



# Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 11:31 am BST

PDB ID : 6V37  
Title : K2P2.1(TREK-1)I110D:RuR:ML335 bound channel structure  
Authors : Pope, L.; Lolicato, M.; Minor, D.L.  
Deposited on : 2019-11-25  
Resolution : 2.80 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](mailto: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.

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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  
buster-report : 1.1.7 (2018)  
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

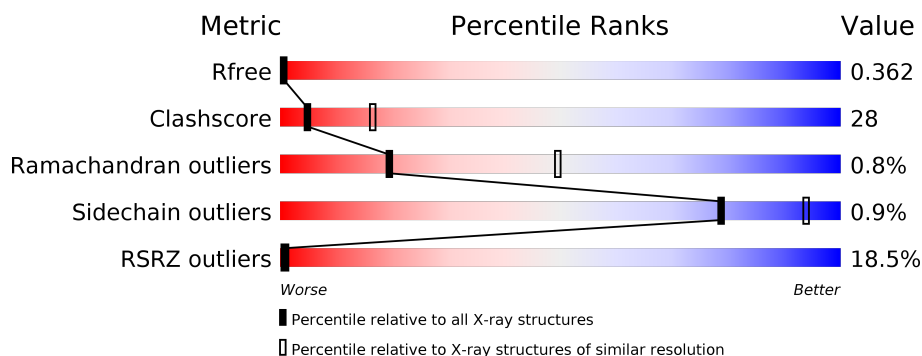
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	312	<div> <div>17%</div> <div>48%</div> <div>35%</div> <div>..</div> <div>16%</div> </div>
1	B	312	<div> <div>15%</div> <div>42%</div> <div>45%</div> <div>13%</div> </div>

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
3	CD	B	906	-	-	-	X
4	D12	A	405	-	-	-	X

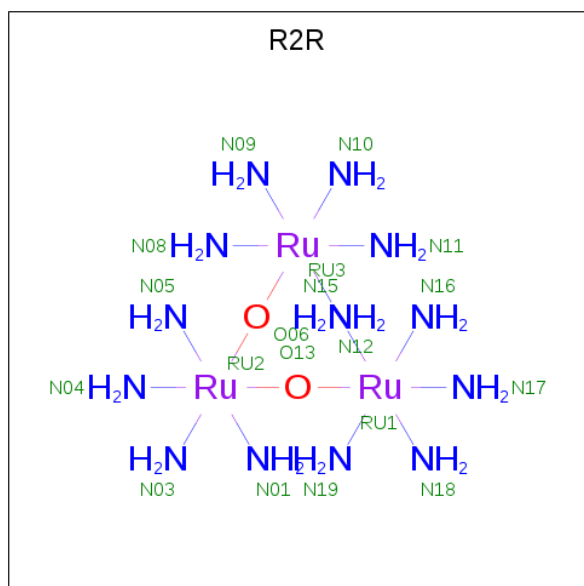


In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Potassium channel subfamily K member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total 2060	C 1372	N 325	O 358	S 5	0	0	0
1	B	271	Total 2120	C 1421	N 333	O 361	S 5	0	1	0

- Molecule 2 is ruthenium(6+) azanide pentaamino(oxido)ruthenium (1/4/2) (three-letter code: R2R) (formula:  $\text{H}_{28}\text{N}_{14}\text{O}_2\text{Ru}_3$ ) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	N	O	Ru	0	0
			19	14	2	3		

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

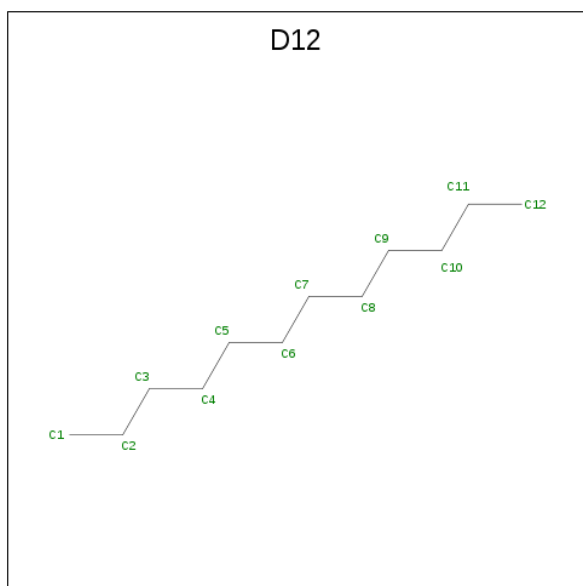
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total Cd 2 2	0	0

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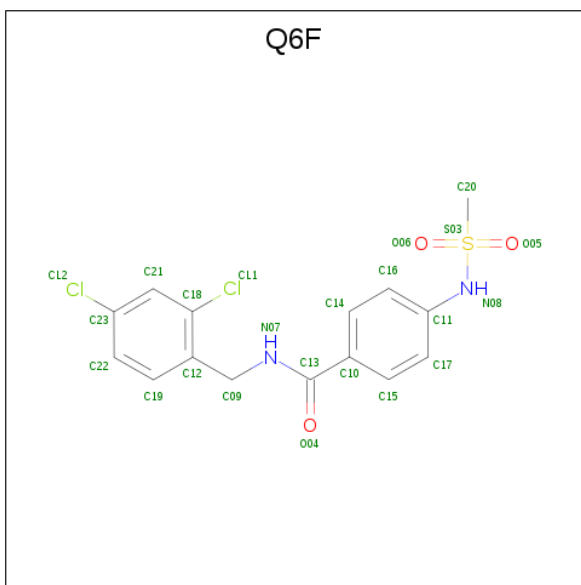
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cd	0	0
			1	1		

- Molecule 4 is DODECANE (three-letter code: D12) (formula:  $C_{12}H_{26}$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	C	0	0
			12	12		
4	A	1	Total	C	0	0
			12	12		
4	A	1	Total	C	0	0
			38	12	26	
4	B	1	Total	C	0	0
			12	12		

- Molecule 5 is N-[(2,4-dichlorophenyl)methyl]-4-[(methylsulfonyl)amino]benzamide (three-letter code: Q6F) (formula:  $C_{15}H_{14}Cl_2N_2O_3S$ ).

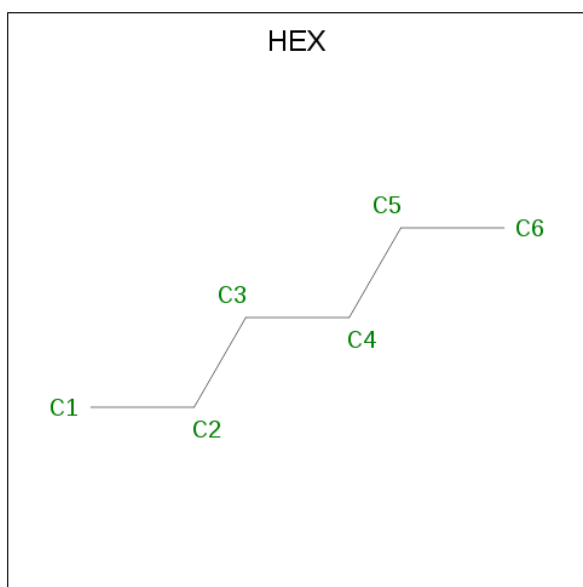


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	Cl	N	O	S	0	0
			23	15	2	2	3	1		
5	B	1	Total	C	Cl	N	O	S	0	0
			23	15	2	2	3	1		

- Molecule 6 is POTASSIUM ION (three-letter code: K) (formula: K).

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

- Molecule 7 is HEXANE (three-letter code: HEX) (formula: C<sub>6</sub>H<sub>14</sub>).



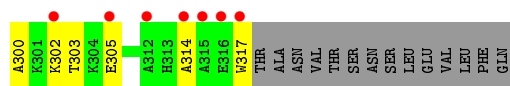
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	H	0	0
			20	6	14		

- Molecule 8 is water.

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







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.02Å 118.73Å 129.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.99 – 2.80 46.48 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.4 (14.99-2.80) 99.8 (46.48-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.20 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, $R_{free}$	0.272 , 0.317 0.291 , 0.362	Depositor DCC
$R_{free}$ test set	1195 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	118.1	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 125.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	4349	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	145.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: D12, Q6F, HEX, CD, K, R2R

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	0.28	0/2110	0.44	0/2867
1	B	0.26	0/2176	0.40	0/2958
All	All	0.27	0/4286	0.42	0/5825

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	262	SER	Peptide

### 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	2060	0	2094	118	0
1	B	2120	0	2174	147	0
2	A	19	0	0	4	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
4	A	36	26	78	0	0
4	B	12	0	26	4	0
5	A	23	0	0	3	0
5	B	23	0	0	3	0
6	A	4	0	0	0	0
6	B	1	0	0	0	0
7	B	6	14	14	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
All	All	4309	40	4386	240	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 28.

All (240) 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:114:ILE:HD12	1:B:116:PRO:HD3	1.30	1.14
1:B:109:ALA:HB1	1:B:114:ILE:HG12	1.33	1.10
1:B:250:THR:O	1:B:251:THR:HG22	1.73	0.89
1:A:110:ASP:OD2	2:A:401:R2R:N01	2.07	0.88
1:B:245:VAL:HG13	1:B:277:TRP:HZ2	1.42	0.84
1:A:88:LEU:HD11	1:A:99:LEU:HD13	1.60	0.82
1:A:89:ARG:O	1:A:92:PRO:HD3	1.80	0.81
1:B:41:VAL:HG12	1:B:45:LYS:HE3	1.60	0.81
1:A:80:ILE:CD1	1:B:114:ILE:HG21	2.10	0.81
1:A:91:HIS:NE2	1:B:95:SER:HB3	1.98	0.79
1:A:106:ILE:HD13	1:B:106:ILE:HD11	1.65	0.76
1:A:133:PHE:HA	1:B:62:ALA:HB2	1.69	0.75
1:B:37:SER:HA	1:B:40:ASN:ND2	2.03	0.74
1:B:40:ASN:HA	1:B:43:LYS:HE3	1.69	0.74
1:A:84:ARG:NH2	1:B:116:PRO:HB2	2.03	0.74
1:A:187:LYS:C	1:A:191:LYS:HE3	2.08	0.73
1:A:185:PHE:O	1:A:189:ILE:HG12	1.89	0.72
1:A:135:PHE:CZ	1:A:150:PRO:HD3	2.23	0.72
1:A:156:LYS:NZ	1:B:240:ASP:OD1	2.23	0.71
1:B:109:ALA:HB1	1:B:114:ILE:CG1	2.18	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:TYR:O	1:A:267:LEU:HD12	1.90	0.71
1:A:189:ILE:HG13	1:A:190:ALA:H	1.56	0.70
1:A:189:ILE:HG13	1:A:190:ALA:N	2.06	0.70
1:A:152:THR:HG22	1:A:154:GLY:H	1.57	0.70
1:B:194:ASP:HA	1:B:197:ILE:HG22	1.75	0.69
1:B:231:LYS:HG3	1:B:236:TRP:O	1.92	0.69
1:A:191:LYS:HE2	1:A:191:LYS:H	1.57	0.69
1:B:111:ASN:HB3	1:B:151:ARG:NH2	2.08	0.68
1:A:106:ILE:HG21	1:B:106:ILE:HD11	1.77	0.67
1:A:80:ILE:HD13	1:B:114:ILE:HG21	1.75	0.67
1:B:84:ARG:CG	1:B:88:LEU:HD13	2.24	0.67
1:B:194:ASP:O	1:B:197:ILE:HG22	1.94	0.66
1:A:103:ILE:HG21	1:B:116:PRO:HG2	1.77	0.65
1:B:84:ARG:HG2	1:B:88:LEU:HD13	1.79	0.65
1:A:75:SER:O	1:A:79:THR:HG23	1.97	0.65
1:B:210:SER:O	1:B:213:ILE:HG22	1.96	0.64
1:B:41:VAL:CG1	1:B:45:LYS:HE3	2.27	0.64
1:A:106:ILE:HG21	1:B:106:ILE:CD1	2.26	0.64
1:B:59:ILE:O	1:B:63:THR:HG23	1.98	0.64
1:A:191:LYS:CE	1:A:191:LYS:H	2.12	0.63
1:B:45:LYS:O	1:B:49:THR:HG23	1.98	0.63
1:B:222:PHE:CE1	1:B:280:VAL:HG12	2.33	0.63
1:A:260:GLY:HA3	5:A:406:Q6F:C14	2.29	0.63
1:B:167:ILE:HB	1:B:168:PRO:HD3	1.80	0.63
1:B:109:ALA:CB	1:B:114:ILE:HG12	2.21	0.62
1:A:249:LEU:HB2	1:A:285:PHE:CE2	2.34	0.62
1:A:129:LEU:HD23	1:B:66:LYS:NZ	2.14	0.62
1:A:250:THR:O	1:A:251:THR:OG1	2.18	0.62
1:A:186:GLY:O	1:A:191:LYS:NZ	2.33	0.62
1:B:249:LEU:HB2	1:B:285:PHE:CE2	2.35	0.61
1:A:190:ALA:H	1:A:191:LYS:HZ3	1.49	0.61
1:B:218:GLY:HA3	1:B:284:TYR:CZ	2.35	0.61
1:A:220:VAL:HA	1:A:224:ALA:HB3	1.83	0.60
1:B:37:SER:HA	1:B:40:ASN:HD22	1.66	0.60
1:A:84:ARG:NH2	1:B:116:PRO:O	2.33	0.60
1:A:41:VAL:C	1:A:43:LYS:H	2.03	0.60
1:B:114:ILE:CD1	1:B:116:PRO:HD3	2.20	0.60
1:A:107:VAL:O	1:A:110:ASP:HB3	2.02	0.59
1:A:190:ALA:N	1:A:191:LYS:HZ3	2.00	0.59
1:B:49:THR:O	1:B:53:VAL:HG23	2.02	0.59
1:B:110:ASP:O	1:B:113:GLY:N	2.30	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:LEU:HG	1:B:96:ASP:OD1	2.03	0.58
1:B:147:ASN:OD1	1:B:148:ILE:N	2.36	0.58
1:A:47:VAL:HG22	1:B:176:GLY:HA3	1.85	0.58
1:A:284:TYR:O	1:A:288:VAL:HG23	2.03	0.58
1:A:134:PHE:CE1	1:A:138:THR:HG23	2.39	0.58
1:B:198:LYS:HE3	1:B:199:TRP:CH2	2.38	0.58
1:A:303:THR:O	1:A:307:VAL:HG23	2.03	0.58
1:B:213:ILE:HD11	1:B:217:PHE:CE2	2.38	0.58
1:B:224:ALA:O	1:B:228:VAL:HG23	2.03	0.57
1:B:188:GLY:HA2	1:B:191:LYS:NZ	2.19	0.57
1:A:91:HIS:CD2	1:B:95:SER:HB3	2.38	0.57
1:A:147:ASN:OD1	1:A:148:ILE:N	2.38	0.57
1:B:84:ARG:NH2	1:B:100:ASP:OD1	2.37	0.57
1:B:98:GLU:HA	1:B:101:GLU:HG2	1.85	0.57
1:B:206:ILE:HA	1:B:209:ILE:HD13	1.87	0.57
1:A:87:PHE:CD2	1:A:99:LEU:HD12	2.39	0.57
1:A:190:ALA:HA	1:A:193:GLU:CG	2.35	0.56
1:A:138:THR:HB	1:A:144:GLY:HA2	1.86	0.56
1:B:135:PHE:CZ	1:B:150:PRO:HD3	2.41	0.55
1:B:60:ILE:O	1:B:64:VAL:HG23	2.06	0.55
1:A:258:VAL:HG12	1:A:260:GLY:H	1.71	0.55
1:B:260:GLY:HA3	5:B:904:Q6F:C14	2.36	0.55
1:A:302:LYS:O	1:A:305:GLU:HG2	2.07	0.55
1:B:41:VAL:O	1:B:45:LYS:HG3	2.07	0.55
1:A:179:ASP:O	1:A:183:THR:HG23	2.06	0.55
1:B:84:ARG:HG3	1:B:88:LEU:HD13	1.89	0.54
1:B:194:ASP:HA	1:B:197:ILE:CG2	2.37	0.54
1:B:213:ILE:HD11	1:B:217:PHE:HE2	1.71	0.54
1:B:216:LEU:O	1:B:220:VAL:HG23	2.08	0.54
1:A:62:ALA:HB2	1:B:133:PHE:HA	1.90	0.54
1:B:84:ARG:HG3	1:B:88:LEU:CD1	2.38	0.54
1:B:148:ILE:O	1:B:148:ILE:HG13	2.08	0.54
1:B:111:ASN:HB3	1:B:151:ARG:HH22	1.71	0.54
1:B:93:CYS:O	1:B:95:SER:N	2.40	0.54
1:A:131:SER:HA	5:A:406:Q6F:C17	2.37	0.54
1:A:135:PHE:O	1:A:139:VAL:HG23	2.08	0.53
1:A:135:PHE:HZ	1:A:150:PRO:HD3	1.69	0.53
1:B:245:VAL:HG13	1:B:277:TRP:CZ2	2.33	0.53
1:A:218:GLY:HA3	1:A:284:TYR:CZ	2.43	0.53
1:B:243:TYR:O	1:B:247:ILE:HG12	2.09	0.53
1:A:229:ILE:CG2	1:A:233:ILE:HD12	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:LEU:HD11	1:B:283:ALA:N	2.24	0.53
1:B:39:ILE:HD12	1:B:40:ASN:N	2.24	0.53
1:A:100:ASP:O	1:A:104:GLN:HG3	2.08	0.53
1:B:194:ASP:CA	1:B:197:ILE:HG22	2.39	0.52
1:A:240:ASP:OD1	1:B:156:LYS:NZ	2.42	0.52
1:B:56:LEU:HD21	1:B:60:ILE:HD11	1.91	0.52
1:B:195:THR:O	1:B:198:LYS:HE2	2.10	0.52
1:B:98:GLU:O	1:B:101:GLU:HB2	2.10	0.52
1:B:186:GLY:HA2	1:B:189:ILE:HG22	1.90	0.52
1:A:134:PHE:HE1	1:A:138:THR:CG2	2.23	0.52
1:A:190:ALA:HA	1:A:193:GLU:HG3	1.91	0.52
1:B:302:LYS:O	1:B:305:GLU:HG2	2.09	0.52
1:A:292:ILE:O	1:A:296:LEU:HG	2.10	0.52
1:A:42:MET:O	1:A:44:TRP:HE3	1.92	0.52
1:A:274:VAL:O	1:A:277:TRP:HB3	2.11	0.51
1:A:106:ILE:HD13	1:B:106:ILE:CD1	2.37	0.51
1:B:106:ILE:O	1:B:110:ASP:N	2.43	0.51
1:A:267:LEU:O	1:A:268:ASP:OD1	2.28	0.51
1:A:259:ALA:HB3	5:A:406:Q6F:CL1	2.47	0.51
1:A:254:PHE:HD1	1:B:146:GLY:HA2	1.76	0.50
1:B:218:GLY:HA3	1:B:284:TYR:OH	2.10	0.50
1:A:188:GLY:O	1:A:192:VAL:HG22	2.12	0.50
1:A:191:LYS:N	1:A:191:LYS:HD3	2.27	0.50
1:A:299:ILE:O	1:A:303:THR:HG23	2.11	0.50
1:B:314:ALA:HB3	1:B:317:TRP:C	2.32	0.50
1:A:134:PHE:CE1	1:A:138:THR:CG2	2.95	0.49
2:A:401:R2R:N19	2:A:401:R2R:N01	2.60	0.49
1:A:88:LEU:CD1	1:A:99:LEU:HD13	2.38	0.49
1:A:134:PHE:HA	1:A:275:TRP:CH2	2.47	0.49
1:A:129:LEU:HD23	1:B:66:LYS:HZ2	1.77	0.49
1:B:296:LEU:O	1:B:300:ALA:N	2.41	0.49
1:B:56:LEU:CD2	4:B:903:D12:H31	2.43	0.49
1:B:173:LEU:O	1:B:177:VAL:HG23	2.12	0.49
1:B:181:LEU:HD22	1:B:280:VAL:HG13	1.95	0.48
1:B:100:ASP:O	1:B:104:GLN:HG3	2.13	0.48
1:B:84:ARG:O	1:B:88:LEU:HD13	2.12	0.48
1:A:181:LEU:HD22	1:A:280:VAL:HG13	1.95	0.48
1:A:70:GLN:N	1:A:71:PRO:CD	2.77	0.48
1:B:298:VAL:O	1:B:302:LYS:HG2	2.14	0.48
1:A:247:ILE:HD12	1:A:254:PHE:CE2	2.49	0.48
1:A:313:HIS:O	1:A:317:TRP:N	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:PHE:HD2	1:A:99:LEU:HD12	1.79	0.48
1:A:58:LEU:HD21	1:B:140:ILE:CD1	2.43	0.48
1:B:299:ILE:O	1:B:303:THR:HG23	2.13	0.48
1:A:180:GLN:OE1	1:B:43:LYS:NZ	2.47	0.48
1:A:261:GLY:O	1:A:262:SER:OG	2.26	0.47
1:B:213:ILE:HG13	1:B:217:PHE:CD2	2.49	0.47
1:A:142:THR:HA	1:B:252:ILE:HG12	1.94	0.47
2:A:401:R2R:N10	2:A:401:R2R:N05	2.61	0.47
1:B:149:SER:HB2	1:B:150:PRO:HD2	1.96	0.47
1:B:98:GLU:HA	1:B:101:GLU:CG	2.44	0.47
1:B:249:LEU:HB2	1:B:285:PHE:CD2	2.50	0.47
1:A:266:TYR:C	1:A:267:LEU:HD12	2.34	0.47
1:A:42:MET:O	1:A:44:TRP:CE3	2.68	0.47
1:B:209:ILE:HD12	1:B:209:ILE:H	1.78	0.47
1:B:250:THR:O	1:B:251:THR:CG2	2.55	0.47
1:B:70:GLN:N	1:B:71:PRO:CD	2.78	0.47
1:A:195:THR:O	1:A:196:PHE:HB2	2.15	0.47
1:A:147:ASN:ND2	1:A:258:VAL:HG21	2.30	0.47
1:A:43:LYS:O	1:A:43:LYS:HD2	2.14	0.46
1:A:140:ILE:HG22	1:A:166:GLY:HA3	1.95	0.46
2:A:401:R2R:N05	2:A:401:R2R:N12	2.64	0.46
1:A:130:GLY:O	1:A:133:PHE:HB3	2.15	0.46
1:A:191:LYS:N	1:A:191:LYS:CD	2.78	0.46
1:A:222:PHE:CE1	1:A:280:VAL:HG12	2.51	0.46
1:B:221:LEU:O	1:B:226:PRO:HD3	2.16	0.46
1:B:56:LEU:HD22	4:B:903:D12:H31	1.98	0.46
1:A:249:LEU:HD22	1:A:281:GLY:C	2.35	0.46
4:B:903:D12:H42	4:B:903:D12:H72	1.47	0.46
1:B:288:VAL:HA	1:B:291:MET:HE3	1.98	0.46
1:A:221:LEU:O	1:A:226:PRO:HD3	2.15	0.45
1:A:44:TRP:NE1	1:A:46:THR:OG1	2.49	0.45
1:B:225:LEU:N	1:B:226:PRO:HD2	2.31	0.45
1:A:173:LEU:HD22	1:B:54:VAL:HG21	1.98	0.45
1:B:107:VAL:O	1:B:110:ASP:HB3	2.16	0.45
1:B:197:ILE:HG13	1:B:197:ILE:O	2.17	0.45
1:B:262:SER:OG	1:B:264:ILE:HG12	2.17	0.45
1:B:110:ASP:C	1:B:113:GLY:H	2.17	0.45
1:B:188:GLY:HA2	1:B:191:LYS:HZ3	1.80	0.45
1:B:261:GLY:HA2	5:B:904:Q6F:C20	2.46	0.45
1:A:180:GLN:HG2	1:B:44:TRP:CZ3	2.52	0.45
1:B:145:PHE:HD1	1:B:255:GLY:HA2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:ILE:HG13	1:B:193:GLU:CD	2.37	0.45
1:B:40:ASN:CA	1:B:43:LYS:HE3	2.44	0.45
1:A:306:ALA:O	1:A:309:GLU:HB3	2.16	0.44
1:B:213:ILE:HG13	1:B:217:PHE:HD2	1.82	0.44
1:A:134:PHE:HA	1:A:275:TRP:HH2	1.82	0.44
1:A:220:VAL:HA	1:A:224:ALA:CB	2.47	0.44
1:B:249:LEU:HD22	1:B:281:GLY:C	2.37	0.44
1:A:127:TRP:O	1:B:66:LYS:HG3	2.17	0.44
1:B:114:ILE:O	1:B:116:PRO:HD3	2.18	0.44
1:B:46:THR:O	1:B:50:ILE:HG13	2.17	0.44
1:A:246:VAL:O	1:A:250:THR:HG23	2.18	0.44
1:B:233:ILE:HD13	1:B:270:TYR:CD1	2.52	0.44
1:A:209:ILE:HD12	1:A:209:ILE:H	1.83	0.44
1:B:50:ILE:O	1:B:54:VAL:HG23	2.18	0.44
1:A:249:LEU:HB2	1:A:285:PHE:CD2	2.53	0.43
1:B:270:TYR:O	1:B:274:VAL:HG23	2.18	0.43
1:B:160:ILE:O	1:B:164:LEU:HG	2.17	0.43
1:A:208:ILE:O	1:A:212:ILE:HG13	2.17	0.43
1:B:194:ASP:C	1:B:197:ILE:HG22	2.39	0.43
1:A:141:THR:O	1:A:142:THR:OG1	2.22	0.43
1:B:79:THR:HA	1:B:82:ILE:HD12	1.99	0.43
4:B:903:D12:H62	4:B:903:D12:H92	1.45	0.43
1:B:208:ILE:O	1:B:212:ILE:HG13	2.19	0.43
1:A:160:ILE:HG23	1:B:246:VAL:HG11	1.99	0.43
1:B:104:GLN:HA	1:B:107:VAL:HG12	2.00	0.43
1:A:41:VAL:C	1:A:43:LYS:N	2.70	0.43
1:A:79:THR:O	1:A:83:GLN:HG3	2.18	0.43
1:B:75:SER:O	1:B:79:THR:HG23	2.18	0.43
1:B:196:PHE:CD2	1:B:209:ILE:HG21	2.54	0.42
1:A:216:LEU:O	1:A:220:VAL:HG23	2.19	0.42
1:A:58:LEU:HD21	1:B:140:ILE:HD11	2.00	0.42
1:B:140:ILE:HG22	1:B:166:GLY:HA3	2.01	0.42
1:B:222:PHE:CE1	1:B:280:VAL:CG1	3.02	0.42
1:A:187:LYS:CA	1:A:191:LYS:HE3	2.49	0.42
1:B:258:VAL:HG12	1:B:260:GLY:H	1.83	0.42
1:B:88:LEU:O	1:B:92:PRO:HB3	2.19	0.42
1:B:89:ARG:O	1:B:92:PRO:HD3	2.19	0.42
1:A:277:TRP:CE3	1:A:278:ILE:HG13	2.55	0.42
1:A:249:LEU:CB	1:A:285:PHE:CD2	3.02	0.42
1:B:78:THR:O	1:B:82:ILE:HG13	2.19	0.41
1:A:207:ARG:O	1:A:211:THR:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:249:LEU:CB	1:B:285:PHE:CD2	3.03	0.41
1:B:103:ILE:HA	1:B:106:ILE:HG12	2.03	0.41
1:B:241:ALA:O	1:B:245:VAL:HG23	2.20	0.41
1:B:271:LYS:HB2	1:B:272:PRO:HD3	2.02	0.41
1:B:136:ALA:HB1	1:B:162:TYR:CE2	2.55	0.41
1:A:254:PHE:HZ	1:B:139:VAL:HG22	1.85	0.41
1:B:269:PHE:C	1:B:272:PRO:HD2	2.41	0.41
1:A:126:HIS:N	1:B:73:GLU:OE2	2.54	0.41
1:A:126:HIS:HB2	1:B:73:GLU:OE2	2.21	0.41
1:B:278:ILE:CD1	5:B:904:Q6F:C21	2.99	0.41
1:A:99:LEU:O	1:A:103:ILE:HG12	2.21	0.41
1:A:229:ILE:HG23	1:A:233:ILE:HD12	2.03	0.41
1:A:191:LYS:HE2	1:A:191:LYS:HB2	1.78	0.41
1:B:222:PHE:HE1	1:B:280:VAL:CG1	2.34	0.40
1:A:189:ILE:O	1:A:193:GLU:CG	2.70	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	257/312 (82%)	237 (92%)	17 (7%)	3 (1%)	13	39
1	B	266/312 (85%)	247 (93%)	18 (7%)	1 (0%)	34	66
All	All	523/624 (84%)	484 (92%)	35 (7%)	4 (1%)	19	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	MET
1	A	44	TRP
1	A	43	LYS

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Mol	Chain	Res	Type
1	B	264	ILE

### 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	215/260 (83%)	212 (99%)	3 (1%)	67	90
1	B	221/260 (85%)	219 (99%)	2 (1%)	78	94
All	All	436/520 (84%)	431 (99%)	5 (1%)	78	92

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	MET
1	A	43	LYS
1	A	134	PHE
1	B	134[A]	PHE
1	B	134[B]	PHE

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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
4	D12	A	405	-	11,11,11	0.30	0	10,10,10	0.83	0
2	R2R	A	401	-	14,18,18	18.06	14 (100%)	-		
4	D12	B	903	-	11,11,11	0.30	0	10,10,10	0.85	0
5	Q6F	B	904	-	24,24,24	2.06	5 (20%)	34,34,34	3.15	12 (35%)
4	D12	A	404	-	11,11,11	0.31	0	10,10,10	0.81	0
7	HEX	B	901	-	5,5,5	0.32	0	4,4,4	0.56	0
4	D12	A	403	-	11,11,11	0.32	0	10,10,10	0.79	0
5	Q6F	A	406	-	24,24,24	2.04	7 (29%)	34,34,34	3.16	13 (38%)

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
4	D12	A	405	-	-	2/9/9/9	-
4	D12	B	903	-	-	5/9/9/9	-
5	Q6F	B	904	-	-	1/14/14/14	0/2/2/2
4	D12	A	404	-	-	2/9/9/9	-
7	HEX	B	901	-	-	0/3/3/3	-
4	D12	A	403	-	-	5/9/9/9	-
5	Q6F	A	406	-	-	1/14/14/14	0/2/2/2

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	R2R	RU2-N01	-19.17	1.79	2.11
2	A	401	R2R	RU2-N04	-18.72	1.80	2.11
2	A	401	R2R	RU2-N05	-18.66	1.80	2.11
2	A	401	R2R	RU2-N03	-18.52	1.80	2.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	R2R	RU1-N19	-17.95	1.80	2.10
2	A	401	R2R	RU3-N12	-17.82	1.80	2.10
2	A	401	R2R	RU1-N16	-17.78	1.80	2.10
2	A	401	R2R	RU3-N10	-17.76	1.80	2.10
2	A	401	R2R	RU3-N11	-17.76	1.80	2.10
2	A	401	R2R	RU3-N09	-17.73	1.80	2.10
2	A	401	R2R	RU1-N15	-17.73	1.80	2.10
2	A	401	R2R	RU1-N17	-17.72	1.80	2.10
2	A	401	R2R	RU1-N18	-17.71	1.80	2.10
2	A	401	R2R	RU3-N08	-17.68	1.80	2.10
5	B	904	Q6F	S03-N08	5.94	1.71	1.63
5	A	406	Q6F	S03-N08	5.84	1.71	1.63
5	B	904	Q6F	C13-N07	5.54	1.46	1.33
5	A	406	Q6F	C13-N07	5.48	1.45	1.33
5	B	904	Q6F	C20-S03	2.72	1.81	1.75
5	A	406	Q6F	C20-S03	2.70	1.81	1.75
5	B	904	Q6F	C23-CL2	2.29	1.79	1.74
5	A	406	Q6F	C18-CL1	2.24	1.79	1.73
5	A	406	Q6F	C23-CL2	2.16	1.79	1.74
5	A	406	Q6F	O05-S03	2.07	1.47	1.43
5	A	406	Q6F	O04-C13	-2.06	1.19	1.23
5	B	904	Q6F	O04-C13	-2.04	1.19	1.23

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	406	Q6F	O06-S03-O05	-12.71	100.58	118.85
5	B	904	Q6F	O06-S03-O05	-12.61	100.71	118.85
5	A	406	Q6F	C16-C11-N08	-6.38	106.29	120.09
5	A	406	Q6F	C17-C11-N08	6.35	133.82	120.09
5	B	904	Q6F	C17-C11-N08	6.28	133.67	120.09
5	B	904	Q6F	C16-C11-N08	-6.28	106.52	120.09
5	A	406	Q6F	C21-C18-C12	-3.52	119.29	122.42
5	B	904	Q6F	C22-C23-CL2	3.36	124.60	119.35
5	A	406	Q6F	C20-S03-N08	3.32	110.42	106.63
5	B	904	Q6F	C21-C18-C12	-3.18	119.59	122.42
5	B	904	Q6F	C21-C23-CL2	-3.14	115.23	119.15
5	B	904	Q6F	C12-C09-N07	-3.11	106.47	113.03
5	A	406	Q6F	C12-C09-N07	-3.09	106.51	113.03
5	B	904	Q6F	C20-S03-N08	3.08	110.14	106.63
5	A	406	Q6F	C22-C23-CL2	2.96	123.98	119.35
5	B	904	Q6F	C19-C12-C18	2.73	121.40	116.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	406	Q6F	C09-C12-C19	-2.63	115.01	121.00
5	A	406	Q6F	C21-C23-CL2	-2.60	115.90	119.15
5	A	406	Q6F	C19-C12-C18	2.57	121.13	116.91
5	B	904	Q6F	C09-C12-C19	-2.52	115.26	121.00
5	B	904	Q6F	O06-S03-N08	2.23	111.66	107.10
5	A	406	Q6F	O06-S03-N08	2.08	111.34	107.10
5	B	904	Q6F	O05-S03-N08	2.07	111.33	107.10
5	A	406	Q6F	O05-S03-C20	2.05	111.58	108.28
5	A	406	Q6F	C09-C12-C18	2.05	123.84	121.13

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	903	D12	C4-C5-C6-C7
4	B	903	D12	C6-C7-C8-C9
4	A	403	D12	C4-C5-C6-C7
4	B	903	D12	C7-C8-C9-C10
4	B	903	D12	C3-C4-C5-C6
4	A	405	D12	C4-C5-C6-C7
4	B	903	D12	C2-C3-C4-C5
4	A	403	D12	C5-C6-C7-C8
4	A	403	D12	C1-C2-C3-C4
4	A	405	D12	C3-C4-C5-C6
4	A	403	D12	C2-C3-C4-C5
4	A	404	D12	C2-C3-C4-C5
5	B	904	Q6F	C11-N08-S03-O05
5	A	406	Q6F	C11-N08-S03-O05
4	A	403	D12	C3-C4-C5-C6
4	A	404	D12	C6-C7-C8-C9

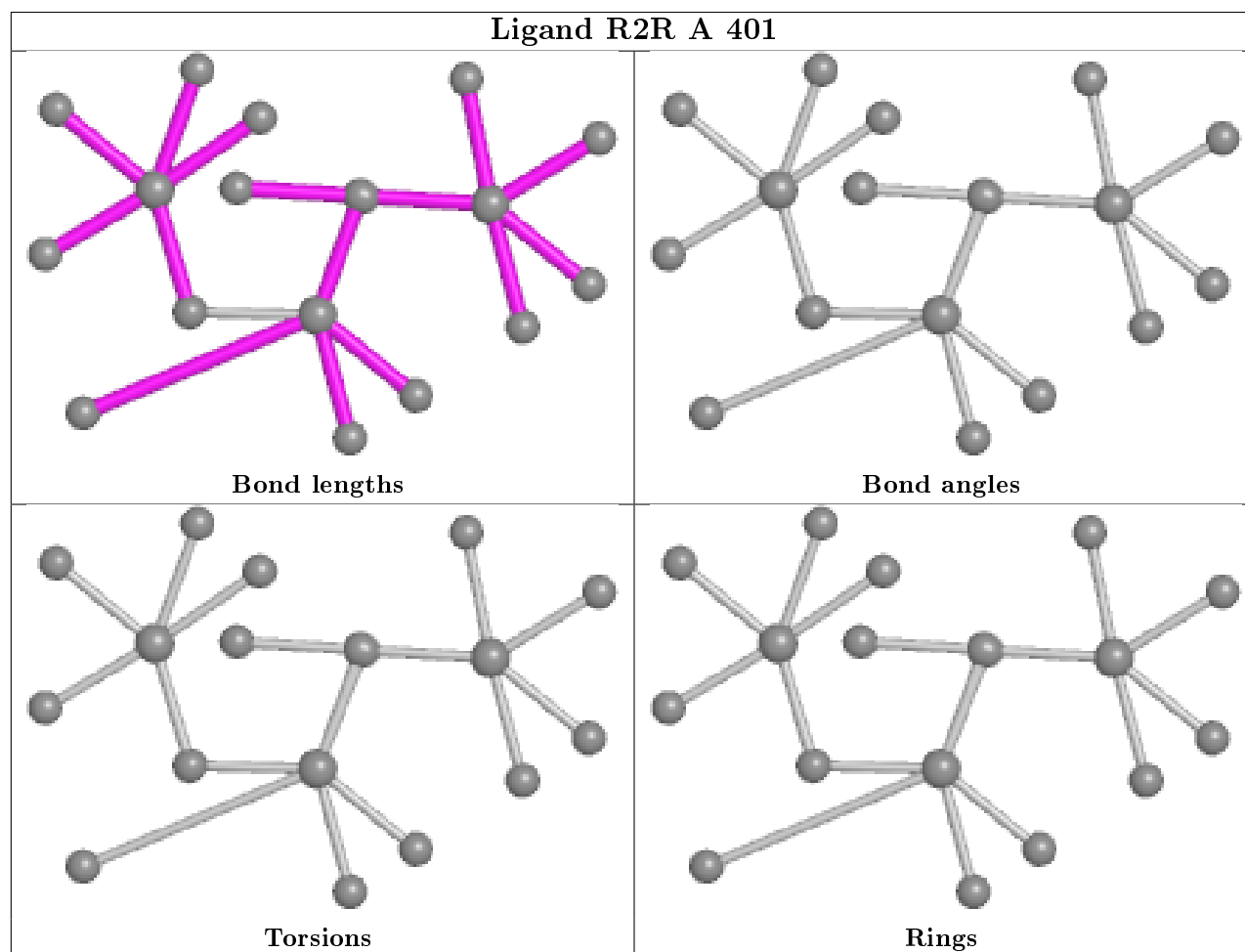
There are no ring outliers.

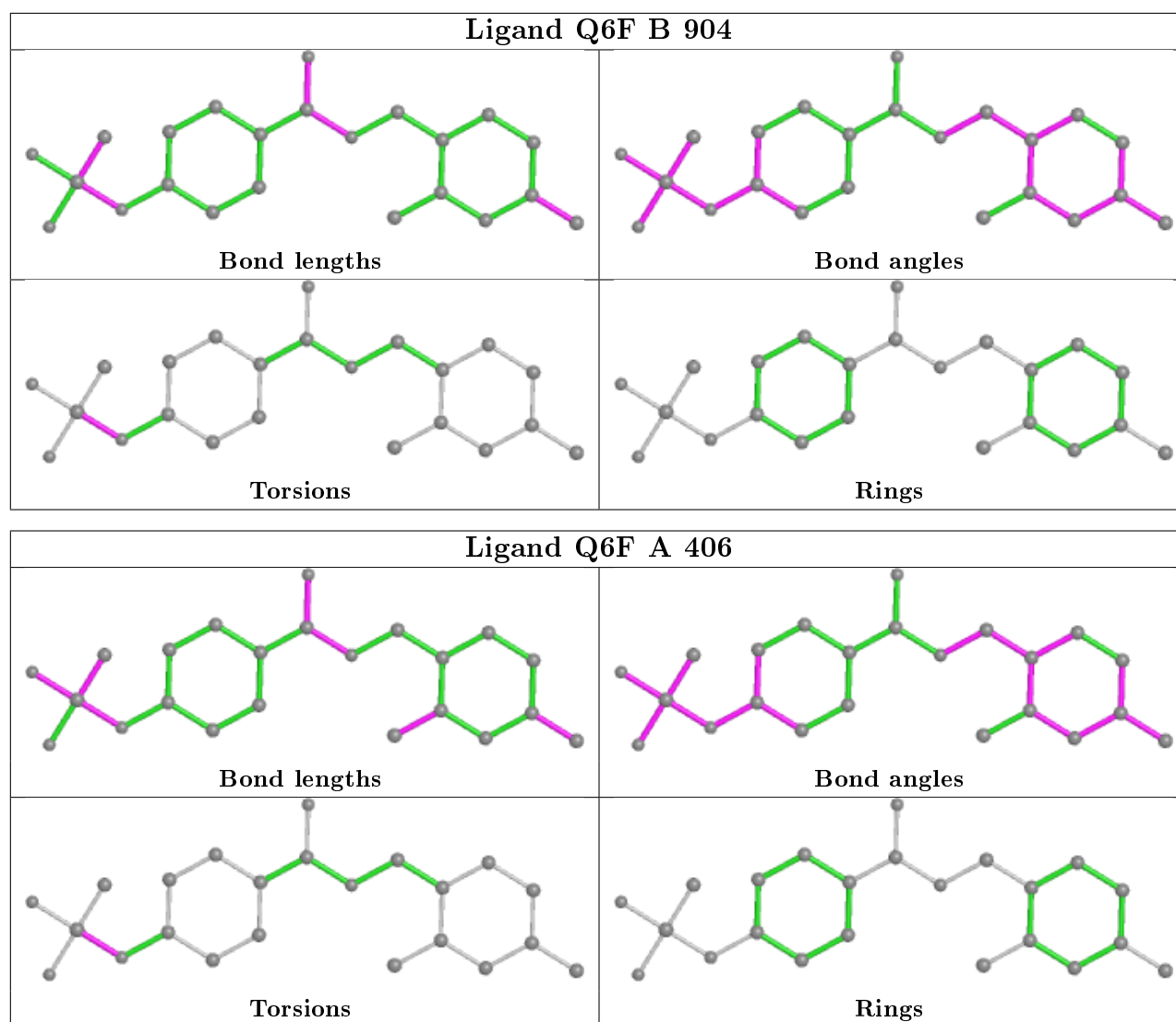
4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	R2R	4	0
4	B	903	D12	4	0
5	B	904	Q6F	3	0
5	A	406	Q6F	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	263/312 (84%)	1.04	53 (20%) 1 0	88, 137, 202, 230	0
1	B	271/312 (86%)	0.99	46 (16%) 1 1	89, 136, 213, 238	0
All	All	534/624 (85%)	1.01	99 (18%) 1 1	88, 137, 209, 238	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	196	PHE	13.1
1	A	85	GLU	8.9
1	B	192	VAL	7.9
1	B	195	THR	7.2
1	B	317	TRP	5.9
1	A	82	ILE	5.7
1	B	206	ILE	5.6
1	A	155	GLY	5.5
1	B	87	PHE	5.3
1	A	154	GLY	5.3
1	B	199	TRP	5.0
1	A	161	ILE	5.0
1	A	206	ILE	5.0
1	B	209	ILE	4.7
1	B	69	GLU	4.5
1	A	311	ARG	4.5
1	B	204	THR	4.4
1	B	185	PHE	4.4
1	A	158	PHE	4.4
1	B	205	LYS	4.3
1	B	193	GLU	4.3
1	A	263	ASP	4.1
1	A	296	LEU	4.0
1	B	265	GLU	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	153	GLU	3.9
1	A	96	ASP	3.8
1	A	81	VAL	3.8
1	A	156	LYS	3.7
1	B	89	ARG	3.7
1	A	187	LYS	3.7
1	B	70	GLN	3.6
1	A	193	GLU	3.6
1	A	94	VAL	3.5
1	A	42	MET	3.5
1	B	85	GLU	3.5
1	A	196	PHE	3.5
1	A	316	GLU	3.4
1	A	266	TYR	3.4
1	A	303	THR	3.3
1	B	198	LYS	3.3
1	B	261	GLY	3.2
1	A	157	ILE	3.2
1	B	315	ALA	3.2
1	A	41	VAL	3.2
1	A	102	LEU	3.2
1	A	141	THR	3.1
1	B	203	GLN	3.1
1	B	100	ASP	3.0
1	A	291	MET	3.0
1	B	77	ARG	3.0
1	B	305	GLU	3.0
1	A	312	ALA	2.9
1	A	315	ALA	2.9
1	B	95	SER	2.8
1	A	308	GLY	2.8
1	A	268	ASP	2.7
1	B	314	ALA	2.7
1	A	307	VAL	2.7
1	A	310	PHE	2.6
1	B	188	GLY	2.6
1	A	189	ILE	2.6
1	B	80	ILE	2.6
1	B	83	GLN	2.6
1	A	87	PHE	2.5
1	A	171	GLY	2.5
1	A	138	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	81	VAL	2.5
1	A	264	ILE	2.5
1	A	185	PHE	2.5
1	B	37	SER	2.5
1	B	239	LEU	2.5
1	A	220	VAL	2.5
1	B	74	ILE	2.4
1	B	65	PHE	2.4
1	A	88	LEU	2.4
1	A	160	ILE	2.4
1	B	157	ILE	2.4
1	A	40	ASN	2.3
1	B	316	GLU	2.3
1	B	266	TYR	2.3
1	A	278	ILE	2.3
1	A	170	PHE	2.3
1	A	205	LYS	2.2
1	B	191	LYS	2.2
1	B	240	ASP	2.2
1	A	143	ILE	2.2
1	B	187	LYS	2.2
1	A	295	TRP	2.1
1	B	312	ALA	2.1
1	B	212	ILE	2.1
1	A	217	PHE	2.1
1	B	282	LEU	2.1
1	A	209	ILE	2.1
1	B	283	ALA	2.1
1	B	302	LYS	2.0
1	A	176	GLY	2.0
1	A	152	THR	2.0
1	B	251	THR	2.0
1	A	127	TRP	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

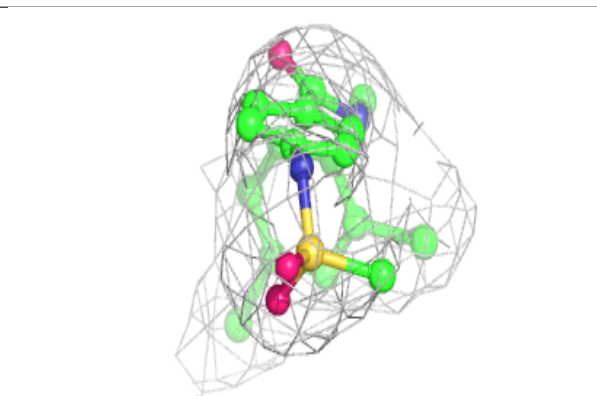
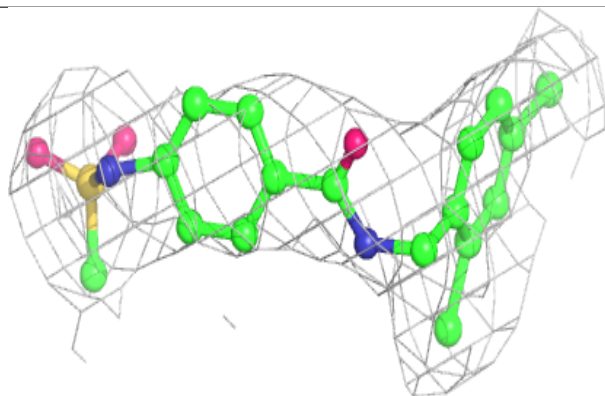
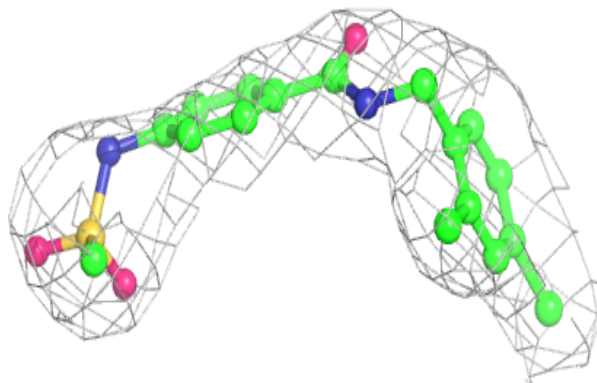
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	CD	B	906	1/1	0.39	1.22	360,360,360,360	0
3	CD	A	402	1/1	0.74	0.24	329,329,329,329	0
4	D12	A	405	12/12	0.74	0.52	111,170,190,192	0
7	HEX	B	901	6/6	0.79	0.23	127,160,182,182	0
4	D12	A	403	12/12	0.80	0.23	99,127,145,152	0
4	D12	A	404	12/12	0.84	0.21	111,139,147,148	0
6	K	A	410	1/1	0.85	0.12	107,107,107,107	0
5	Q6F	B	904	23/23	0.86	0.25	105,131,151,152	0
6	K	A	408	1/1	0.87	0.18	106,106,106,106	0
6	K	A	409	1/1	0.88	0.80	127,127,127,127	0
6	K	B	905	1/1	0.89	0.25	106,106,106,106	0
4	D12	B	903	12/12	0.89	0.29	96,116,133,136	0
6	K	A	407	1/1	0.90	0.07	128,128,128,128	0
5	Q6F	A	406	23/23	0.91	0.22	118,128,150,160	0
3	CD	B	902	1/1	0.94	0.15	172,172,172,172	0
2	R2R	A	401	19/19	0.97	0.14	208,228,244,244	0

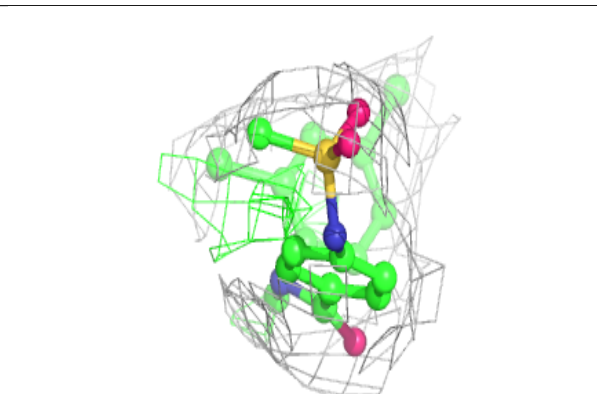
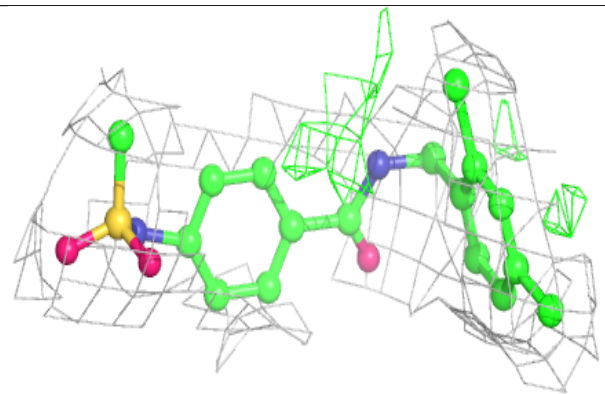
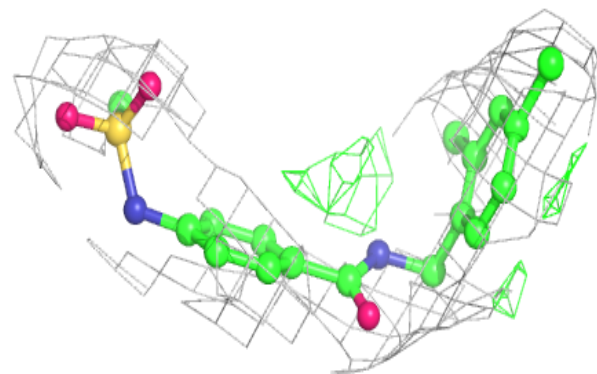
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around Q6F B 904:**

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

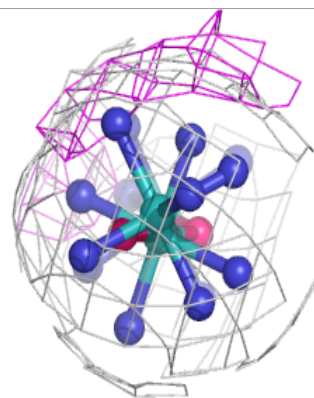
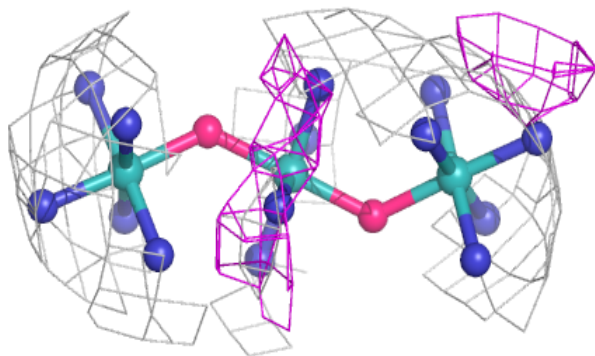
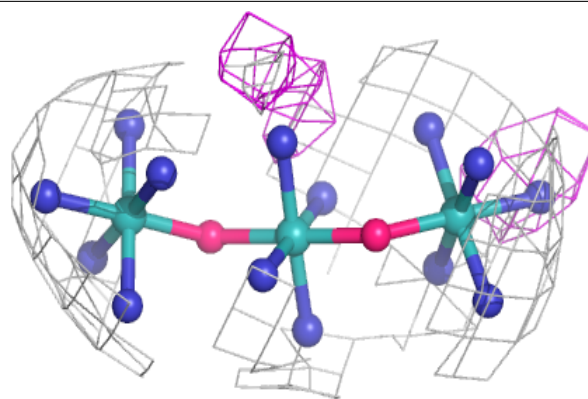
**Electron density around Q6F A 406:**

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



**Electron density around R2R A 401:**

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



## 6.5 Other polymers [i](#)

There are no such residues in this entry.