



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 12:26 am GMT

PDB ID : 3CTH
Title : Crystal structure of the tyrosine kinase domain of the hepatocyte growth factor receptor c-met in complex with a aminopyridine based inhibitor
Authors : Sack, J.
Deposited on : 2008-04-14
Resolution : 2.30 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 : trunk28620
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) : recalc28949

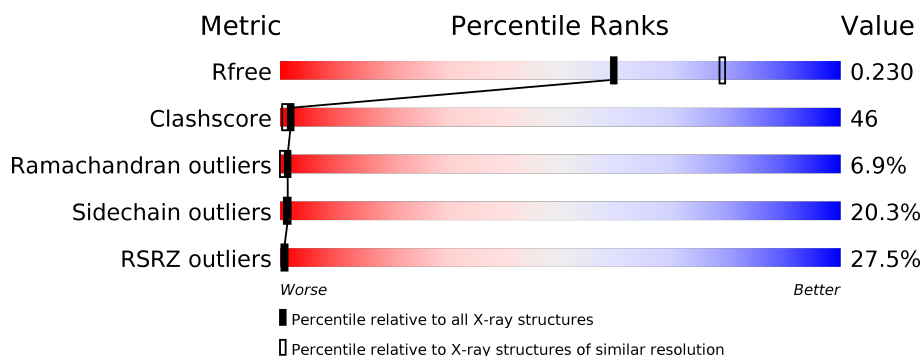
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	314	<div> <div>26%</div> <div>49%</div> <div>28%</div> <div>15%</div> <div>6%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2461 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.

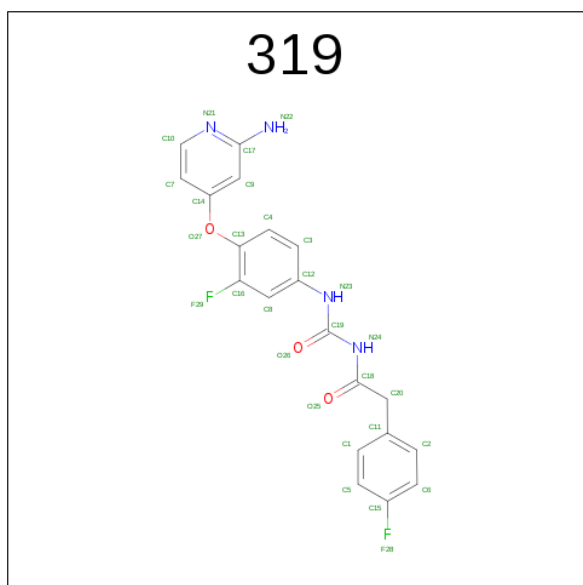
- Molecule 1 is a protein called Hepatocyte growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	295	2341	1511	402	414	14	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1047	GLY	-	EXPRESSION TAG	UNP P08581
A	1048	ALA	-	EXPRESSION TAG	UNP P08581
A	1194	PHE	TYR	ENGINEERED	UNP P08581
A	1234	PHE	TYR	ENGINEERED	UNP P08581
A	1235	ASP	TYR	ENGINEERED	UNP P08581

- Molecule 2 is N-({4-[(2-AMINOPYRIDIN-4-YL)OXY]-3-FLUOROPHENYL}CARBAMOYL)-2-(4-FLUOROPHENYL)ACETAMIDE (three-letter code: 319) (formula: C₂₀H₁₆F₂N₄O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			29	20	2	4	3		

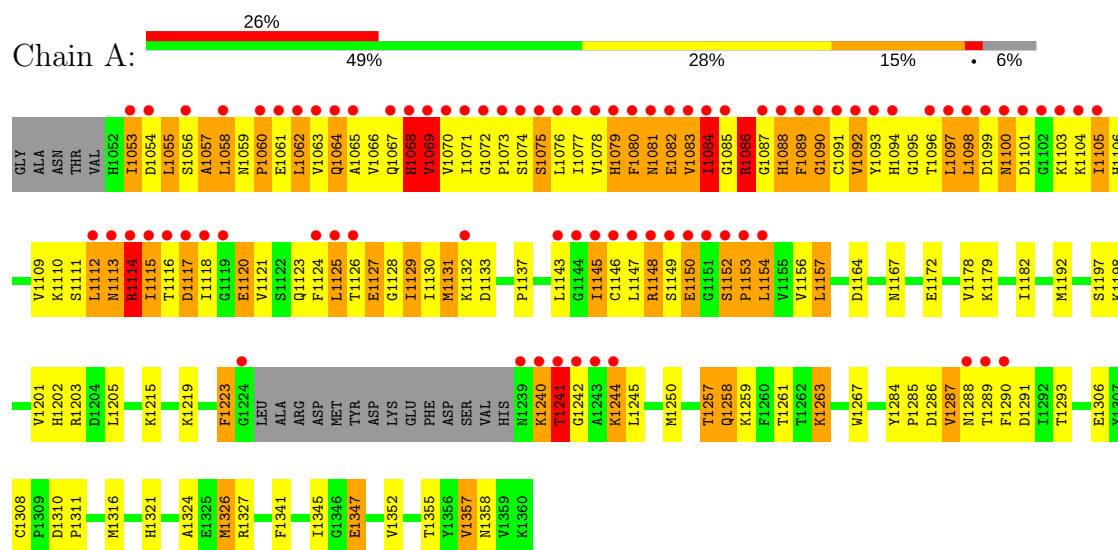
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	91	Total	O	0	0
			91	91		

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 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: Hepatocyte growth factor receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	42.75Å 47.64Å 152.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.25 – 2.30 34.76 – 2.31	Depositor EDS
% Data completeness (in resolution range)	87.3 (15.25-2.30) 87.4 (34.76-2.31)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.82 (at 2.31Å)	Xtriage
Refinement program	BUSTER-TNT 2.1.1	Depositor
R, R_{free}	0.207 , 0.273 0.218 , 0.230	Depositor DCC
R_{free} test set	762 reflections (6.44%)	DCC
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.445	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 86.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2461	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.65% 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: 319

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	0.41	0/2399	0.58	0/3249

There are no bond length outliers.

There are no bond angle outliers.

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	2341	0	2366	217	0
2	A	29	0	16	5	0
3	A	91	0	0	3	0
All	All	2461	0	2382	218	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 46.

All (218) 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:1147:LEU:HD11	1:A:1154:LEU:H	1.09	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1085:GLY:HA3	1:A:1092:VAL:HG23	1.29	1.14
1:A:1058:LEU:HB2	1:A:1063:VAL:HG11	1.13	1.10
1:A:1099:ASP:HB3	1:A:1103:LYS:HB2	1.32	1.10
1:A:1086:ARG:HH11	1:A:1086:ARG:HA	1.18	1.08
1:A:1076:LEU:HD13	1:A:1097:LEU:HD12	1.35	1.02
1:A:1076:LEU:HD22	1:A:1097:LEU:HD11	1.43	0.98
1:A:1092:VAL:HG13	1:A:1110:LYS:HB3	1.47	0.96
1:A:1086:ARG:NH1	1:A:1086:ARG:HA	1.83	0.94
1:A:1109:VAL:HG12	1:A:1156:VAL:HG22	1.52	0.92
1:A:1053:ILE:HG13	1:A:1055:LEU:HD13	1.55	0.88
1:A:1076:LEU:HB2	1:A:1097:LEU:HG	1.52	0.88
1:A:1053:ILE:HD12	1:A:1054:ASP:N	1.90	0.85
1:A:1085:GLY:HA3	1:A:1092:VAL:CG2	2.06	0.85
1:A:1091:CYS:HG	1:A:1093:TYR:HE1	1.25	0.84
1:A:1147:LEU:HD11	1:A:1154:LEU:N	1.90	0.84
1:A:1066:VAL:HG21	1:A:1125:LEU:HD11	1.60	0.84
1:A:1147:LEU:CD1	1:A:1154:LEU:H	1.91	0.83
1:A:1058:LEU:HB2	1:A:1063:VAL:CG1	2.04	0.82
1:A:1079:HIS:CE1	1:A:1095:GLY:HA2	2.15	0.82
1:A:1117:ASP:O	1:A:1118:ILE:HD12	1.82	0.80
1:A:1067:GLN:O	1:A:1070:VAL:HG12	1.81	0.79
1:A:1145:ILE:HD12	1:A:1146:CYS:N	1.97	0.79
1:A:1069:VAL:HG12	1:A:1145:ILE:HG22	1.65	0.79
1:A:1085:GLY:CA	1:A:1092:VAL:HG23	2.13	0.79
1:A:1069:VAL:HG12	1:A:1145:ILE:CG2	2.13	0.78
1:A:1092:VAL:HA	1:A:1109:VAL:O	1.84	0.77
1:A:1289:THR:HG23	1:A:1291:ASP:H	1.49	0.77
1:A:1257:THR:OG1	1:A:1259:LYS:HG2	1.84	0.77
1:A:1053:ILE:HG13	1:A:1055:LEU:CD1	2.14	0.77
1:A:1120:GLU:HA	1:A:1123:GLN:NE2	2.02	0.75
1:A:1145:ILE:HD12	1:A:1146:CYS:H	1.49	0.74
1:A:1076:LEU:HD22	1:A:1097:LEU:CD1	2.18	0.72
1:A:1059:ASN:HB2	1:A:1063:VAL:HG12	1.71	0.71
1:A:1131:MET:SD	2:A:2001:319:H2O	2.29	0.71
1:A:1118:ILE:HA	1:A:1121:VAL:HB	1.72	0.71
1:A:1240:LYS:HB3	1:A:1258:GLN:HB3	1.71	0.71
1:A:1064:GLN:O	1:A:1067:GLN:HG3	1.89	0.71
1:A:1076:LEU:HD13	1:A:1097:LEU:CD1	2.19	0.71
1:A:1058:LEU:HD12	1:A:1063:VAL:HG13	1.71	0.70
1:A:1152:SER:CB	1:A:1153:PRO:HD2	2.20	0.70
1:A:1289:THR:HG23	1:A:1291:ASP:N	2.06	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1079:HIS:ND1	1:A:1095:GLY:HA2	2.06	0.69
1:A:1112:LEU:HD11	1:A:1124:PHE:CD1	2.27	0.69
1:A:1240:LYS:O	1:A:1258:GLN:HG2	1.93	0.69
1:A:1126:THR:HA	1:A:1129:ILE:CG2	2.23	0.69
1:A:1096:THR:HA	1:A:1105:ILE:O	1.93	0.68
1:A:1066:VAL:HG13	1:A:1145:ILE:HG21	1.76	0.68
1:A:1055:LEU:O	1:A:1057:ALA:N	2.28	0.67
1:A:1059:ASN:HB2	1:A:1063:VAL:CG1	2.25	0.67
1:A:1201:VAL:HG12	1:A:1203:ARG:HG3	1.77	0.66
1:A:1070:VAL:HG23	1:A:1145:ILE:O	1.96	0.65
1:A:1289:THR:HG23	1:A:1290:PHE:N	2.11	0.65
1:A:1067:GLN:O	1:A:1069:VAL:N	2.29	0.65
1:A:1066:VAL:HG22	1:A:1125:LEU:HD21	1.79	0.65
1:A:1075:SER:O	1:A:1098:LEU:N	2.30	0.65
1:A:1059:ASN:HB3	1:A:1062:LEU:CB	2.27	0.64
1:A:1259:LYS:HG3	1:A:1259:LYS:O	1.96	0.64
1:A:1284:TYR:HB3	1:A:1287:VAL:HG13	1.80	0.64
1:A:1109:VAL:HB	1:A:1154:LEU:HD21	1.80	0.63
1:A:1257:THR:O	1:A:1259:LYS:N	2.28	0.63
1:A:1152:SER:OG	1:A:1153:PRO:HD2	1.98	0.63
1:A:1066:VAL:CG2	1:A:1125:LEU:HD21	2.29	0.63
1:A:1053:ILE:HD12	1:A:1054:ASP:H	1.63	0.63
1:A:1093:TYR:HB2	1:A:1109:VAL:CG2	2.29	0.63
1:A:1059:ASN:N	1:A:1063:VAL:HG12	2.15	0.62
1:A:1065:ALA:HA	1:A:1068:HIS:NE2	2.15	0.61
1:A:1127:GLU:HG3	1:A:1128:GLY:N	2.15	0.61
1:A:1289:THR:HG23	1:A:1290:PHE:H	1.66	0.61
1:A:1090:GLY:HA2	1:A:1111:SER:O	2.02	0.60
1:A:1128:GLY:O	1:A:1131:MET:HG3	2.02	0.60
1:A:1242:GLY:HA3	3:A:91:HOH:O	2.03	0.59
1:A:1068:HIS:O	1:A:1069:VAL:HG23	2.03	0.59
1:A:1053:ILE:HG23	1:A:1055:LEU:HD13	1.85	0.59
1:A:1077:ILE:HG22	1:A:1078:VAL:H	1.66	0.59
1:A:1240:LYS:C	1:A:1258:GLN:HG2	2.23	0.59
1:A:1146:CYS:O	1:A:1147:LEU:HD12	2.04	0.58
1:A:1192:MET:CE	1:A:1205:LEU:HD22	2.33	0.58
1:A:1112:LEU:HD21	1:A:1124:PHE:CE1	2.38	0.58
1:A:1070:VAL:HA	1:A:1145:ILE:O	2.03	0.58
1:A:1125:LEU:O	1:A:1129:ILE:HG22	2.03	0.58
1:A:1059:ASN:CB	1:A:1063:VAL:HG12	2.34	0.58
1:A:1070:VAL:HA	1:A:1145:ILE:CG2	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1240:LYS:CB	1:A:1258:GLN:HB3	2.34	0.56
1:A:1287:VAL:O	1:A:1287:VAL:HG22	2.06	0.56
1:A:1066:VAL:O	1:A:1070:VAL:HB	2.06	0.56
1:A:1059:ASN:HB3	1:A:1062:LEU:HB2	1.87	0.55
1:A:1202:HIS:O	1:A:1203:ARG:HB2	2.07	0.55
1:A:1076:LEU:CD1	1:A:1097:LEU:HD12	2.22	0.55
1:A:1197:SER:OG	1:A:1198:LYS:HE3	2.07	0.55
1:A:1070:VAL:HG22	1:A:1071:ILE:N	2.21	0.54
1:A:1097:LEU:O	1:A:1098:LEU:HD13	2.07	0.54
1:A:1077:ILE:HG22	1:A:1078:VAL:N	2.23	0.54
1:A:1069:VAL:HG12	1:A:1145:ILE:HG21	1.89	0.54
1:A:1080:PHE:CD1	1:A:1080:PHE:N	2.76	0.54
1:A:1124:PHE:CD2	1:A:1153:PRO:HB2	2.43	0.54
1:A:1126:THR:HA	1:A:1129:ILE:HG22	1.90	0.54
1:A:1070:VAL:HA	1:A:1145:ILE:HG23	1.90	0.54
1:A:1083:VAL:HG12	1:A:1093:TYR:HD1	1.73	0.53
1:A:1060:PRO:O	1:A:1064:GLN:HG2	2.09	0.53
1:A:1080:PHE:CD2	1:A:1093:TYR:CE2	2.96	0.53
1:A:1147:LEU:HD21	1:A:1153:PRO:HA	1.89	0.53
1:A:1111:SER:O	1:A:1112:LEU:HD23	2.08	0.53
1:A:1105:ILE:HG13	1:A:1106:HIS:N	2.24	0.52
1:A:1357:VAL:O	1:A:1358:ASN:HB2	2.09	0.52
1:A:1059:ASN:HB2	1:A:1063:VAL:CB	2.40	0.52
1:A:1091:CYS:SG	1:A:1093:TYR:HE1	2.32	0.52
1:A:1132:LYS:HG3	1:A:1133:ASP:OD2	2.10	0.52
1:A:1104:LYS:C	1:A:1105:ILE:HG22	2.29	0.52
1:A:1310:ASP:N	1:A:1311:PRO:HD2	2.23	0.52
1:A:1066:VAL:HG21	1:A:1125:LEU:CD1	2.35	0.52
1:A:1059:ASN:CA	1:A:1063:VAL:HG12	2.40	0.52
1:A:1063:VAL:C	1:A:1065:ALA:H	2.11	0.52
1:A:1075:SER:O	1:A:1097:LEU:HA	2.10	0.52
1:A:1126:THR:HA	1:A:1129:ILE:HG21	1.92	0.52
1:A:1070:VAL:C	1:A:1071:ILE:HG12	2.30	0.52
1:A:1088:HIS:CE1	1:A:1089:PHE:CD2	2.98	0.51
1:A:1066:VAL:HG13	1:A:1145:ILE:CG2	2.41	0.51
1:A:1065:ALA:HA	1:A:1068:HIS:CD2	2.46	0.51
1:A:1116:THR:HG23	1:A:1116:THR:O	2.11	0.51
1:A:1092:VAL:HG13	1:A:1110:LYS:CB	2.32	0.51
1:A:1089:PHE:CD1	1:A:1112:LEU:HB3	2.46	0.51
1:A:1070:VAL:O	1:A:1071:ILE:HG12	2.11	0.51
1:A:1120:GLU:OE2	1:A:1121:VAL:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1094:HIS:HD1	1:A:1106:HIS:CE1	2.28	0.50
1:A:1131:MET:HE3	2:A:2001:319:H2	1.93	0.50
1:A:1112:LEU:O	1:A:1113:ASN:O	2.29	0.50
1:A:1058:LEU:CD1	1:A:1063:VAL:HG13	2.42	0.49
1:A:1099:ASP:HB3	1:A:1103:LYS:CB	2.23	0.49
1:A:1114:ARG:HG3	1:A:1152:SER:HB2	1.93	0.49
1:A:1093:TYR:HB2	1:A:1109:VAL:HG23	1.93	0.49
1:A:1157:LEU:HD13	2:A:2001:319:C8	2.41	0.49
1:A:1112:LEU:HG	1:A:1124:PHE:CE1	2.48	0.49
1:A:1058:LEU:HD12	1:A:1063:VAL:CG1	2.39	0.49
1:A:1179:LYS:HE3	1:A:1341:PHE:CZ	2.47	0.49
1:A:1324:ALA:O	1:A:1327:ARG:HB2	2.13	0.49
1:A:1059:ASN:HB2	1:A:1063:VAL:HB	1.94	0.49
1:A:1111:SER:HA	1:A:1124:PHE:CZ	2.47	0.49
1:A:1063:VAL:O	1:A:1063:VAL:HG22	2.13	0.49
1:A:1080:PHE:CD2	1:A:1093:TYR:HE2	2.30	0.49
1:A:1086:ARG:NH1	1:A:1091:CYS:HA	2.29	0.48
1:A:1263:LYS:HG3	1:A:1324:ALA:O	2.14	0.48
1:A:1053:ILE:HG23	1:A:1055:LEU:CD1	2.43	0.48
1:A:1081:ASN:HD22	1:A:1082:GLU:CD	2.17	0.48
1:A:1114:ARG:C	1:A:1116:THR:H	2.17	0.48
1:A:1244:LYS:HB2	3:A:91:HOH:O	2.13	0.48
1:A:1083:VAL:O	1:A:1084:ILE:HG23	2.14	0.48
1:A:1137:PRO:O	1:A:1219:LYS:HD3	2.14	0.48
1:A:1086:ARG:CZ	1:A:1091:CYS:HA	2.44	0.47
1:A:1112:LEU:HD21	1:A:1124:PHE:HE1	1.79	0.47
1:A:1080:PHE:HD1	1:A:1080:PHE:H	1.60	0.47
1:A:1127:GLU:CG	1:A:1128:GLY:N	2.76	0.47
1:A:1117:ASP:C	1:A:1118:ILE:HD12	2.35	0.47
1:A:1192:MET:HE2	1:A:1205:LEU:CD2	2.45	0.47
1:A:1241:THR:HG23	1:A:1241:THR:O	2.14	0.46
1:A:1053:ILE:HD12	1:A:1053:ILE:C	2.33	0.46
1:A:1089:PHE:HB3	1:A:1112:LEU:HD22	1.97	0.46
1:A:1126:THR:CA	1:A:1129:ILE:HG22	2.46	0.46
1:A:1321:HIS:CD2	1:A:1326:MET:HB3	2.51	0.46
1:A:1112:LEU:CD1	1:A:1124:PHE:CD1	2.98	0.45
1:A:1059:ASN:N	1:A:1063:VAL:CG1	2.79	0.45
1:A:1178:VAL:HG12	1:A:1182:ILE:HD12	1.98	0.45
1:A:1068:HIS:CD2	1:A:1068:HIS:N	2.84	0.45
1:A:1070:VAL:CG2	1:A:1071:ILE:N	2.79	0.45
1:A:1127:GLU:HG3	1:A:1128:GLY:H	1.78	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1257:THR:HG1	1:A:1259:LYS:HG2	1.81	0.45
1:A:1068:HIS:CD2	1:A:1068:HIS:H	2.34	0.45
1:A:1101:ASP:O	1:A:1101:ASP:OD1	2.34	0.45
1:A:1081:ASN:ND2	1:A:1082:GLU:OE2	2.50	0.45
1:A:1112:LEU:CD2	1:A:1124:PHE:CE1	3.00	0.45
1:A:1223:PHE:CE2	2:A:2001:319:C4	2.99	0.44
1:A:1059:ASN:HB3	1:A:1062:LEU:HB3	1.98	0.44
1:A:1072:GLY:C	1:A:1074:SER:H	2.21	0.44
1:A:1124:PHE:CE2	1:A:1153:PRO:HB2	2.52	0.44
1:A:1285:PRO:HG2	1:A:1286:ASP:H	1.81	0.44
1:A:1111:SER:C	1:A:1112:LEU:HD23	2.38	0.44
1:A:1114:ARG:O	1:A:1116:THR:N	2.51	0.44
1:A:1087:GLY:O	1:A:1088:HIS:HB2	2.18	0.44
1:A:1069:VAL:C	1:A:1145:ILE:HG22	2.37	0.44
1:A:1109:VAL:HG12	1:A:1156:VAL:CG2	2.36	0.43
1:A:1120:GLU:HA	1:A:1123:GLN:CD	2.38	0.43
1:A:1130:ILE:HA	1:A:1130:ILE:HD12	1.89	0.43
1:A:1250:MET:HA	3:A:50:HOH:O	2.18	0.43
1:A:1083:VAL:HG12	1:A:1093:TYR:CD1	2.53	0.43
1:A:1257:THR:O	1:A:1257:THR:OG1	2.34	0.43
1:A:1064:GLN:HG3	1:A:1064:GLN:H	1.45	0.43
1:A:1076:LEU:HA	1:A:1096:THR:O	2.19	0.43
1:A:1126:THR:C	1:A:1129:ILE:HG22	2.39	0.43
1:A:1259:LYS:NZ	1:A:1261:THR:HG21	2.34	0.43
1:A:1080:PHE:HD2	1:A:1093:TYR:HE2	1.64	0.43
1:A:1090:GLY:H	1:A:1112:LEU:CD2	2.32	0.43
1:A:1259:LYS:HE2	1:A:1261:THR:HG23	2.01	0.42
1:A:1164:ASP:OD2	1:A:1167:ASN:ND2	2.53	0.42
1:A:1267:TRP:CD1	1:A:1267:TRP:C	2.93	0.42
1:A:1345:ILE:HG12	1:A:1347:GLU:OE2	2.19	0.42
1:A:1070:VAL:HA	1:A:1145:ILE:HG22	2.01	0.42
1:A:1111:SER:HA	1:A:1124:PHE:HZ	1.83	0.42
1:A:1192:MET:HE2	1:A:1205:LEU:HD22	2.02	0.42
1:A:1127:GLU:O	1:A:1130:ILE:HG22	2.20	0.41
2:A:2001:319:H8	2:A:2001:319:O26	2.19	0.41
1:A:1058:LEU:CD1	1:A:1063:VAL:CG1	2.98	0.41
1:A:1096:THR:HG23	1:A:1105:ILE:O	2.21	0.41
1:A:1143:LEU:HA	1:A:1143:LEU:HD23	1.80	0.41
1:A:1058:LEU:HG	1:A:1058:LEU:H	1.62	0.41
1:A:1112:LEU:CG	1:A:1124:PHE:CE1	3.04	0.41
1:A:1053:ILE:CG2	1:A:1055:LEU:CD1	2.99	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1284:TYR:HB3	1:A:1287:VAL:CG1	2.49	0.41
1:A:1308:CYS:SG	1:A:1316:MET:CE	3.09	0.41
1:A:1065:ALA:O	1:A:1069:VAL:HB	2.21	0.41
1:A:1073:PRO:C	1:A:1075:SER:H	2.25	0.41
1:A:1114:ARG:HB3	1:A:1115:ILE:H	1.61	0.41
1:A:1069:VAL:O	1:A:1145:ILE:HG22	2.20	0.41
1:A:1192:MET:HG3	1:A:1205:LEU:HD21	2.03	0.41
1:A:1078:VAL:CG1	1:A:1080:PHE:CZ	3.04	0.40
1:A:1066:VAL:O	1:A:1066:VAL:HG12	2.21	0.40
1:A:1071:ILE:O	1:A:1146:CYS:SG	2.77	0.40
1:A:1120:GLU:CA	1:A:1123:GLN:NE2	2.78	0.40
1:A:1240:LYS:CB	1:A:1258:GLN:CB	3.00	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	291/314 (93%)	239 (82%)	32 (11%)	20 (7%)	1 0

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1056	SER
1	A	1060	PRO
1	A	1084	ILE
1	A	1088	HIS
1	A	1113	ASN
1	A	1150	GLU
1	A	1258	GLN
1	A	1057	ALA
1	A	1069	VAL

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Mol	Chain	Res	Type
1	A	1100	ASN
1	A	1115	ILE
1	A	1241	THR
1	A	1068	HIS
1	A	1086	ARG
1	A	1090	GLY
1	A	1114	ARG
1	A	1153	PRO
1	A	1223	PHE
1	A	1148	ARG
1	A	1288	ASN

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	261/277 (94%)	208 (80%)	53 (20%)	1 1

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1053	ILE
1	A	1055	LEU
1	A	1058	LEU
1	A	1061	GLU
1	A	1062	LEU
1	A	1064	GLN
1	A	1068	HIS
1	A	1069	VAL
1	A	1075	SER
1	A	1079	HIS
1	A	1080	PHE
1	A	1081	ASN
1	A	1082	GLU
1	A	1083	VAL
1	A	1084	ILE

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Mol	Chain	Res	Type
1	A	1086	ARG
1	A	1089	PHE
1	A	1092	VAL
1	A	1097	LEU
1	A	1098	LEU
1	A	1100	ASN
1	A	1105	ILE
1	A	1112	LEU
1	A	1114	ARG
1	A	1117	ASP
1	A	1120	GLU
1	A	1125	LEU
1	A	1127	GLU
1	A	1129	ILE
1	A	1131	MET
1	A	1145	ILE
1	A	1148	ARG
1	A	1149	SER
1	A	1150	GLU
1	A	1152	SER
1	A	1154	LEU
1	A	1157	LEU
1	A	1172	GLU
1	A	1215	LYS
1	A	1240	LYS
1	A	1241	THR
1	A	1244	LYS
1	A	1245	LEU
1	A	1257	THR
1	A	1263	LYS
1	A	1287	VAL
1	A	1293	THR
1	A	1306	GLU
1	A	1326	MET
1	A	1347	GLU
1	A	1352	VAL
1	A	1355	THR
1	A	1357	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1064	GLN
1	A	1068	HIS
1	A	1081	ASN
1	A	1088	HIS
1	A	1123	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](#)

1 ligand is 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	$\# Z > 2$
2	319	A	2001	-	31,31,31	0.85	1 (3%)	41,42,42	1.10	3 (7%)

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	319	A	2001	-	-	0/16/16/16	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2001	319	C17-N22	2.14	1.42	1.35

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	319	C20-C18-N24	-3.42	109.91	114.34
2	A	2001	319	F29-C16-C13	2.07	121.20	118.11
2	A	2001	319	C12-C8-C16	2.23	120.38	118.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	319	5	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 ⓘ

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	295/314 (93%)	1.37	81 (27%) 1 1	17, 38, 123, 136	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1062	LEU	10.6
1	A	1060	PRO	10.3
1	A	1097	LEU	9.6
1	A	1116	THR	9.5
1	A	1074	SER	9.0
1	A	1113	ASN	8.5
1	A	1073	PRO	8.4
1	A	1077	ILE	8.2
1	A	1070	VAL	8.1
1	A	1104	LYS	7.6
1	A	1069	VAL	7.5
1	A	1239	ASN	7.4
1	A	1241	THR	7.4
1	A	1091	CYS	7.1
1	A	1076	LEU	6.9
1	A	1080	PHE	6.9
1	A	1101	ASP	6.1
1	A	1243	ALA	6.0
1	A	1240	LYS	5.9
1	A	1071	ILE	5.8
1	A	1118	ILE	5.8
1	A	1088	HIS	5.7
1	A	1067	GLN	5.7
1	A	1242	GLY	5.6
1	A	1075	SER	5.5
1	A	1125	LEU	5.4
1	A	1147	LEU	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	1092	VAL	5.2
1	A	1061	GLU	5.1
1	A	1288	ASN	4.9
1	A	1089	PHE	4.9
1	A	1105	ILE	4.9
1	A	1098	LEU	4.8
1	A	1100	ASN	4.7
1	A	1114	ARG	4.6
1	A	1083	VAL	4.5
1	A	1099	ASP	4.5
1	A	1117	ASP	4.4
1	A	1103	LYS	4.4
1	A	1112	LEU	4.4
1	A	1290	PHE	4.3
1	A	1068	HIS	4.2
1	A	1093	TYR	4.1
1	A	1064	GLN	4.1
1	A	1126	THR	4.0
1	A	1082	GLU	4.0
1	A	1145	ILE	3.9
1	A	1079	HIS	3.9
1	A	1081	ASN	3.9
1	A	1149	SER	3.8
1	A	1053	ILE	3.8
1	A	1065	ALA	3.8
1	A	1084	ILE	3.8
1	A	1063	VAL	3.7
1	A	1115	ILE	3.6
1	A	1072	GLY	3.6
1	A	1102	GLY	3.5
1	A	1224	GLY	3.5
1	A	1244	LYS	3.4
1	A	1150	GLU	3.3
1	A	1148	ARG	3.1
1	A	1151	GLY	3.1
1	A	1144	GLY	3.0
1	A	1056	SER	2.9
1	A	1096	THR	2.9
1	A	1085	GLY	2.8
1	A	1153	PRO	2.8
1	A	1124	PHE	2.7
1	A	1289	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	1143	LEU	2.6
1	A	1152	SER	2.5
1	A	1087	GLY	2.5
1	A	1054	ASP	2.4
1	A	1146	CYS	2.4
1	A	1090	GLY	2.4
1	A	1058	LEU	2.4
1	A	1132	LYS	2.3
1	A	1119	GLY	2.2
1	A	1094	HIS	2.1
1	A	1078	VAL	2.1
1	A	1154	LEU	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
2	319	A	2001	29/29	0.92	0.13	-0.66	34,37,41,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.