors	:	Kim, Y.; Chhor, G.; Jedrzejczak, R.; Winans, S.C.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2017-05-19
Resolution	:	2.28 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

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

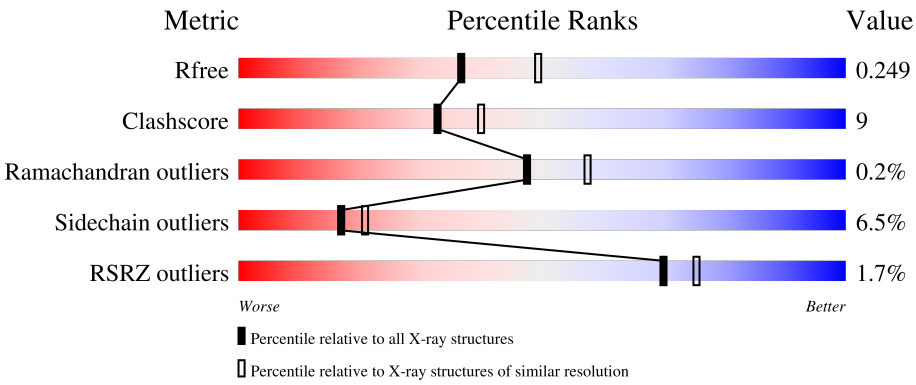
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.28 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	210	80% 16% • •
1	B	210	3% 75% 19% • •
1	C	210	% 75% 20% • •

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 4933 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.

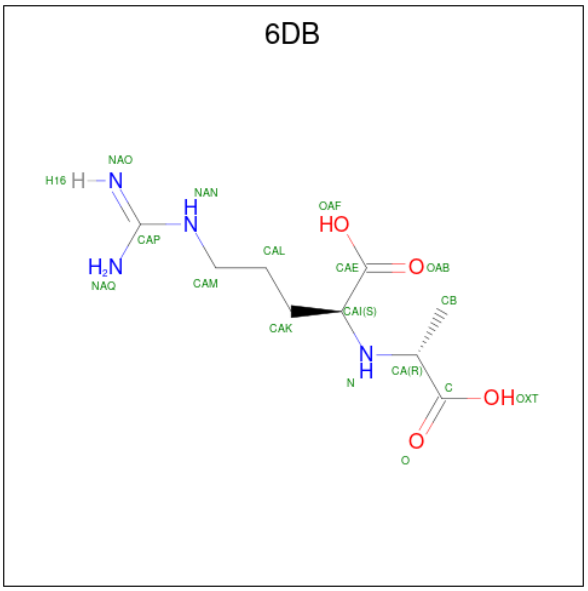
- Molecule 1 is a protein called Octopine catabolism/uptake operon regulatory protein OccR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	208	1634	1034	297	294	9	0	4	0
1	B	203	1569	995	280	286	8	0	2	0
1	C	204	1568	995	281	284	8	0	1	0

There are 12 discrepancies between the modelled and reference sequences:

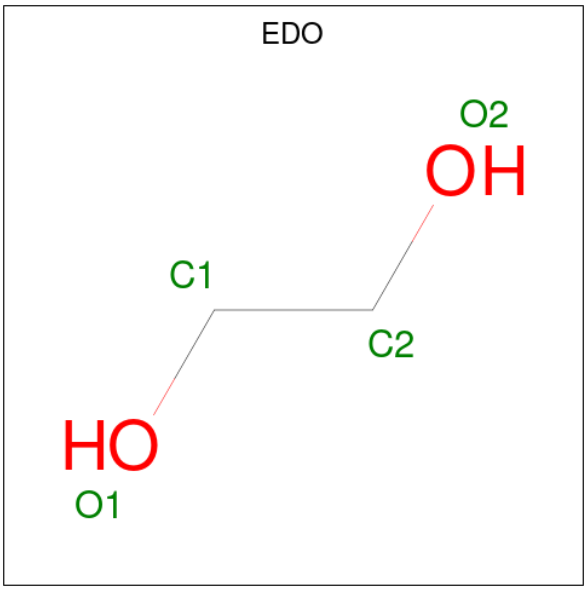
Chain	Residue	Modelled	Actual	Comment	Reference
A	89	SER	-	expression tag	UNP P0A4T4
A	90	ASN	-	expression tag	UNP P0A4T4
A	91	ALA	-	expression tag	UNP P0A4T4
A	261	GLU	GLY	conflict	UNP P0A4T4
B	89	SER	-	expression tag	UNP P0A4T4
B	90	ASN	-	expression tag	UNP P0A4T4
B	91	ALA	-	expression tag	UNP P0A4T4
B	261	GLU	GLY	conflict	UNP P0A4T4
C	89	SER	-	expression tag	UNP P0A4T4
C	90	ASN	-	expression tag	UNP P0A4T4
C	91	ALA	-	expression tag	UNP P0A4T4
C	261	GLU	GLY	conflict	UNP P0A4T4

- Molecule 2 is octopine (three-letter code: 6DB) (formula: $C_9H_{18}N_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	17	9	4	4	0	0
2	B	1	17	9	4	4	0	0
2	C	1	17	9	4	4	0	0

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	4	2	2	0	0

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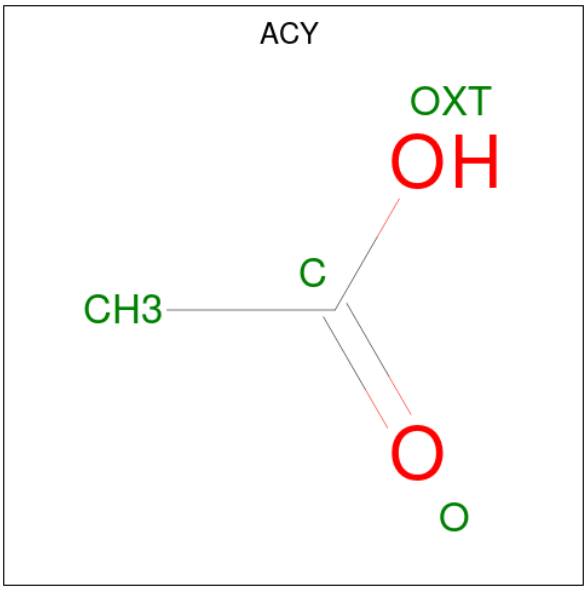
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	B	1	4	2	2	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
4	B	1	1	1	0	0

- Molecule 5 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	C	1	4	2	2	0	0

- Molecule 6 is water.

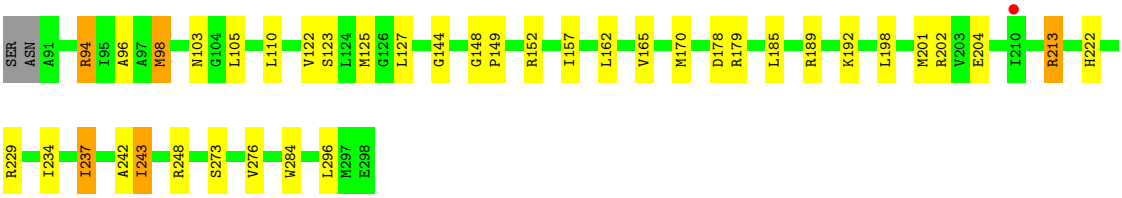
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	42	42	42	0	0
			Total	O		
6	B	36	36	36	0	0
			Total	O		
6	C	20	20	20	0	0

3 Residue-property plots (i)

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

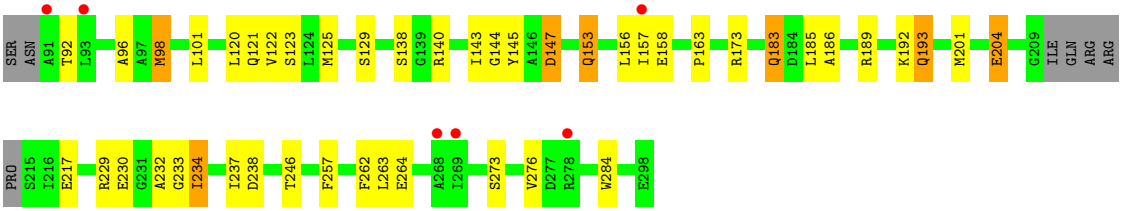
• Molecule 1: Octopine catabolism/uptake operon regulatory protein OccR

Chain A:



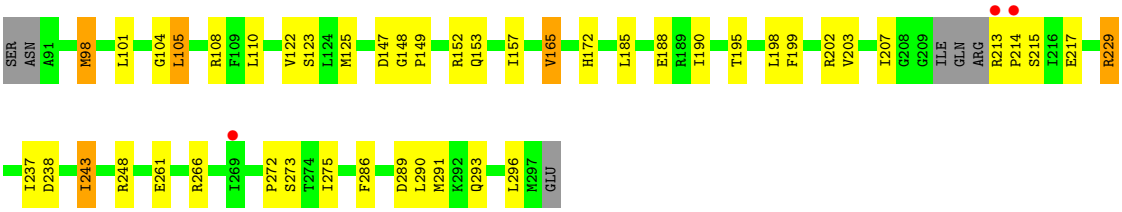
• Molecule 1: Octopine catabolism/uptake operon regulatory protein OccR

Chain B:



• Molecule 1: Octopine catabolism/uptake operon regulatory protein OccR

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	58.42Å 101.98Å 112.71Å 90.00° 98.14° 90.00°	Depositor
Resolution (Å)	31.84 – 2.28 48.13 – 2.28	Depositor EDS
% Data completeness (in resolution range)	89.0 (31.84-2.28) 89.0 (48.13-2.28)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	1.76 (at 2.27Å)	Xtriage
Refinement program	PHENIX DEV_1839	Depositor
R, R_{free}	0.192 , 0.249 0.192 , 0.249	Depositor DCC
R_{free} test set	1414 reflections (5.30%)	wwPDB-VP
Wilson B-factor (Å^2)	29.6	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 48.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4933	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

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

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 6DB, EDO, ACY, 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	0.49	0/1655	0.62	0/2226
1	B	0.48	0/1588	0.66	1/2136 (0.0%)
1	C	0.41	0/1587	0.57	0/2135
All	All	0.46	0/4830	0.62	1/6497 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	147	ASP	CB-CG-OD2	6.48	124.13	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1634	0	1667	31	0
1	B	1569	0	1594	33	0
1	C	1568	0	1602	27	0
2	A	17	0	0	0	0
2	B	17	0	0	0	0
2	C	17	0	0	1	0
3	A	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	6	1	0
4	B	1	0	0	0	0
5	C	4	0	3	0	0
6	A	42	0	0	3	0
6	B	36	0	0	5	0
6	C	20	0	0	0	0
All	All	4933	0	4878	85	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 9.

All (85) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:GLY:HA2	3:B:302:EDO:H11	1.64	0.78
1:A:229:ARG:HH11	1:A:248:ARG:NH1	1.89	0.70
1:B:156:LEU:HD12	1:B:157:ILE:H	1.59	0.67
1:A:103:ASN:HD22	1:A:222[B]:HIS:CD2	2.12	0.66
1:A:152:ARG:HG2	1:A:157:ILE:HD12	1.80	0.64
1:C:152:ARG:HD3	1:C:157:ILE:HD12	1.78	0.64
1:C:266[A]:ARG:HH22	1:C:272:PRO:HA	1.64	0.62
1:A:123:SER:HB2	1:A:125:MSE:HE3	1.83	0.59
1:A:94[A]:ARG:HG3	6:A:402:HOH:O	2.03	0.58
1:A:94[A]:ARG:NH2	6:A:402:HOH:O	2.37	0.57
1:C:98:MSE:SE	1:C:101:LEU:HG	2.55	0.57
1:C:290:LEU:HD22	1:C:291:MSE:HE2	1.88	0.56
1:A:222[B]:HIS:HD2	6:A:410:HOH:O	1.90	0.55
1:B:192:LYS:HG2	1:B:193:GLN:O	2.07	0.54
1:A:201:MSE:SE	1:B:204:GLU:HG2	2.58	0.54
1:A:149:PRO:HD3	1:A:198:LEU:HD23	1.91	0.53
1:B:125:MSE:HE1	1:C:217:GLU:HG2	1.90	0.53
1:B:186:ALA:O	6:B:401:HOH:O	2.19	0.53
1:A:110:LEU:HD11	1:A:122:VAL:HG11	1.92	0.52
1:B:129:SER:OG	1:B:147:ASP:N	2.35	0.51
1:A:229:ARG:HH11	1:A:248:ARG:HH12	1.56	0.51
1:A:237:ILE:HD13	1:A:242:ALA:HB2	1.94	0.50
1:C:213:ARG:HD3	1:C:214:PRO:HD2	1.94	0.50
1:B:183:GLN:HG3	6:B:432:HOH:O	2.12	0.50
1:C:105:LEU:HD23	1:C:286:PHE:HE2	1.76	0.50
1:C:273:SER:OG	1:C:275:ILE:HG22	2.12	0.49
1:A:94[A]:ARG:NH2	1:A:123:SER:OG	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:GLY:HA3	1:C:261:GLU:HG2	1.93	0.49
1:B:138:SER:HB3	1:B:140:ARG:HG3	1.95	0.49
1:B:173:ARG:NH1	6:B:405:HOH:O	2.31	0.49
1:A:189:ARG:HB2	1:A:213:ARG:HH11	1.77	0.48
1:C:238:ASP:N	1:C:238:ASP:OD1	2.42	0.48
1:C:203:VAL:O	1:C:207:ILE:HG12	2.14	0.48
1:A:165:VAL:HG11	1:A:296:LEU:HD22	1.95	0.48
1:B:238:ASP:HB2	6:B:417:HOH:O	2.15	0.47
1:C:198:LEU:HD21	1:C:202:ARG:HE	1.79	0.47
1:C:199:PHE:O	1:C:203:VAL:HG23	2.14	0.47
1:B:157:ILE:HD13	1:B:263:LEU:HD13	1.96	0.47
1:B:217:GLU:HG2	1:C:125:MSE:HE1	1.97	0.47
1:B:121:GLN:NE2	1:C:215:SER:O	2.47	0.47
1:C:172:HIS:NE2	1:C:188:GLU:OE2	2.48	0.46
1:A:185:LEU:HD22	1:A:234:ILE:HD11	1.97	0.46
1:A:148:GLY:HA2	1:A:149:PRO:HA	1.64	0.46
1:B:145:TYR:HA	1:B:263:LEU:O	2.15	0.46
1:C:243:ILE:HA	1:C:243:ILE:HD12	1.59	0.46
1:C:266[A]:ARG:HE	1:C:266[A]:ARG:HB3	1.52	0.46
1:A:170:MSE:HE2	1:A:248:ARG:O	2.15	0.46
1:A:96:ALA:O	1:A:144:GLY:HA2	2.16	0.45
1:B:229:ARG:NH2	1:B:230:GLU:OE2	2.50	0.45
1:B:98:MSE:HG3	1:B:147:ASP:OD1	2.16	0.45
1:A:98:MSE:HA	1:A:127:LEU:O	2.16	0.45
1:B:96:ALA:O	1:B:144:GLY:HA2	2.16	0.45
1:B:273:SER:HB3	1:B:276:VAL:HB	1.99	0.45
1:C:110:LEU:HD11	1:C:122:VAL:HG11	1.99	0.45
1:B:153:GLN:HE21	1:B:153:GLN:HB3	1.57	0.45
1:A:198:LEU:O	1:A:202:ARG:HG3	2.17	0.44
1:B:98:MSE:HE1	1:B:262:PHE:HB2	2.00	0.44
1:B:230:GLU:HG3	1:C:108:ARG:HA	1.99	0.44
1:A:229:ARG:NH1	1:A:248:ARG:HH12	2.15	0.44
1:A:201:MSE:HE1	1:B:204:GLU:HG2	1.98	0.44
1:B:138:SER:OG	1:B:140:ARG:NH2	2.51	0.44
1:B:189:ARG:HD2	1:B:232:ALA:O	2.18	0.43
1:C:190:ILE:O	1:C:215:SER:OG	2.30	0.43
1:C:98:MSE:HG2	1:C:147:ASP:OD2	2.19	0.43
1:A:94[B]:ARG:HE	1:A:94[B]:ARG:HB2	1.60	0.42
1:C:148:GLY:HA2	1:C:149:PRO:HA	1.79	0.42
1:A:105:LEU:HD23	1:A:105:LEU:HA	1.86	0.42
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:TRP:HA	1:A:284:TRP:CE3	2.54	0.42
1:C:98:MSE:HB3	2:C:301:6DB:OXT	2.20	0.42
1:B:185:LEU:HD22	1:B:234:ILE:HD11	2.02	0.41
1:C:229:ARG:HD2	1:C:248:ARG:CZ	2.51	0.41
1:B:193:GLN:HG2	6:B:418:HOH:O	2.21	0.41
1:A:192:LYS:NZ	1:A:204:GLU:OE1	2.52	0.41
1:C:165:VAL:HG11	1:C:296:LEU:HD22	2.03	0.41
1:A:273:SER:HB3	1:A:276:VAL:HB	2.02	0.41
1:B:120:LEU:O	1:B:122:VAL:HG23	2.21	0.41
1:B:163:PRO:HB3	1:B:257:PHE:CD1	2.56	0.41
1:B:237:ILE:HD12	1:B:237:ILE:HA	1.81	0.41
1:A:213:ARG:HB3	1:A:213:ARG:CZ	2.51	0.40
1:A:243:ILE:HD12	1:A:243:ILE:HA	1.54	0.40
1:B:156:LEU:HD12	1:B:157:ILE:N	2.29	0.40
1:B:183:GLN:HG3	1:B:183:GLN:H	1.69	0.40
1:B:284:TRP:CE3	1:B:284:TRP:HA	2.57	0.40
1:C:237:ILE:HD12	1:C:237:ILE:HA	1.92	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	210/210 (100%)	204 (97%)	6 (3%)	0	100	100
1	B	201/210 (96%)	197 (98%)	4 (2%)	0	100	100
1	C	201/210 (96%)	194 (96%)	6 (3%)	1 (0%)	29	34
All	All	612/630 (97%)	595 (97%)	16 (3%)	1 (0%)	47	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	104	GLY

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	171/161 (106%)	162 (95%)	9 (5%)	22	29
1	B	164/161 (102%)	150 (92%)	14 (8%)	10	12
1	C	164/161 (102%)	153 (93%)	11 (7%)	16	19
All	All	499/483 (103%)	465 (93%)	34 (7%)	17	19

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94[A]	ARG
1	A	94[B]	ARG
1	A	98	MSE
1	A	178	ASP
1	A	179[A]	ARG
1	A	179[B]	ARG
1	A	213	ARG
1	A	237	ILE
1	A	243	ILE
1	B	92	THR
1	B	98	MSE
1	B	101	LEU
1	B	123	SER
1	B	143	ILE
1	B	153	GLN
1	B	158	GLU
1	B	183	GLN
1	B	193	GLN
1	B	201	MSE
1	B	204	GLU
1	B	234	ILE
1	B	246	THR

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Mol	Chain	Res	Type
1	B	264	GLU
1	C	98	MSE
1	C	105	LEU
1	C	123	SER
1	C	153	GLN
1	C	165	VAL
1	C	185	LEU
1	C	195	THR
1	C	229	ARG
1	C	243	ILE
1	C	289	ASP
1	C	293	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	153	GLN
1	C	293	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\#\|Z\| > 2$
5	ACY	C	302	-	3,3,3	0.69	0	3,3,3	1.13	0
3	EDO	A	302	-	3,3,3	0.44	0	2,2,2	0.39	0
2	6DB	B	301	1	16,16,16	1.10	1 (6%)	17,20,20	2.13	6 (35%)
2	6DB	A	301	-	16,16,16	1.00	1 (6%)	17,20,20	1.90	5 (29%)
2	6DB	C	301	-	16,16,16	1.01	1 (6%)	17,20,20	1.93	4 (23%)
3	EDO	B	302	-	3,3,3	0.40	0	2,2,2	0.42	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
3	EDO	A	302	-	-	0/1/1/1	-
2	6DB	B	301	1	-	5/19/19/19	-
2	6DB	A	301	-	-	6/19/19/19	-
2	6DB	C	301	-	-	5/19/19/19	-
3	EDO	B	302	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	6DB	CAP-NAN	2.53	1.38	1.33
2	A	301	6DB	CAP-NAN	2.52	1.38	1.33
2	B	301	6DB	CAP-NAN	2.36	1.38	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	301	6DB	CAK-CAI-CAE	-5.43	97.26	110.35
2	A	301	6DB	CB-CA-N	4.93	119.28	108.85
2	B	301	6DB	NAN-CAP-NAO	-4.39	112.98	120.70
2	B	301	6DB	CAK-CAI-CAE	-3.83	101.12	110.35
2	B	301	6DB	CB-CA-N	2.91	115.00	108.85
2	C	301	6DB	CB-CA-C	-2.68	104.15	110.20
2	B	301	6DB	OXT-C-O	2.63	130.07	124.09
2	C	301	6DB	OXT-C-O	2.60	130.00	124.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	301	6DB	NAQ-CAP-NAN	-2.42	113.62	119.19
2	B	301	6DB	O-C-CA	-2.35	114.19	121.83
2	A	301	6DB	CAK-CAI-CAE	-2.28	104.85	110.35
2	A	301	6DB	OAB-CAE-CAI	-2.26	114.82	122.26
2	A	301	6DB	CAL-CAK-CAI	-2.15	107.30	113.92
2	C	301	6DB	OAB-CAE-CAI	-2.15	115.21	122.26
2	B	301	6DB	CB-CA-C	-2.15	105.36	110.20

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	6DB	O-C-CA-N
2	A	301	6DB	C-CA-N-CAI
2	B	301	6DB	O-C-CA-N
2	B	301	6DB	OXT-C-CA-CB
2	B	301	6DB	OXT-C-CA-N
2	B	301	6DB	C-CA-N-CAI
2	C	301	6DB	C-CA-N-CAI
2	B	301	6DB	O-C-CA-CB
2	C	301	6DB	O-C-CA-CB
3	B	302	EDO	O1-C1-C2-O2
2	A	301	6DB	OXT-C-CA-N
2	A	301	6DB	CB-CA-N-CAI
2	A	301	6DB	O-C-CA-CB
2	A	301	6DB	OXT-C-CA-CB
2	C	301	6DB	OXT-C-CA-CB
2	C	301	6DB	O-C-CA-N
2	C	301	6DB	OXT-C-CA-N

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	301	6DB	1	0
3	B	302	EDO	1	0

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	200/210 (95%)	-0.12	1 (0%)	91	93	11, 30, 68, 91	0
1	B	195/210 (92%)	0.15	6 (3%)	49	54	14, 39, 73, 90	0
1	C	196/210 (93%)	0.15	3 (1%)	73	78	20, 48, 84, 99	0
All	All	591/630 (93%)	0.06	10 (1%)	70	75	11, 39, 76, 99	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	268	ALA	3.2
1	C	214	PRO	2.9
1	B	91	ALA	2.6
1	B	278	ARG	2.6
1	B	157	ILE	2.4
1	C	269	ILE	2.4
1	C	213	ARG	2.4
1	A	210	ILE	2.3
1	B	93	LEU	2.3
1	B	269	ILE	2.2

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
5	ACY	C	302	4/4	0.88	0.33	47,47,47,47	0
4	CL	B	303	1/1	0.89	0.16	73,73,73,73	0
2	6DB	B	301	17/17	0.91	0.16	20,24,27,27	0
3	EDO	A	302	4/4	0.91	0.22	28,28,28,28	4
2	6DB	C	301	17/17	0.94	0.15	28,31,32,32	0
2	6DB	A	301	17/17	0.96	0.12	16,18,19,20	0
3	EDO	B	302	4/4	0.96	0.15	39,39,39,39	0

6.5 Other polymers (i)

There are no such residues in this entry.