



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 02:12 am BST

PDB ID : 5FS0
Title : crystal structure of the phenol-responsive sensory domain of the transcription activator PoxR with 2,4-dichlorophenol
Authors : Patil, V.V.; Woo, E.J.
Deposited on : 2015-12-23
Resolution : 2.40 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

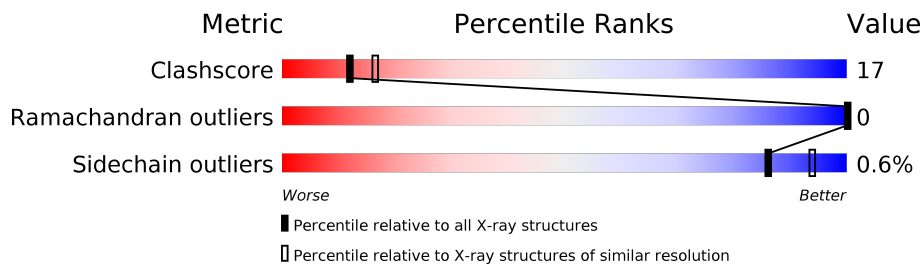
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	211	
1	B	211	

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
2	5JC	A	1209	-	-	X	-

2 Entry composition [i](#)

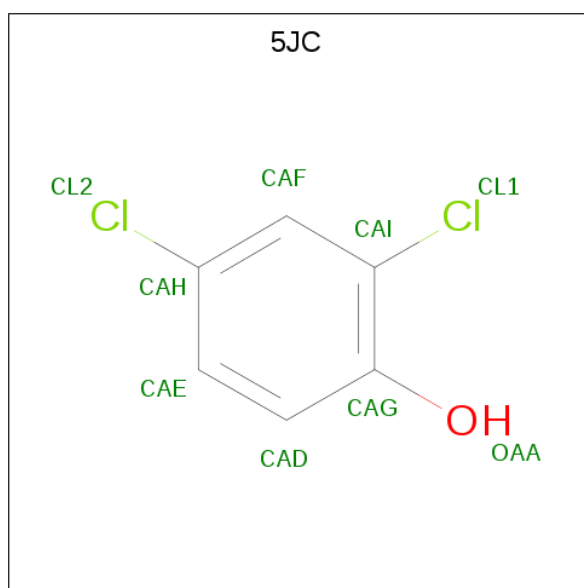
There are 4 unique types of molecules in this entry. The entry contains 3158 atoms, of which 8 are hydrogens and 0 are deuteriums.

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 POSITIVE PHENOL-DEGRADATIVE GENE REGULATOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	196	Total 1531	C 966	N 274	O 275	S 16	0	0	0
1	B	196	Total 1531	C 966	N 274	O 275	S 16	0	0	0

- Molecule 2 is 2,4-dichlorophenol (three-letter code: 5JC) (formula: C₆H₄Cl₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	H	O		
2	A	1	Total 13	C 6	Cl 2	H 4	O 1	0	0
2	B	1	Total 13	C 6	Cl 2	H 4	O 1	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Zn 1	0	0
3	A	1	Total 1	Zn 1	0	0

- Molecule 4 is water.

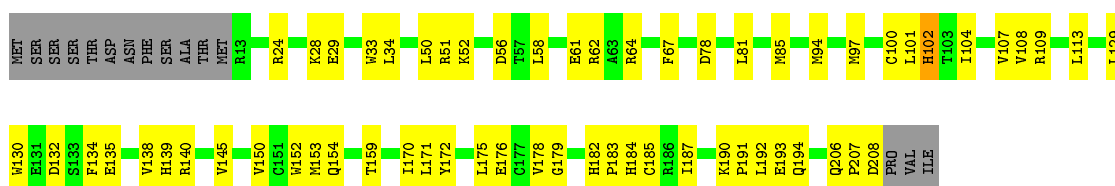
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	43	Total 43	O 43	0	0
4	B	25	Total 25	O 25	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

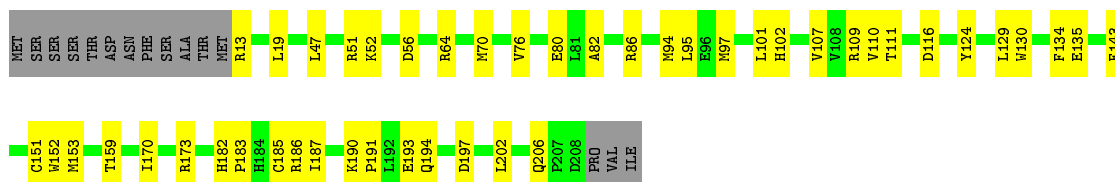
- Molecule 1: POSITIVE PHENOL-DEGRADATIVE GENE REGULATOR

Chain A: 



- Molecule 1: POSITIVE PHENOL-DEGRADATIVE GENE REGULATOR

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.69Å 70.87Å 104.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.35 – 2.40 31.35 – 1.79	Depositor EDS
% Data completeness (in resolution range)	99.1 (31.35-2.40) 80.1 (31.35-1.79)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.88 (at 1.79Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE: 1.9_1692)	Depositor
R, R_{free}	0.181 , 0.258 0.213 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtrriage
Anisotropy	0.319	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3158	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 5JC

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	0.47	0/1564	0.58	1/2107 (0.0%)
1	B	0.44	0/1564	0.51	0/2107
All	All	0.46	0/3128	0.55	1/4214 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	GLN	C-N-CD	5.71	140.38	128.40

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	1531	0	1505	62	0
1	B	1531	0	1505	46	0
2	A	9	4	0	4	0
2	B	9	4	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	43	0	0	3	0
4	B	25	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3150	8	3010	105	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 17.

All (105) 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:107:VAL:HG13	1:B:134:PHE:CE1	1.74	1.21
1:A:107:VAL:HG13	1:A:134:PHE:HE1	1.13	1.09
1:A:107:VAL:HG13	1:A:134:PHE:CE1	1.87	1.07
1:A:140:ARG:HH21	1:A:145:VAL:HG22	1.38	0.89
1:B:107:VAL:HG13	1:B:134:PHE:HE1	1.16	0.88
1:B:19:LEU:HD11	1:B:97:MET:HG2	1.55	0.88
1:A:78:ASP:HB3	1:A:101:LEU:HD13	1.57	0.87
1:A:130:TRP:HB2	1:A:185:CYS:HB2	1.56	0.86
1:B:135:GLU:OE2	1:B:153:MET:N	2.09	0.85
1:A:51:ARG:NH2	1:A:135:GLU:OE2	2.15	0.79
1:A:109:ARG:NH2	1:A:132:ASP:OD2	2.15	0.79
1:A:207:PRO:O	1:A:208:ASP:HB3	1.84	0.78
1:B:82:ALA:O	1:B:94:MET:HE1	1.84	0.78
1:B:107:VAL:CG1	1:B:134:PHE:HE1	1.97	0.77
1:B:130:TRP:HB2	1:B:185:CYS:HB2	1.65	0.76
1:A:52:LYS:NZ	1:A:56:ASP:OD2	2.18	0.75
1:A:134:PHE:O	1:A:138:VAL:HG23	1.88	0.74
1:B:143:PHE:O	4:B:2018:HOH:O	2.06	0.74
1:A:81:LEU:HD11	1:A:85:MET:HE1	1.68	0.74
1:B:173:ARG:NH1	1:B:197:ASP:OD2	2.21	0.74
1:A:182:HIS:HB3	1:A:183:PRO:HD2	1.69	0.73
1:B:107:VAL:HG13	1:B:134:PHE:CD1	2.22	0.73
1:A:139:HIS:CD2	1:A:150:VAL:HG11	2.25	0.71
1:B:86:ARG:O	1:B:86:ARG:HG2	1.90	0.70
1:A:107:VAL:CG1	1:A:134:PHE:HE1	2.00	0.69
1:B:202:LEU:O	4:B:2025:HOH:O	2.13	0.66
1:B:64:ARG:HG3	1:B:152:TRP:CD1	2.30	0.65
1:B:82:ALA:O	1:B:94:MET:CE	2.44	0.65
1:B:194:GLN:OE1	4:B:2015:HOH:O	2.15	0.65
1:B:130:TRP:CE3	1:B:187:ILE:HD11	2.33	0.64
1:A:78:ASP:HB3	1:A:101:LEU:CD1	2.27	0.63
1:A:51:ARG:HH22	1:A:135:GLU:CG	2.11	0.63
1:A:78:ASP:CB	1:A:101:LEU:HD13	2.28	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:LYS:O	1:A:29:GLU:HG3	2.01	0.61
1:A:34:LEU:HD21	1:A:101:LEU:HD22	1.85	0.58
1:B:107:VAL:O	1:B:134:PHE:HD1	1.86	0.58
1:B:13:ARG:HG3	4:B:2001:HOH:O	2.03	0.58
1:A:134:PHE:CG	1:A:135:GLU:N	2.71	0.57
1:A:178:VAL:HG13	1:A:184:HIS:HA	1.85	0.57
1:A:51:ARG:NH2	1:A:135:GLU:HG3	2.20	0.57
1:B:76:VAL:O	1:B:80:GLU:HG3	2.04	0.57
1:A:102:HIS:ND1	2:A:1209:5JC:OAA	2.34	0.56
1:B:187:ILE:HD12	1:B:187:ILE:N	2.20	0.55
1:B:19:LEU:HD11	1:B:97:MET:CG	2.35	0.55
1:B:51:ARG:NH2	1:B:135:GLU:HG2	2.23	0.54
1:A:130:TRP:CB	1:A:185:CYS:HB2	2.34	0.54
1:A:187:ILE:HD11	2:A:1209:5JC:CL1	2.44	0.54
1:B:111:THR:HB	1:B:129:LEU:HB3	1.90	0.54
1:B:64:ARG:HG3	1:B:152:TRP:CG	2.43	0.54
1:A:113:LEU:HD11	1:A:129:LEU:CD2	2.39	0.52
1:A:94:MET:HA	1:A:97:MET:HE2	1.91	0.52
1:A:107:VAL:O	1:A:134:PHE:HD1	1.93	0.52
1:A:113:LEU:HD21	1:A:129:LEU:HG	1.92	0.51
1:A:175:LEU:O	1:A:176:GLU:HG3	2.11	0.51
1:A:51:ARG:NH2	1:A:135:GLU:CG	2.72	0.51
1:B:52:LYS:NZ	1:B:56:ASP:OD2	2.43	0.51
1:A:207:PRO:HD3	4:A:2014:HOH:O	2.10	0.51
1:A:51:ARG:HH22	1:A:135:GLU:CD	2.13	0.51
1:A:29:GLU:O	1:B:109:ARG:HB2	2.11	0.50
1:A:130:TRP:NE1	2:A:1209:5JC:CL1	2.82	0.49
1:A:190:LYS:HB2	1:A:191:PRO:HD2	1.94	0.49
1:B:135:GLU:HB3	1:B:151:CYS:HA	1.93	0.49
1:B:124:TYR:O	1:B:191:PRO:HD3	2.12	0.49
1:A:135:GLU:OE1	1:A:154:GLN:HG3	2.13	0.49
1:A:190:LYS:HE3	1:A:194:GLN:O	2.13	0.49
1:A:50:LEU:HD22	1:B:70:MET:CE	2.43	0.48
1:B:206:GLN:HG2	4:B:2025:HOH:O	2.14	0.48
1:A:64:ARG:HA	1:A:152:TRP:CE3	2.49	0.47
1:A:176:GLU:O	1:A:185:CYS:HA	2.15	0.47
1:A:178:VAL:CG1	1:A:184:HIS:HA	2.45	0.47
1:B:153:MET:HB2	1:B:153:MET:HE2	1.68	0.46
1:B:186:ARG:C	1:B:187:ILE:HD12	2.36	0.46
1:B:110:VAL:HG22	1:B:130:TRP:CD1	2.51	0.46
1:B:95:LEU:HD13	1:B:95:LEU:C	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:ARG:HG3	1:A:152:TRP:CG	2.51	0.46
1:B:47:LEU:HA	1:B:47:LEU:HD23	1.71	0.45
1:A:172:TYR:CD2	1:A:187:ILE:HD12	2.52	0.45
1:B:116:ASP:O	1:B:124:TYR:HA	2.17	0.45
1:B:64:ARG:HA	1:B:152:TRP:CE3	2.52	0.44
1:A:29:GLU:HA	1:B:109:ARG:HE	1.81	0.43
1:A:108:VAL:HG22	1:A:109:ARG:N	2.33	0.43
1:A:187:ILE:CD1	2:A:1209:5JC:CL1	3.04	0.43
1:A:193:GLU:HG2	1:A:193:GLU:O	2.19	0.43
1:A:58:LEU:HD22	1:A:62:ARG:NH1	2.34	0.43
1:B:193:GLU:OE1	1:B:193:GLU:N	2.48	0.43
1:A:67:PHE:HD1	1:A:153:MET:HE2	1.84	0.43
1:A:61:GLU:OE1	4:A:2012:HOH:O	2.21	0.42
1:A:175:LEU:HD12	4:A:2038:HOH:O	2.19	0.42
1:A:178:VAL:HA	1:A:182:HIS:O	2.19	0.42
1:B:135:GLU:OE2	1:B:152:TRP:N	2.53	0.42
1:B:159:THR:HG23	1:B:170:ILE:O	2.20	0.42
1:A:150:VAL:HG22	1:A:179:GLY:HA3	2.02	0.42
1:B:130:TRP:CD2	1:B:187:ILE:HD11	2.54	0.41
1:A:67:PHE:CD1	1:A:153:MET:HE2	2.55	0.41
1:A:28:LYS:C	1:A:29:GLU:HG3	2.41	0.41
1:A:24:ARG:HG2	1:A:33:TRP:HB2	2.02	0.41
1:A:171:LEU:HD11	1:A:192:LEU:HD13	2.01	0.41
1:B:134:PHE:CG	1:B:135:GLU:N	2.88	0.41
1:B:182:HIS:HB3	1:B:183:PRO:HD2	2.03	0.41
1:A:178:VAL:HG13	1:A:184:HIS:CA	2.50	0.41
1:A:100:CYS:O	1:A:104:ILE:HG13	2.20	0.40
1:A:94:MET:HA	1:A:97:MET:CE	2.50	0.40
1:A:159:THR:HG23	1:A:170:ILE:O	2.21	0.40
1:B:19:LEU:HD22	1:B:101:LEU:HD21	2.02	0.40
1:B:190:LYS:HB2	1:B:191:PRO:CD	2.51	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	194/211 (92%)	182 (94%)	12 (6%)	0	100	100
1	B	194/211 (92%)	188 (97%)	6 (3%)	0	100	100
All	All	388/422 (92%)	370 (95%)	18 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	157/171 (92%)	156 (99%)	1 (1%)	86	94
1	B	157/171 (92%)	156 (99%)	1 (1%)	86	94
All	All	314/342 (92%)	312 (99%)	2 (1%)	86	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	HIS
1	B	102	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	154	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	5JC	B	1210	-	9,9,9	4.55	6 (66%)	12,12,12	0.88	0
2	5JC	A	1209	-	9,9,9	4.52	6 (66%)	12,12,12	0.67	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	5JC	B	1210	-	-	-	0/1/1/1
2	5JC	A	1209	-	-	-	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1210	5JC	CAF-CAI	7.50	1.50	1.38
2	A	1209	5JC	CAE-CAH	7.49	1.52	1.38
2	B	1210	5JC	CAE-CAH	7.26	1.51	1.38
2	A	1209	5JC	CAF-CAI	7.07	1.50	1.38
2	B	1210	5JC	CAD-CAG	7.00	1.51	1.39
2	A	1209	5JC	CAD-CAG	6.52	1.51	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1209	5JC	CAG-CAI	-4.62	1.35	1.39
2	B	1210	5JC	CAG-CAI	-3.64	1.36	1.39
2	B	1210	5JC	OAA-CAG	2.48	1.41	1.36
2	A	1209	5JC	CAF-CAH	-2.39	1.34	1.38
2	B	1210	5JC	CAF-CAH	-2.25	1.34	1.38
2	A	1209	5JC	OAA-CAG	2.18	1.40	1.36

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1209	5JC	4	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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.