



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 16, 2017 – 10:10 PM EDT

PDB ID : 5V0L
Title : Crystal structure of the AHR-ARNT heterodimer in complex with the DRE
Authors : Seok, S.-H.; Lee, W.; Jiang, L.; Bradfield, C.A.; Xing, Y.
Deposited on : unknown
Resolution : 4.00 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20029824

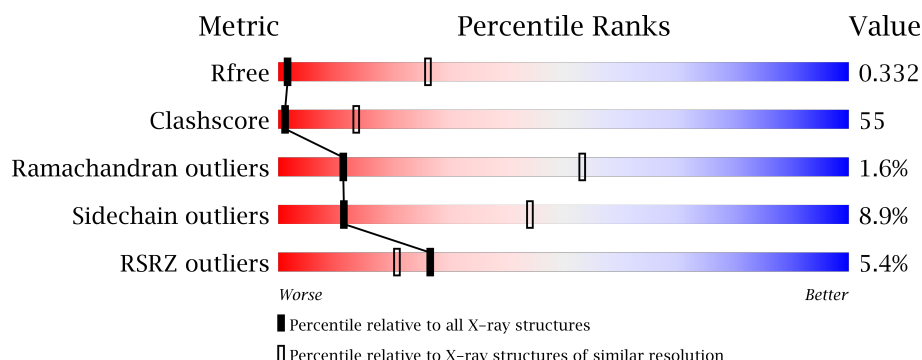
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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}	100719	1088 (4.40-3.60)
Clashscore	112137	1187 (4.40-3.60)
Ramachandran outliers	110173	1139 (4.40-3.60)
Sidechain outliers	110143	1126 (4.40-3.60)
RSRZ outliers	101464	1099 (4.40-3.60)

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	279	
2	B	241	
3	C	17	
4	D	14	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3418 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 Aryl hydrocarbon receptor nuclear translocator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	S	0	0	0
			1391	865	252	263	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	68	GLY	-	expression tag	UNP P27540
A	69	SER	-	expression tag	UNP P27540

- Molecule 2 is a protein called Aryl hydrocarbon receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	166	Total	C	N	O	S	0	0	0
			1347	866	232	245	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	27	GLY	-	expression tag	UNP P30561
B	28	SER	-	expression tag	UNP P30561

- Molecule 3 is a DNA chain called DNA (5'-D(P*GP*GP*AP*TP*TP*GP*CP*GP*TP*GP*AP*GP*AP*AP*CP*TP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	17	Total	C	N	O	P	0	0	0
			356	168	69	102	17			

- Molecule 4 is a DNA chain called DNA (5'-D(P*AP*GP*TP*TP*CP*TP*CP*AP*CP*GP*CP*AP*AP*T)-3').

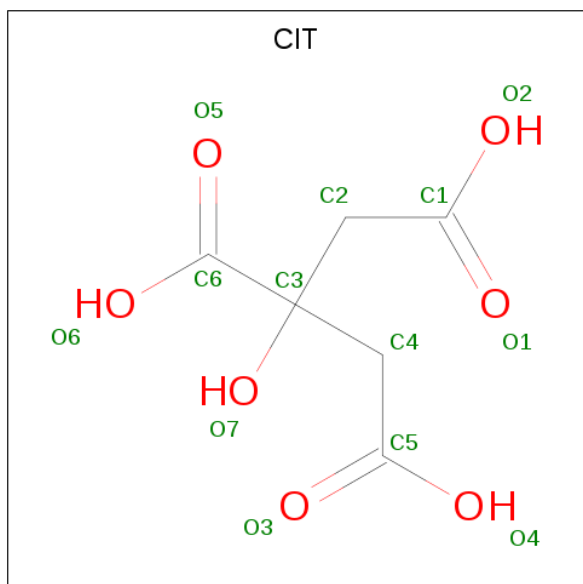
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	14	Total	C	N	O	P	0	0	0
			284	136	50	84	14			

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



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

- Molecule 3: DNA (5'-D(P*GP*GP*AP*TP*TP*GP*CP*GP*TP*GP*AP*GP*AP*AP*CP*T
P*G)-3')

Chain C:  12% 100%

G4	G5	A6	T7	T8	G9	C10	G11	T12	G13	A14	G15	A16	A17	C18	T19	G20
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- Molecule 4: DNA (5'-D(P*AP*GP*TP*TP*CP*TP*CP*AP*CP*GP*CP*AP*AP*T)-3'
)

Chain D:  14% 7% 93%

A4	G5	T6	T7	C8	T9	C10	A11	C12	G13	C14	A15	A16	T17
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	78.20 Å 64.36 Å 157.72 Å 90.00° 100.14° 90.00°	Depositor
Resolution (Å)	39.69 – 4.00 39.68 – 4.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.69-4.00) 99.5 (39.68-4.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 4.00 Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, R_{free}	0.285 , 0.323 0.306 , 0.332	Depositor DCC
R_{free} test set	332 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	194.2	Xtriage
Anisotropy	0.490	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 235.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3418	wwPDB-VP
Average B, all atoms (Å ²)	221.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% 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: PEG, CIT

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	0.32	0/1412	0.61	5/1896 (0.3%)
2	B	0.27	0/1372	0.51	2/1852 (0.1%)
3	C	0.45	0/400	0.76	0/617
4	D	0.44	0/317	0.87	0/486
All	All	0.33	0/3501	0.63	7/4851 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	191	ASP	O-C-N	-8.66	108.85	122.70
2	B	81	PHE	CB-CA-C	7.69	125.78	110.40
1	A	191	ASP	CA-C-N	6.23	130.90	117.20
1	A	125	LYS	C-N-CD	5.52	139.98	128.40
1	A	191	ASP	C-N-CA	5.51	135.49	121.70
2	B	82	PHE	N-CA-C	5.21	125.06	111.00
1	A	302	HIS	N-CA-CB	-5.16	101.31	110.60

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	1391	0	1370	209	2
2	B	1347	0	1337	148	1
3	C	356	0	192	19	0
4	D	284	0	159	17	1
5	A	7	0	10	0	0
5	B	7	0	10	0	0
6	B	13	0	5	0	0
6	C	13	0	5	0	0
All	All	3418	0	3088	359	2

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 55.

All (359) 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:180:SER:CB	1:A:183:THR:CG2	2.05	1.34
1:A:86:LYS:HA	1:A:90:ALA:CB	1.58	1.34
1:A:180:SER:HB3	1:A:183:THR:CG2	1.62	1.30
1:A:86:LYS:O	1:A:90:ALA:HB3	1.32	1.26
1:A:331:GLN:NE2	1:A:332:GLY:O	1.68	1.25
1:A:180:SER:CB	1:A:183:THR:HG23	1.70	1.18
1:A:86:LYS:C	1:A:90:ALA:HB3	1.69	1.13
1:A:180:SER:OG	1:A:183:THR:CG2	2.00	1.09
1:A:331:GLN:CD	1:A:332:GLY:O	1.92	1.08
1:A:86:LYS:HA	1:A:90:ALA:HB2	1.08	1.04
1:A:86:LYS:CA	1:A:90:ALA:CB	2.36	1.02
1:A:169:LEU:HD22	1:A:192:SER:HB3	1.45	0.98
1:A:86:LYS:CA	1:A:90:ALA:HB3	1.94	0.98
1:A:180:SER:OG	1:A:183:THR:HG22	1.60	0.98
1:A:164:LEU:HD22	1:A:167:LEU:HD23	1.46	0.97
1:A:119:CYS:O	1:A:122:LEU:HD23	1.66	0.96
1:A:331:GLN:OE1	1:A:332:GLY:N	1.97	0.96
2:B:80:SER:OG	2:B:84:VAL:CG1	2.16	0.94
1:A:169:LEU:HD22	1:A:192:SER:CB	1.99	0.92
2:B:80:SER:OG	2:B:84:VAL:HG11	1.69	0.92
1:A:213:VAL:HA	1:A:267:MET:HA	1.52	0.92
2:B:145:GLY:HA3	2:B:216:CYS:HB3	1.49	0.91
1:A:121:ALA:C	1:A:122:LEU:HD22	1.89	0.90
1:A:190:SER:O	1:A:193:VAL:HG23	1.71	0.89
2:B:124:VAL:HA	2:B:262:ILE:HG22	1.54	0.89
1:A:136:VAL:HG22	2:B:49:LEU:HD21	1.55	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:CYS:O	1:A:122:LEU:CD2	2.22	0.88
1:A:180:SER:CB	1:A:183:THR:HG21	2.00	0.88
1:A:180:SER:HB3	1:A:183:THR:HG23	0.89	0.87
1:A:331:GLN:NE2	1:A:334:LYS:HB2	1.89	0.86
1:A:108:TYR:HA	1:A:111:GLU:HB3	1.55	0.86
3:C:14:DA:H1'	3:C:15:DG:H5'	1.59	0.84
1:A:139:MET:HG2	2:B:49:LEU:HD22	1.60	0.83
2:B:176:LEU:HD13	2:B:237:LEU:HD13	1.59	0.83
1:A:213:VAL:HG21	1:A:221:LEU:HD22	1.59	0.83
2:B:124:VAL:O	2:B:137:SER:OG	2.00	0.80
2:B:114:GLU:HB3	2:B:117:LEU:HD21	1.61	0.80
1:A:184:GLY:O	1:A:208:THR:HB	1.81	0.79
1:A:108:TYR:O	1:A:112:LEU:N	2.16	0.79
1:A:107:ALA:O	1:A:111:GLU:N	2.14	0.78
1:A:128:LYS:HA	1:A:131:ILE:HD12	1.65	0.77
2:B:137:SER:HB2	2:B:140:ILE:HB	1.67	0.76
1:A:169:LEU:CD2	1:A:192:SER:HB3	2.15	0.76
1:A:89:LEU:HA	1:A:93:ASN:HB3	1.67	0.76
2:B:176:LEU:HG	2:B:257:LEU:HD22	1.67	0.75
1:A:94:HIS:O	1:A:98:GLU:HG3	1.85	0.75
2:B:38:HIS:HA	2:B:41:ARG:HG2	1.68	0.74
2:B:215:ARG:HG2	2:B:225:SER:HB3	1.68	0.74
1:A:168:ILE:HG21	2:B:116:LEU:HD22	1.70	0.73
2:B:117:LEU:O	2:B:139:THR:HG21	1.89	0.73
2:B:205:GLU:OE2	2:B:236:ARG:HG3	1.90	0.72
1:A:122:LEU:N	1:A:122:LEU:HD22	2.04	0.72
1:A:179:VAL:HG12	1:A:186:VAL:HG22	1.71	0.71
2:B:117:LEU:HD12	2:B:118:GLN:N	2.05	0.71
2:B:39:ARG:NH1	3:C:9:DG:O6	2.24	0.71
1:A:164:LEU:HA	1:A:167:LEU:HB3	1.70	0.71
1:A:136:VAL:CG1	1:A:140:LYS:HE3	2.21	0.71
1:A:224:GLN:HG2	1:A:263:PHE:HB3	1.73	0.71
1:A:307:HIS:HB3	1:A:344:GLN:NE2	2.07	0.70
2:B:167:GLU:O	2:B:170:ARG:HG2	1.92	0.70
1:A:163:GLU:O	1:A:167:LEU:N	2.24	0.70
1:A:176:LEU:O	1:A:190:SER:HB2	1.91	0.70
1:A:130:THR:O	1:A:134:MET:HG2	1.92	0.70
2:B:80:SER:OG	2:B:84:VAL:HG12	1.92	0.69
1:A:94:HIS:HA	1:A:97:ILE:HG12	1.74	0.69
4:D:8:DC:H2'	4:D:9:DT:C6	2.27	0.69
2:B:124:VAL:CA	2:B:262:ILE:HG22	2.23	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:116:LEU:O	2:B:120:LEU:HG	1.93	0.69
2:B:138:SER:O	2:B:141:GLN:NE2	2.21	0.69
1:A:169:LEU:HD13	1:A:192:SER:HB3	1.74	0.69
3:C:9:DG:N2	4:D:14:DC:O2	2.26	0.69
1:A:164:LEU:CD2	1:A:167:LEU:HD23	2.23	0.68
1:A:136:VAL:HG12	1:A:140:LYS:HE3	1.75	0.68
1:A:127:ASP:O	1:A:131:ILE:HG13	1.93	0.68
1:A:86:LYS:O	1:A:90:ALA:CB	2.27	0.68
1:A:343:LEU:HD22	1:A:343:LEU:H	1.58	0.67
1:A:162:GLN:HA	1:A:165:LYS:HE2	1.76	0.67
1:A:131:ILE:O	1:A:135:ALA:HB2	1.95	0.66
2:B:237:LEU:HD23	2:B:259:LEU:HD23	1.77	0.66
1:A:331:GLN:OE1	1:A:332:GLY:O	2.14	0.66
1:A:119:CYS:C	1:A:122:LEU:HD21	2.15	0.66
1:A:180:SER:OG	1:A:183:THR:HG21	1.90	0.66
1:A:169:LEU:HD22	1:A:192:SER:OG	1.95	0.66
2:B:264:THR:HB	2:B:265:PRO:HD2	1.78	0.66
3:C:19:DT:N3	3:C:20:DG:N7	2.44	0.66
2:B:202:LEU:O	2:B:205:GLU:N	2.29	0.65
2:B:159:LEU:O	2:B:214:PHE:HB3	1.97	0.65
2:B:215:ARG:NH1	2:B:216:CYS:SG	2.69	0.65
1:A:168:ILE:HG21	2:B:116:LEU:CD2	2.26	0.64
2:B:151:VAL:HG21	2:B:159:LEU:HD21	1.78	0.64
1:A:224:GLN:OE1	1:A:224:GLN:N	2.31	0.64
1:A:112:LEU:O	1:A:112:LEU:HD23	1.99	0.63
1:A:86:LYS:CA	1:A:90:ALA:HB2	2.03	0.63
1:A:136:VAL:O	1:A:140:LYS:N	2.31	0.63
1:A:164:LEU:O	1:A:168:ILE:HG22	1.98	0.63
1:A:304:VAL:HG23	1:A:343:LEU:HD12	1.81	0.62
1:A:300:GLU:HB2	1:A:301:PRO:HD3	1.81	0.62
2:B:57:GLN:OE1	2:B:60:ILE:HG21	2.00	0.62
1:A:111:GLU:HG2	2:B:73:VAL:HG11	1.80	0.62
1:A:132:LEU:HD21	2:B:45:GLU:HG3	1.81	0.62
1:A:213:VAL:HG21	1:A:221:LEU:CD2	2.28	0.62
1:A:121:ALA:O	1:A:122:LEU:HD13	2.00	0.62
1:A:113:SER:HB3	1:A:131:ILE:HD13	1.82	0.61
1:A:331:GLN:OE1	1:A:332:GLY:CA	2.47	0.61
1:A:158:PHE:HB2	2:B:75:TYR:CD1	2.36	0.61
1:A:186:VAL:HG21	1:A:204:TRP:CE3	2.35	0.61
2:B:202:LEU:HB2	2:B:203:PRO:HD2	1.83	0.61
1:A:174:GLY:HA2	1:A:341:GLY:O	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LEU:HD21	2:B:45:GLU:CB	2.30	0.61
1:A:166:HIS:HE1	2:B:82:PHE:CD1	2.19	0.61
4:D:5:DG:H3'	4:D:6:DT:C7	2.31	0.61
2:B:240:LEU:HD13	2:B:241:HIS:N	2.15	0.60
2:B:202:LEU:HB2	2:B:203:PRO:CD	2.29	0.60
4:D:6:DT:H73	4:D:6:DT:OP2	2.00	0.60
1:A:108:TYR:HB3	2:B:69:LEU:HG	1.83	0.60
1:A:119:CYS:C	1:A:122:LEU:CD2	2.69	0.60
1:A:331:GLN:NE2	1:A:334:LYS:CB	2.64	0.60
1:A:185:ARG:HA	1:A:208:THR:HG22	1.84	0.59
1:A:176:LEU:HD11	1:A:338:VAL:HG23	1.83	0.59
2:B:257:LEU:O	2:B:257:LEU:HD12	2.01	0.59
1:A:331:GLN:HE22	1:A:334:LYS:CB	2.16	0.58
1:A:121:ALA:N	1:A:122:LEU:HD22	2.18	0.58
2:B:137:SER:O	2:B:140:ILE:HG22	2.03	0.58
1:A:186:VAL:HG21	1:A:204:TRP:HE3	1.68	0.58
1:A:267:MET:SD	1:A:306:VAL:HG21	2.44	0.58
1:A:95:SER:HA	1:A:98:GLU:CD	2.23	0.58
2:B:211:GLU:O	2:B:230:ALA:HA	2.03	0.58
1:A:89:LEU:HA	1:A:93:ASN:CB	2.33	0.58
2:B:124:VAL:O	2:B:125:LEU:HD23	2.03	0.57
1:A:158:PHE:CE1	1:A:159:LEU:HG	2.40	0.57
2:B:68:VAL:O	2:B:72:SER:N	2.30	0.57
1:A:158:PHE:HA	2:B:54:PRO:HB2	1.86	0.57
1:A:133:ARG:HG3	2:B:45:GLU:OE2	2.04	0.57
2:B:65:LYS:HG2	4:D:10:DC:OP2	2.05	0.57
2:B:127:VAL:HG11	2:B:168:PHE:CE1	2.40	0.56
1:A:94:HIS:HA	1:A:97:ILE:CG1	2.35	0.56
1:A:172:ALA:HA	2:B:236:ARG:HE	1.71	0.56
1:A:155:LYS:HD3	1:A:156:PRO:N	2.20	0.56
1:A:169:LEU:CD1	1:A:192:SER:HB3	2.35	0.56
2:B:38:HIS:NE2	3:C:6:DA:H5'	2.21	0.56
1:A:125:LYS:O	1:A:126:PRO:C	2.40	0.56
3:C:5:DG:N2	3:C:6:DA:N1	2.53	0.56
1:A:127:ASP:HB2	1:A:130:THR:OG1	2.06	0.56
2:B:172:LEU:HD23	2:B:257:LEU:HD11	1.87	0.56
1:A:119:CYS:O	1:A:122:LEU:HD21	2.03	0.56
1:A:136:VAL:HG12	1:A:140:LYS:HG3	1.87	0.56
1:A:310:GLY:HA3	1:A:339:ALA:HB2	1.88	0.55
4:D:11:DA:H2'	4:D:12:DC:C6	2.41	0.55
2:B:116:LEU:HD12	2:B:117:LEU:H	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:67:SER:O	2:B:71:LEU:HG	2.07	0.55
1:A:95:SER:OG	1:A:99:ARG:NH2	2.39	0.55
4:D:6:DT:H2'	4:D:7:DT:C6	2.41	0.55
2:B:140:ILE:HG13	2:B:144:LEU:HD12	1.89	0.55
1:A:124:ARG:O	1:A:125:LYS:HG3	2.07	0.54
1:A:142:LEU:HD22	2:B:79:LYS:HG3	1.89	0.54
2:B:167:GLU:HA	2:B:170:ARG:NE	2.22	0.54
1:A:167:LEU:HD12	2:B:260:PHE:CD2	2.43	0.54
2:B:213:CYS:O	2:B:228:PHE:HA	2.07	0.54
2:B:259:LEU:HD11	2:B:261:ALA:HB2	1.90	0.54
2:B:235:GLY:HA3	2:B:259:LEU:HD21	1.90	0.54
2:B:128:THR:OG1	2:B:132:LEU:HD12	2.08	0.54
2:B:126:VAL:CG1	2:B:134:PHE:HB3	2.37	0.53
2:B:148:GLN:O	2:B:152:ILE:HB	2.08	0.53
2:B:78:ALA:O	2:B:81:PHE:HB3	2.08	0.53
2:B:127:VAL:HG11	2:B:168:PHE:CZ	2.43	0.53
1:A:166:HIS:CD2	1:A:170:GLU:HG2	2.43	0.53
2:B:157:TYR:CD1	2:B:165:ARG:HD3	2.43	0.53
2:B:240:LEU:HD23	2:B:260:PHE:CZ	2.43	0.53
1:A:138:HIS:HA	1:A:141:SER:OG	2.08	0.53
1:A:337:LEU:C	1:A:337:LEU:HD23	2.28	0.53
1:A:132:LEU:HD23	1:A:133:ARG:N	2.23	0.53
2:B:176:LEU:CD1	2:B:237:LEU:HD13	2.32	0.53
2:B:79:LYS:HE3	2:B:84:VAL:HG21	1.90	0.53
1:A:142:LEU:HD22	2:B:79:LYS:CG	2.40	0.52
1:A:113:SER:HB3	1:A:131:ILE:CG2	2.40	0.52
1:A:342:ARG:O	1:A:344:GLN:HG3	2.10	0.52
2:B:176:LEU:CG	2:B:257:LEU:HD22	2.36	0.52
1:A:116:VAL:HG13	1:A:138:HIS:HD2	1.74	0.52
1:A:86:LYS:C	1:A:90:ALA:CB	2.57	0.52
1:A:158:PHE:HD2	2:B:75:TYR:HD1	1.56	0.52
1:A:116:VAL:HG13	1:A:138:HIS:CD2	2.44	0.52
1:A:193:VAL:HG12	1:A:193:VAL:O	2.09	0.52
1:A:139:MET:HG2	2:B:49:LEU:CD2	2.36	0.51
2:B:116:LEU:HD12	2:B:117:LEU:N	2.25	0.51
1:A:122:LEU:N	1:A:122:LEU:CD2	2.73	0.51
1:A:81:GLN:O	1:A:85:ASP:HB2	2.09	0.51
1:A:136:VAL:HA	1:A:139:MET:HB3	1.93	0.51
2:B:161:HIS:HB2	2:B:212:ARG:HB2	1.92	0.51
1:A:158:PHE:CD2	2:B:75:TYR:HD1	2.28	0.51
4:D:10:DC:H2''	4:D:11:DA:C8	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:LYS:HD3	1:A:156:PRO:CD	2.41	0.51
1:A:260:ARG:HG3	1:A:310:GLY:O	2.10	0.51
1:A:337:LEU:HD23	1:A:338:VAL:N	2.25	0.51
2:B:126:VAL:HG12	2:B:134:PHE:HB3	1.92	0.51
1:A:155:LYS:HB3	1:A:156:PRO:HD3	1.93	0.51
1:A:307:HIS:HB3	1:A:344:GLN:CD	2.30	0.51
1:A:139:MET:CG	2:B:49:LEU:HD22	2.36	0.50
1:A:134:MET:O	1:A:137:SER:OG	2.24	0.50
1:A:261:ARG:O	1:A:309:THR:HA	2.11	0.50
1:A:158:PHE:HA	2:B:54:PRO:CB	2.42	0.50
2:B:173:HIS:HB3	2:B:174:TRP:CE3	2.46	0.50
1:A:218:VAL:HG13	1:A:219:ASP:OD2	2.11	0.50
2:B:48:ARG:O	2:B:52:LEU:HD23	2.12	0.50
1:A:104:LYS:HD2	1:A:108:TYR:CZ	2.46	0.50
1:A:136:VAL:HG11	1:A:140:LYS:HE3	1.94	0.50
1:A:132:LEU:HD21	2:B:45:GLU:CG	2.40	0.50
1:A:264:ILE:O	1:A:264:ILE:HG23	2.12	0.50
2:B:235:GLY:HA2	2:B:261:ALA:HB2	1.93	0.50
2:B:262:ILE:O	2:B:262:ILE:HG13	2.11	0.50
2:B:161:HIS:HB2	2:B:212:ARG:CB	2.42	0.50
2:B:71:LEU:HA	2:B:74:SER:OG	2.12	0.49
2:B:206:ASN:ND2	2:B:206:ASN:O	2.45	0.49
1:A:311:TYR:CZ	1:A:338:VAL:HG11	2.47	0.49
2:B:117:LEU:HB2	2:B:139:THR:CG2	2.43	0.49
2:B:42:LEU:O	2:B:45:GLU:N	2.42	0.49
2:B:159:LEU:HD12	2:B:159:LEU:N	2.28	0.49
2:B:176:LEU:HD23	2:B:176:LEU:N	2.27	0.49
1:A:105:MET:O	1:A:109:ILE:N	2.25	0.49
2:B:38:HIS:HA	2:B:41:ARG:CG	2.39	0.49
4:D:13:DG:H2"	4:D:14:DC:C6	2.48	0.49
2:B:230:ALA:O	2:B:266:LEU:HB2	2.12	0.49
1:A:337:LEU:HD21	1:A:339:ALA:CB	2.43	0.49
2:B:53:LEU:HG	2:B:55:PHE:CD2	2.48	0.49
2:B:44:THR:O	2:B:48:ARG:HG3	2.13	0.48
1:A:304:VAL:HG23	1:A:343:LEU:CD1	2.43	0.48
1:A:172:ALA:HB3	1:A:342:ARG:CD	2.43	0.48
2:B:114:GLU:O	2:B:117:LEU:HG	2.13	0.48
2:B:170:ARG:HA	2:B:174:TRP:CE3	2.48	0.48
3:C:12:DT:H2"	3:C:13:DG:C8	2.47	0.48
2:B:59:VAL:HG22	2:B:59:VAL:O	2.13	0.48
1:A:203:GLU:O	1:A:207:SER:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:75:TYR:O	2:B:78:ALA:HB3	2.13	0.48
3:C:18:DC:H2''	3:C:19:DT:O5'	2.12	0.48
2:B:234:GLN:O	2:B:234:GLN:HG3	2.13	0.48
1:A:179:VAL:HA	1:A:186:VAL:HA	1.96	0.48
1:A:95:SER:O	1:A:99:ARG:HG2	2.13	0.48
2:B:207:ALA:O	2:B:234:GLN:HA	2.13	0.48
2:B:206:ASN:HA	2:B:235:GLY:O	2.14	0.48
1:A:178:ILE:O	1:A:187:VAL:N	2.32	0.48
2:B:64:ASP:OD1	2:B:67:SER:HB2	2.13	0.48
3:C:17:DA:H2''	3:C:18:DC:C6	2.49	0.48
1:A:136:VAL:O	1:A:140:LYS:HG3	2.14	0.48
3:C:19:DT:H3'	3:C:20:DG:C5'	2.44	0.48
3:C:15:DG:H2''	3:C:16:DA:OP2	2.14	0.47
1:A:108:TYR:CA	1:A:111:GLU:HB3	2.38	0.47
1:A:176:LEU:HD11	1:A:338:VAL:CG2	2.44	0.47
2:B:157:TYR:CD1	2:B:165:ARG:HB2	2.49	0.47
3:C:7:DT:H6	3:C:7:DT:H3'	1.78	0.47
1:A:264:ILE:HD11	1:A:305:VAL:CG1	2.44	0.47
1:A:179:VAL:O	1:A:336:CYS:HB2	2.15	0.47
1:A:172:ALA:HB3	1:A:342:ARG:HD2	1.96	0.47
1:A:201:GLN:O	1:A:205:PHE:HB2	2.14	0.47
1:A:304:VAL:HG22	1:A:304:VAL:O	2.15	0.47
1:A:331:GLN:HE22	1:A:334:LYS:HB2	1.69	0.47
3:C:10:DC:H2''	3:C:11:DG:OP1	2.15	0.47
2:B:235:GLY:CA	2:B:261:ALA:HB2	2.45	0.47
2:B:114:GLU:HB3	2:B:117:LEU:CD2	2.38	0.47
2:B:38:HIS:CD2	3:C:6:DA:H5'	2.50	0.46
1:A:105:MET:SD	1:A:106:THR:N	2.89	0.46
1:A:207:SER:HB3	1:A:211:ASP:CB	2.44	0.46
1:A:125:LYS:O	1:A:127:ASP:N	2.48	0.46
1:A:156:PRO:HG2	1:A:157:SER:H	1.80	0.46
2:B:59:VAL:HG22	2:B:62:LYS:HB3	1.96	0.46
1:A:181:CYS:O	1:A:225:LEU:HD12	2.16	0.46
1:A:189:VAL:O	1:A:189:VAL:HG23	2.15	0.46
2:B:144:LEU:HD13	2:B:146:PHE:HE1	1.80	0.46
2:B:44:THR:HA	2:B:47:ASP:OD2	2.16	0.46
4:D:5:DG:H3'	4:D:6:DT:H72	1.98	0.46
1:A:155:LYS:HD3	1:A:156:PRO:HD3	1.96	0.46
1:A:264:ILE:HD11	1:A:305:VAL:HG11	1.98	0.46
2:B:117:LEU:HD13	2:B:139:THR:HB	1.97	0.46
2:B:172:LEU:N	2:B:172:LEU:HD12	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:259:LEU:HD11	2:B:261:ALA:CB	2.46	0.46
3:C:17:DA:H2''	3:C:18:DC:C5	2.50	0.46
1:A:185:ARG:O	1:A:185:ARG:HG2	2.16	0.45
1:A:214:HIS:CE1	1:A:268:ARG:HB3	2.51	0.45
2:B:53:LEU:HG	2:B:55:PHE:HD2	1.82	0.45
1:A:158:PHE:CA	2:B:54:PRO:HB2	2.46	0.45
1:A:188:TYR:CZ	1:A:190:SER:HA	2.52	0.45
2:B:125:LEU:CD1	2:B:263:ALA:HB3	2.46	0.45
1:A:331:GLN:OE1	1:A:332:GLY:C	2.55	0.45
1:A:337:LEU:HD21	1:A:339:ALA:HB2	1.99	0.45
1:A:216:ASP:OD2	1:A:266:ARG:NH1	2.50	0.45
1:A:113:SER:HB2	1:A:119:CYS:SG	2.57	0.45
2:B:203:PRO:HA	2:B:204:PRO:HA	1.59	0.45
2:B:71:LEU:HD11	2:B:149:SER:HB3	1.98	0.45
2:B:82:PHE:CD2	2:B:83:ASP:HB2	2.52	0.45
4:D:8:DC:H2''	4:D:9:DT:O5'	2.15	0.45
1:A:180:SER:HB2	1:A:183:THR:HG21	1.94	0.45
2:B:56:PRO:O	2:B:60:ILE:HB	2.17	0.45
2:B:64:ASP:OD2	2:B:67:SER:N	2.49	0.45
1:A:180:SER:CB	1:A:183:THR:HG22	2.18	0.45
1:A:115:MET:CE	1:A:115:MET:HA	2.47	0.44
2:B:211:GLU:HG2	2:B:212:ARG:H	1.82	0.44
1:A:91:ARG:NE	3:C:12:DT:OP1	2.50	0.44
4:D:16:DA:C3'	4:D:17:DT:H5''	2.47	0.44
2:B:87:LYS:HD3	2:B:87:LYS:HA	1.65	0.44
1:A:117:PRO:HD2	1:A:138:HIS:NE2	2.33	0.44
1:A:181:CYS:SG	1:A:182:GLU:N	2.87	0.44
2:B:57:GLN:HA	2:B:60:ILE:CG2	2.47	0.44
4:D:8:DC:H4'	4:D:9:DT:OP1	2.18	0.44
2:B:152:ILE:O	2:B:152:ILE:HG12	2.17	0.44
4:D:4:DA:H2'	4:D:4:DA:N3	2.31	0.44
1:A:168:ILE:HD12	1:A:171:ALA:HB3	2.00	0.44
1:A:115:MET:O	1:A:117:PRO:HD3	2.18	0.43
1:A:211:ASP:OD1	1:A:211:ASP:N	2.51	0.43
1:A:224:GLN:HG2	1:A:263:PHE:CB	2.46	0.43
1:A:331:GLN:CD	1:A:332:GLY:N	2.67	0.43
2:B:159:LEU:HD23	2:B:214:PHE:CE1	2.53	0.43
2:B:176:LEU:HG	2:B:257:LEU:CD2	2.42	0.43
2:B:235:GLY:HA2	2:B:261:ALA:CB	2.49	0.43
3:C:14:DA:H1'	3:C:15:DG:C5'	2.42	0.43
1:A:132:LEU:HA	1:A:135:ALA:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:ASP:OD2	1:A:328:GLU:O	2.35	0.43
2:B:161:HIS:HB3	2:B:164:ASP:OD1	2.18	0.43
2:B:240:LEU:HD13	2:B:241:HIS:H	1.80	0.43
3:C:4:DG:H2''	3:C:5:DG:O5'	2.18	0.43
1:A:311:TYR:CZ	1:A:338:VAL:CG1	3.02	0.43
1:A:108:TYR:CB	2:B:69:LEU:HG	2.48	0.42
2:B:141:GLN:OE1	2:B:141:GLN:N	2.52	0.42
1:A:136:VAL:HG21	2:B:45:GLU:OE1	2.20	0.42
2:B:176:LEU:HD22	2:B:237:LEU:HD13	2.01	0.42
1:A:167:LEU:HD12	2:B:260:PHE:HD2	1.84	0.42
2:B:136:ALA:HB1	2:B:140:ILE:HG21	2.00	0.42
1:A:260:ARG:NH1	1:A:260:ARG:HB2	2.35	0.42
2:B:129:ALA:HB3	2:B:256:GLN:CD	2.40	0.42
1:A:99:ARG:HG3	1:A:100:ARG:N	2.35	0.42
2:B:224:ASN:ND2	2:B:224:ASN:O	2.53	0.42
1:A:190:SER:O	1:A:193:VAL:CG2	2.56	0.42
2:B:77:ARG:CG	2:B:134:PHE:HE1	2.33	0.42
1:A:105:MET:O	1:A:109:ILE:HG12	2.19	0.42
1:A:331:GLN:HE22	1:A:334:LYS:HG2	1.85	0.42
1:A:124:ARG:O	1:A:125:LYS:CG	2.67	0.42
1:A:158:PHE:O	2:B:75:TYR:HE1	2.03	0.42
1:A:195:PRO:HD2	1:A:199:GLN:OE1	2.19	0.41
2:B:77:ARG:HG2	2:B:134:PHE:HE1	1.84	0.41
2:B:70:ARG:CG	2:B:153:HIS:HB3	2.50	0.41
1:A:94:HIS:CA	1:A:97:ILE:HG12	2.48	0.41
1:A:137:SER:O	1:A:141:SER:N	2.35	0.41
1:A:263:PHE:CZ	1:A:308:CYS:HB2	2.55	0.41
1:A:169:LEU:O	1:A:173:ASP:HA	2.20	0.41
1:A:207:SER:HB3	1:A:211:ASP:HB2	2.01	0.41
2:B:133:VAL:HG22	2:B:156:VAL:HG12	2.02	0.41
1:A:94:HIS:ND1	1:A:97:ILE:HD11	2.35	0.41
1:A:102:ARG:NH2	3:C:8:DT:H2''	2.36	0.41
2:B:167:GLU:OE1	2:B:170:ARG:HD2	2.20	0.41
1:A:219:ASP:OD1	1:A:222:ARG:NH2	2.32	0.41
1:A:121:ALA:CA	1:A:122:LEU:HD22	2.50	0.41
1:A:331:GLN:HE22	1:A:334:LYS:CG	2.34	0.41
2:B:48:ARG:O	2:B:52:LEU:N	2.44	0.41
4:D:7:DT:H2''	4:D:8:DC:C6	2.55	0.41
1:A:104:LYS:HD2	1:A:108:TYR:CE1	2.56	0.40
2:B:140:ILE:CG1	2:B:144:LEU:HD12	2.51	0.40
1:A:328:GLU:HB3	1:A:329:ALA:H	1.70	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:LYS:CD	1:A:156:PRO:HD3	2.51	0.40
1:A:94:HIS:O	1:A:97:ILE:N	2.49	0.40
2:B:46:LEU:HD23	2:B:65:LYS:HD2	2.03	0.40
4:D:13:DG:H2''	4:D:14:DC:C5	2.55	0.40
4:D:4:DA:H3'	4:D:5:DG:C5'	2.51	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:GLU:OE1	2:B:239:TYR:OH[3_545]	1.76	0.44
1:A:81:GLN:N	4:D:17:DT:OP2[4_445]	2.06	0.14

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	164/279 (59%)	131 (80%)	29 (18%)	4 (2%)	7	45
2	B	154/241 (64%)	138 (90%)	15 (10%)	1 (1%)	28	70
All	All	318/520 (61%)	269 (85%)	44 (14%)	5 (2%)	11	53

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	ALA
1	A	156	PRO
1	A	327	PRO
2	B	204	PRO
1	A	330	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	156/245 (64%)	142 (91%)	14 (9%)	11	43
2	B	148/206 (72%)	135 (91%)	13 (9%)	12	44
All	All	304/451 (67%)	277 (91%)	27 (9%)	11	44

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	GLU
1	A	115	MET
1	A	119	CYS
1	A	122	LEU
1	A	132	LEU
1	A	158	PHE
1	A	161	ASP
1	A	183	THR
1	A	185	ARG
1	A	211	ASP
1	A	225	LEU
1	A	302	HIS
1	A	307	HIS
1	A	328	GLU
2	B	77	ARG
2	B	80	SER
2	B	81	PHE
2	B	86	LEU
2	B	87	LYS
2	B	89	THR
2	B	116	LEU
2	B	161	HIS
2	B	205	GLU
2	B	209	PHE
2	B	214	PHE
2	B	239	TYR
2	B	240	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	HIS
1	A	307	HIS
1	A	344	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](#)

4 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	$\# Z > 2$
5	PEG	A	401	-	6,6,6	0.49	0	5,5,5	0.29	0
6	CIT	B	301	-	3,12,12	1.25	0	3,17,17	2.38	2 (66%)
5	PEG	B	302	-	6,6,6	0.50	0	5,5,5	0.27	0
6	CIT	C	101	-	3,12,12	1.24	0	3,17,17	2.43	2 (66%)

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
5	PEG	A	401	-	-	0/4/4/4	0/0/0/0
6	CIT	B	301	-	-	0/6/16/16	0/0/0/0
5	PEG	B	302	-	-	0/4/4/4	0/0/0/0
6	CIT	C	101	-	-	0/6/16/16	0/0/0/0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
6	C	101	CIT	C3-C4-C5	-3.29	109.81	114.95
6	B	301	CIT	C3-C2-C1	-2.88	110.45	114.95
6	B	301	CIT	C3-C4-C5	-2.72	110.71	114.95
6	C	101	CIT	C3-C2-C1	-2.44	111.13	114.95

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 [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	176/279 (63%)	0.00	6 (3%)	46 36	130, 205, 272, 305	0
2	B	166/241 (68%)	0.14	10 (6%)	23 16	150, 213, 268, 331	0
3	C	17/17 (100%)	0.69	2 (11%)	5 6	209, 250, 322, 357	0
4	D	14/14 (100%)	0.44	2 (14%)	3 4	215, 243, 282, 295	0
All	All	373/551 (67%)	0.11	20 (5%)	26 21	130, 213, 275, 357	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	7	DT	5.3
2	B	211	GLU	3.6
4	D	17	DT	3.5
1	A	128	LYS	3.4
2	B	87	LYS	3.1
1	A	331	GLN	2.8
2	B	143	TYR	2.8
2	B	230	ALA	2.7
4	D	4	DA	2.7
2	B	38	HIS	2.7
2	B	57	GLN	2.7
1	A	220	LYS	2.6
1	A	189	VAL	2.6
1	A	81	GLN	2.5
2	B	229	LEU	2.5
2	B	142	ASP	2.2
3	C	19	DT	2.1
2	B	89	THR	2.1
1	A	202	SER	2.1
2	B	79	LYS	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(\AA^2)	Q<0.9
5	PEG	B	302	7/7	0.67	0.42	-	180,207,233,242	0
6	CIT	B	301	13/13	0.63	0.71	-	226,239,244,253	0
5	PEG	A	401	7/7	0.45	0.34	-	184,191,211,215	0
6	CIT	C	101	13/13	0.81	0.21	-	205,238,248,284	0

6.5 Other polymers [i](#)

There are no such residues in this entry.