



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 08:11 PM GMT

PDB ID : 3C5U
Title : P38 ALPHA map kinase complexed with a benzothiazole based inhibitor
Authors : Sack, J.S.
Deposited on : 2008-02-01
Resolution : 2.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

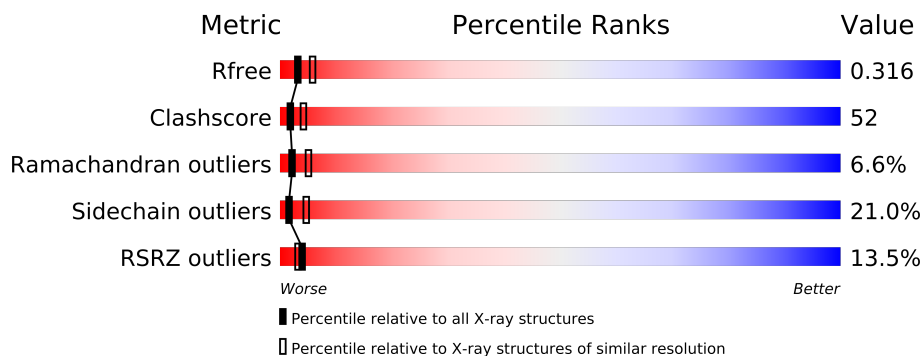
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	366	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 2854 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

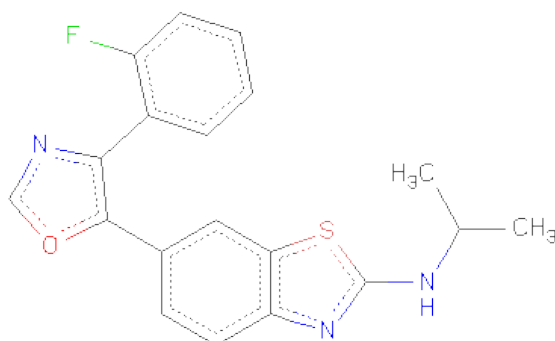
- Molecule 1 is a protein called Mitogen-activated protein kinase 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	338	2728	1751	465	500	12	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	EXPRESSION TAG	UNP Q16539
A	-4	ALA	-	EXPRESSION TAG	UNP Q16539
A	-3	HIS	-	EXPRESSION TAG	UNP Q16539
A	-2	HIS	-	EXPRESSION TAG	UNP Q16539
A	-1	HIS	-	EXPRESSION TAG	UNP Q16539
A	0	HIS	-	EXPRESSION TAG	UNP Q16539
A	1	HIS	-	EXPRESSION TAG	UNP Q16539

- Molecule 2 is 6-[4-(2-FLUOROPHENYL)-1,3-OXAZOL-5-YL]-N-(1-METHYLETHYL)-1,3-BENZOTHAZOL-2-AMINE (three-letter code: P41) (formula: C₁₉H₁₆FN₃OS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			25	19	1	3	1	1		
2	A	1	Total	C	F	N	O	S	0	0
			25	19	1	3	1	1		

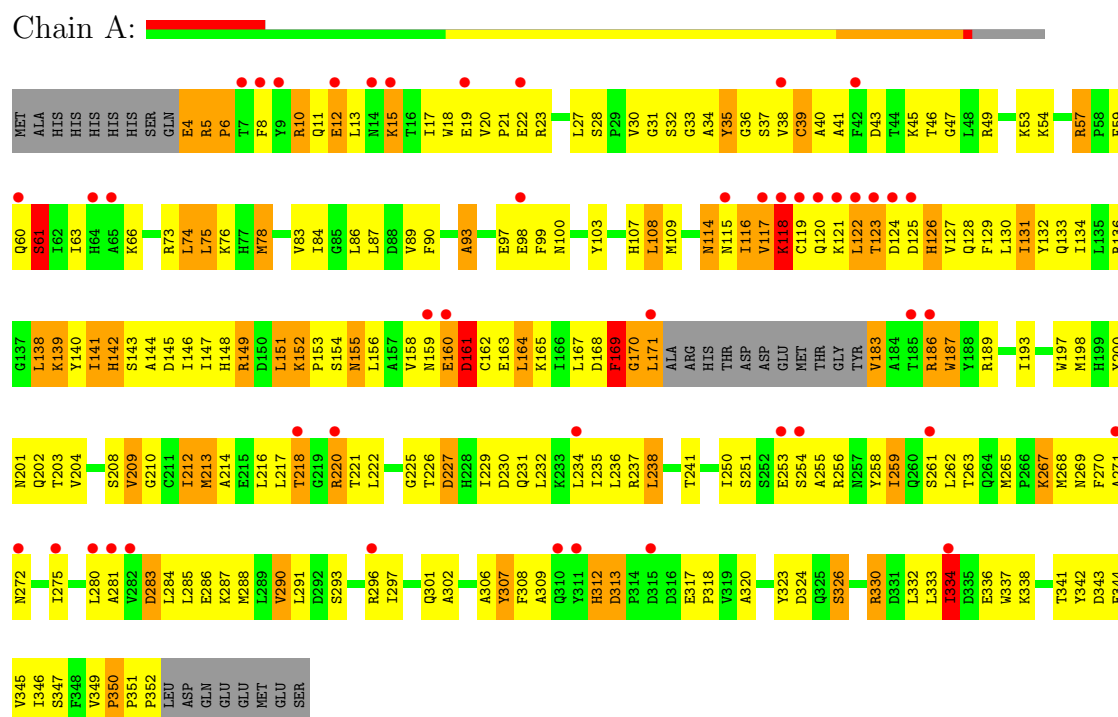
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Mitogen-activated protein kinase 14



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.40Å 80.40Å 225.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	18.26 – 2.80 46.13 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.7 (18.26-2.80) 90.9 (46.13-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.81Å)	Xtriage
Refinement program	BUSTER-TNT 2.1.1	Depositor
R, R_{free}	0.261 , 0.296 0.275 , 0.316	Depositor DCC
R_{free} test set	876 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	67.9	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 84.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 17263 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	2854	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: P41

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	0.36	0/2791	0.62	1/3789 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	197	TRP	N-CA-C	5.17	124.96	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2728	0	2731	285	0
2	A	50	0	32	14	0
3	A	76	0	0	7	0
All	All	2854	0	2763	286	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 52.

All (286) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:147:ILE:HD13	1:A:202:GLN:HA	1.43	1.01
1:A:114:ASN:ND2	1:A:153:PRO:HB2	1.81	0.94
1:A:220:ARG:HD3	1:A:221:THR:N	1.83	0.93
1:A:127:VAL:HG22	1:A:216:LEU:HB3	1.50	0.92
1:A:220:ARG:HD3	1:A:221:THR:H	1.35	0.92
1:A:78:MET:HG2	1:A:169:PHE:HZ	1.35	0.91
1:A:253:GLU:HG3	1:A:256:ARG:NH2	1.85	0.91
1:A:259:ILE:HA	1:A:262:LEU:HD12	1.50	0.90
1:A:186:ARG:CG	1:A:187:TRP:H	1.84	0.90
1:A:250:ILE:HG12	2:A:361:P41:H18B	1.55	0.88
1:A:127:VAL:HG11	1:A:217:LEU:HD21	1.54	0.87
1:A:186:ARG:HG3	1:A:187:TRP:H	1.40	0.86
1:A:59:PHE:CD2	1:A:338:LYS:HD2	2.11	0.85
1:A:213:MET:CE	1:A:284:LEU:HD23	2.06	0.85
1:A:114:ASN:HD22	1:A:153:PRO:HB2	1.42	0.85
1:A:84:ILE:HD13	1:A:167:LEU:HA	1.59	0.82
1:A:5:ARG:HG2	1:A:5:ARG:HH11	1.46	0.81
1:A:124:ASP:O	1:A:128:GLN:HB2	1.80	0.81
1:A:78:MET:HE1	1:A:141:ILE:HG13	1.61	0.81
1:A:259:ILE:HA	1:A:262:LEU:CD1	2.11	0.80
1:A:155:ASN:HB3	1:A:168:ASP:HB2	1.63	0.80
1:A:186:ARG:CD	1:A:187:TRP:H	1.94	0.80
1:A:293:SER:HA	1:A:296:ARG:HG3	1.63	0.80
1:A:117:VAL:HG21	1:A:216:LEU:CD2	2.11	0.80
1:A:10:ARG:HH21	1:A:17:ILE:HD13	1.45	0.80
1:A:108:LEU:HD12	1:A:109:MET:H	1.47	0.78
1:A:117:VAL:O	1:A:120:GLN:HB2	1.83	0.78
1:A:214:ALA:O	1:A:218:THR:HG23	1.84	0.77
1:A:20:VAL:HG11	1:A:27:LEU:HD21	1.67	0.77
1:A:258:TYR:CE1	1:A:262:LEU:HD11	2.20	0.77
1:A:114:ASN:O	1:A:118:LYS:HG2	1.87	0.75
1:A:117:VAL:HG21	1:A:216:LEU:HD22	1.69	0.74
1:A:167:LEU:HD11	2:A:362:P41:C4	2.18	0.74
1:A:27:LEU:HD12	1:A:41:ALA:HB2	1.69	0.74
1:A:10:ARG:NH2	1:A:17:ILE:HD13	2.04	0.73
1:A:87:LEU:HD11	1:A:107:HIS:CE1	2.24	0.73
1:A:54:LYS:HG3	1:A:103:TYR:CE2	2.23	0.73
1:A:226:THR:HG22	1:A:230:ASP:OD2	1.89	0.72
1:A:133:GLN:HE22	1:A:163:GLU:HA	1.55	0.72
1:A:343:ASP:O	1:A:347:SER:HB3	1.90	0.72
1:A:73:ARG:HH11	1:A:73:ARG:HG3	1.54	0.72
1:A:4:GLU:HG3	1:A:5:ARG:N	2.05	0.71
1:A:127:VAL:HG22	1:A:216:LEU:CB	2.20	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:140:TYR:CE1	1:A:320:ALA:HA	2.24	0.71
1:A:164:LEU:O	1:A:164:LEU:HD23	1.90	0.71
2:A:362:P41:S25	2:A:362:P41:H17A	2.30	0.71
1:A:127:VAL:CG2	1:A:216:LEU:HB3	2.21	0.71
1:A:259:ILE:CA	1:A:262:LEU:HD12	2.21	0.70
1:A:186:ARG:HG3	1:A:187:TRP:N	2.05	0.70
1:A:213:MET:HE3	1:A:285:LEU:HG	1.74	0.70
1:A:146:ILE:HD11	1:A:323:TYR:CD2	2.27	0.70
1:A:149:ARG:HG2	1:A:149:ARG:HH11	1.56	0.70
1:A:127:VAL:HG11	1:A:217:LEU:CD2	2.20	0.69
1:A:116:ILE:HD13	1:A:116:ILE:N	2.06	0.69
1:A:318:PRO:HD2	3:A:405:HOH:O	1.92	0.69
1:A:259:ILE:HD11	2:A:361:P41:H5	1.72	0.69
1:A:78:MET:HG2	1:A:169:PHE:CZ	2.23	0.68
1:A:210:GLY:CA	1:A:288:MET:HE3	2.24	0.68
1:A:118:LYS:NZ	1:A:120:GLN:OE1	2.28	0.67
1:A:258:TYR:CD1	1:A:262:LEU:HD11	2.30	0.67
1:A:186:ARG:HD3	1:A:189:ARG:CB	2.24	0.67
1:A:6:PRO:HG2	1:A:8:PHE:CE1	2.29	0.67
1:A:75:LEU:N	1:A:75:LEU:HD23	2.10	0.67
1:A:269:ASN:ND2	1:A:271:ALA:HB3	2.10	0.67
1:A:12:GLU:O	1:A:13:LEU:HG	1.93	0.67
1:A:147:ILE:CD1	1:A:202:GLN:HA	2.21	0.67
1:A:15:LYS:O	1:A:15:LYS:HD3	1.95	0.67
1:A:138:LEU:HD11	1:A:151:LEU:CD2	2.24	0.66
1:A:250:ILE:CG1	2:A:361:P41:H18B	2.24	0.66
1:A:148:HIS:CD2	1:A:169:PHE:HB3	2.31	0.66
1:A:210:GLY:HA2	1:A:288:MET:HE3	1.78	0.66
1:A:144:ALA:HB3	1:A:146:ILE:HG13	1.78	0.65
1:A:147:ILE:H	1:A:171:LEU:HG	1.61	0.65
1:A:31:GLY:HA3	1:A:35:TYR:HE1	1.62	0.65
1:A:159:ASN:O	1:A:161:ASP:N	2.30	0.65
1:A:31:GLY:HA3	1:A:35:TYR:CE1	2.32	0.65
1:A:78:MET:HE3	1:A:140:TYR:CE2	2.32	0.64
1:A:333:LEU:HD12	1:A:336:GLU:OE1	1.96	0.64
1:A:61:SER:HA	1:A:334:ILE:HD11	1.79	0.64
1:A:108:LEU:HD12	1:A:109:MET:N	2.13	0.64
1:A:297:ILE:CG1	1:A:301:GLN:HB2	2.28	0.64
1:A:141:ILE:HG22	1:A:142:HIS:N	2.12	0.64
1:A:76:LYS:NZ	1:A:347:SER:OG	2.31	0.64
1:A:291:LEU:CD2	2:A:361:P41:H17	2.28	0.63
1:A:167:LEU:HD11	2:A:362:P41:C9	2.27	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:210:GLY:HA2	1:A:288:MET:CE	2.28	0.63
1:A:226:THR:HG23	1:A:227:ASP:OD2	1.98	0.63
1:A:30:VAL:HG23	1:A:31:GLY:N	2.14	0.63
1:A:20:VAL:HG23	1:A:21:PRO:O	1.99	0.63
1:A:17:ILE:HD12	3:A:417:HOH:O	1.97	0.63
1:A:31:GLY:O	1:A:33:GLY:N	2.32	0.63
1:A:117:VAL:O	1:A:118:LYS:HG3	1.99	0.62
1:A:186:ARG:HD3	1:A:189:ARG:HB2	1.80	0.62
1:A:213:MET:HE1	1:A:284:LEU:HD23	1.81	0.62
1:A:297:ILE:HD11	1:A:302:ALA:N	2.15	0.62
1:A:158:VAL:HG23	1:A:163:GLU:O	2.00	0.62
1:A:271:ALA:O	1:A:275:ILE:HG12	2.00	0.61
1:A:186:ARG:CD	1:A:189:ARG:HB2	2.29	0.61
1:A:270:PHE:CD2	1:A:286:GLU:HG2	2.35	0.61
1:A:342:TYR:O	1:A:346:ILE:HG12	2.01	0.61
1:A:253:GLU:HG3	1:A:256:ARG:HH22	1.66	0.61
1:A:27:LEU:HD12	1:A:41:ALA:CB	2.30	0.60
1:A:183:VAL:HA	3:A:435:HOH:O	2.02	0.60
1:A:73:ARG:NH1	1:A:73:ARG:HG3	2.10	0.60
1:A:61:SER:CA	1:A:334:ILE:HD11	2.31	0.60
1:A:306:ALA:O	1:A:309:ALA:N	2.34	0.60
1:A:66:LYS:HD2	1:A:337:TRP:CZ2	2.36	0.60
1:A:127:VAL:HG21	1:A:217:LEU:HD23	1.84	0.59
1:A:267:LYS:NZ	1:A:286:GLU:OE1	2.36	0.59
1:A:147:ILE:O	1:A:149:ARG:HD2	2.02	0.59
1:A:35:TYR:O	1:A:38:VAL:HG23	2.03	0.59
1:A:60:GLN:O	1:A:61:SER:HB3	2.03	0.58
1:A:208:SER:O	1:A:212:ILE:HG13	2.03	0.58
1:A:78:MET:CG	1:A:169:PHE:HZ	2.12	0.57
1:A:8:PHE:HD2	1:A:19:GLU:O	1.87	0.57
1:A:134:ILE:HD13	1:A:156:LEU:CD1	2.33	0.57
1:A:117:VAL:CG2	1:A:216:LEU:HD22	2.34	0.57
1:A:259:ILE:CD1	2:A:361:P41:H5	2.35	0.57
1:A:138:LEU:HD11	1:A:151:LEU:HD21	1.85	0.56
1:A:168:ASP:O	1:A:169:PHE:HB3	2.05	0.56
1:A:250:ILE:HD13	1:A:259:ILE:HD11	1.87	0.56
1:A:146:ILE:HG23	1:A:171:LEU:HB2	1.87	0.56
1:A:12:GLU:HG3	1:A:13:LEU:N	2.19	0.56
1:A:133:GLN:HB2	1:A:164:LEU:HD22	1.87	0.56
1:A:57:ARG:HB3	1:A:60:GLN:HG3	1.86	0.56
1:A:186:ARG:HD3	1:A:189:ARG:HB3	1.88	0.56
1:A:220:ARG:HH11	1:A:221:THR:H	1.54	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:225:GLY:CA	1:A:234:LEU:HD12	2.35	0.55
1:A:146:ILE:HA	1:A:171:LEU:HD12	1.87	0.55
1:A:290:VAL:HG12	1:A:296:ARG:HG2	1.88	0.55
1:A:297:ILE:HG13	1:A:301:GLN:HB2	1.87	0.55
1:A:31:GLY:C	1:A:33:GLY:H	2.10	0.55
1:A:270:PHE:CE2	1:A:286:GLU:HA	2.42	0.55
1:A:160:GLU:O	1:A:160:GLU:HG2	2.07	0.55
1:A:131:ILE:HD12	1:A:134:ILE:HD12	1.90	0.55
1:A:229:ILE:HG23	1:A:258:TYR:OH	2.08	0.54
1:A:118:LYS:HA	3:A:364:HOH:O	2.08	0.54
1:A:139:LYS:HE3	1:A:317:GLU:O	2.07	0.54
1:A:158:VAL:HG21	1:A:162:CYS:HA	1.90	0.54
1:A:324:ASP:OD1	1:A:326:SER:HB3	2.08	0.53
1:A:63:ILE:HD13	3:A:437:HOH:O	2.09	0.53
1:A:280:LEU:HD13	1:A:306:ALA:CB	2.38	0.53
1:A:43:ASP:OD2	1:A:46:THR:HG23	2.08	0.53
1:A:288:MET:O	1:A:296:ARG:HD3	2.09	0.53
1:A:134:ILE:HD13	1:A:156:LEU:HD11	1.89	0.53
1:A:10:ARG:HD2	1:A:19:GLU:OE1	2.09	0.53
1:A:20:VAL:HG11	1:A:27:LEU:CD2	2.38	0.53
1:A:226:THR:HG22	1:A:230:ASP:CG	2.30	0.53
1:A:21:PRO:HD3	1:A:90:PHE:CE1	2.43	0.53
1:A:186:ARG:CG	1:A:187:TRP:N	2.56	0.53
1:A:118:LYS:O	1:A:119:CYS:HB2	2.08	0.52
1:A:291:LEU:HD23	2:A:361:P41:H17	1.90	0.52
1:A:164:LEU:HD23	1:A:164:LEU:C	2.29	0.52
1:A:186:ARG:HH22	1:A:235:ILE:HD11	1.73	0.52
1:A:10:ARG:HD2	1:A:19:GLU:CD	2.29	0.52
1:A:148:HIS:O	1:A:149:ARG:HB2	2.10	0.52
1:A:5:ARG:NH1	1:A:5:ARG:HG2	2.20	0.52
1:A:149:ARG:HG2	1:A:149:ARG:NH1	2.24	0.52
1:A:280:LEU:CD1	1:A:306:ALA:HB1	2.40	0.52
1:A:225:GLY:HA2	1:A:234:LEU:CD1	2.39	0.52
1:A:148:HIS:CD2	1:A:169:PHE:CB	2.93	0.52
1:A:281:ALA:O	1:A:284:LEU:HB3	2.10	0.51
1:A:226:THR:HG22	1:A:230:ASP:CB	2.40	0.51
1:A:43:ASP:O	1:A:47:GLY:N	2.37	0.51
1:A:146:ILE:HG23	1:A:171:LEU:CB	2.41	0.51
1:A:147:ILE:H	1:A:171:LEU:CG	2.22	0.51
1:A:76:LYS:HA	1:A:86:LEU:HD12	1.93	0.51
1:A:8:PHE:HA	1:A:21:PRO:HA	1.93	0.51
1:A:20:VAL:HG23	1:A:21:PRO:N	2.26	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:269:ASN:HD21	1:A:271:ALA:HB3	1.74	0.50
1:A:250:ILE:HD11	2:A:361:P41:H17A	1.93	0.50
1:A:213:MET:SD	1:A:284:LEU:HD23	2.51	0.50
1:A:312:HIS:O	1:A:313:ASP:HB2	2.11	0.50
1:A:141:ILE:HG23	1:A:146:ILE:HB	1.93	0.50
1:A:134:ILE:CD1	1:A:156:LEU:HD13	2.42	0.50
1:A:349:VAL:O	1:A:350:PRO:O	2.30	0.50
1:A:250:ILE:CD1	1:A:259:ILE:HD11	2.42	0.50
1:A:291:LEU:HD21	2:A:361:P41:H17	1.94	0.50
1:A:84:ILE:HD12	1:A:165:LYS:HB3	1.94	0.50
1:A:272:ASN:HA	1:A:275:ILE:HD11	1.94	0.50
1:A:186:ARG:CD	1:A:187:TRP:N	2.69	0.50
1:A:186:ARG:CD	1:A:189:ARG:CB	2.88	0.50
1:A:83:VAL:HG11	1:A:141:ILE:HD11	1.94	0.50
1:A:78:MET:HG3	1:A:83:VAL:HG11	1.94	0.49
1:A:86:LEU:HD22	1:A:89:VAL:HG22	1.93	0.49
1:A:146:ILE:HD11	1:A:323:TYR:HD2	1.77	0.49
1:A:169:PHE:HB2	1:A:170:GLY:HA3	1.94	0.49
1:A:78:MET:CE	1:A:140:TYR:CE2	2.95	0.48
1:A:148:HIS:NE2	1:A:168:ASP:O	2.43	0.48
1:A:75:LEU:HD12	2:A:362:P41:H1	1.96	0.48
1:A:241:THR:HG23	1:A:265:MET:H	1.79	0.48
1:A:119:CYS:SG	1:A:220:ARG:NH2	2.87	0.48
1:A:236:LEU:HD13	1:A:241:THR:HG22	1.96	0.47
1:A:306:ALA:O	1:A:308:PHE:N	2.46	0.47
1:A:253:GLU:HG3	1:A:256:ARG:HH21	1.71	0.47
1:A:213:MET:HG2	1:A:285:LEU:HD21	1.95	0.47
1:A:159:ASN:OD1	1:A:159:ASN:O	2.33	0.47
1:A:297:ILE:HD11	1:A:301:GLN:C	2.34	0.47
1:A:297:ILE:HG13	1:A:301:GLN:CB	2.43	0.47
1:A:333:LEU:O	1:A:336:GLU:N	2.46	0.47
1:A:270:PHE:CD1	1:A:270:PHE:N	2.80	0.47
1:A:35:TYR:HB2	1:A:38:VAL:HG21	1.97	0.47
1:A:127:VAL:HG12	1:A:307:TYR:OH	2.15	0.47
1:A:141:ILE:HG22	1:A:142:HIS:HD2	1.80	0.47
1:A:297:ILE:HG12	1:A:301:GLN:HB2	1.97	0.47
1:A:31:GLY:CA	1:A:35:TYR:HE1	2.28	0.47
1:A:270:PHE:N	1:A:270:PHE:HD1	2.12	0.47
1:A:126:HIS:O	1:A:130:LEU:HG	2.15	0.47
1:A:186:ARG:NH2	1:A:235:ILE:HD11	2.30	0.46
1:A:30:VAL:CG2	1:A:31:GLY:N	2.78	0.46
1:A:97:GLU:CD	1:A:97:GLU:H	2.18	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:140:TYR:CE1	1:A:320:ALA:CA	2.95	0.46
1:A:141:ILE:CG2	1:A:142:HIS:N	2.78	0.46
1:A:15:LYS:HD3	1:A:15:LYS:C	2.33	0.46
1:A:141:ILE:O	1:A:143:SER:N	2.49	0.46
1:A:155:ASN:CB	1:A:168:ASP:HB2	2.42	0.45
1:A:40:ALA:HB1	1:A:108:LEU:CD2	2.47	0.45
1:A:78:MET:CE	1:A:140:TYR:HE2	2.29	0.45
1:A:213:MET:HE2	1:A:284:LEU:HD23	1.94	0.45
1:A:49:ARG:HD2	3:A:413:HOH:O	2.16	0.45
1:A:209:VAL:HG12	1:A:288:MET:CE	2.46	0.45
1:A:349:VAL:O	1:A:350:PRO:C	2.53	0.45
1:A:5:ARG:CG	1:A:5:ARG:NH1	2.76	0.45
1:A:312:HIS:CG	1:A:313:ASP:N	2.85	0.45
1:A:201:ASN:OD1	1:A:203:THR:HG23	2.16	0.45
1:A:155:ASN:OD1	1:A:155:ASN:N	2.48	0.45
1:A:75:LEU:CD2	1:A:75:LEU:N	2.78	0.45
1:A:117:VAL:HG21	1:A:216:LEU:HD23	1.95	0.44
1:A:238:LEU:HD11	1:A:270:PHE:CD1	2.52	0.44
1:A:280:LEU:HD11	1:A:306:ALA:HB1	1.98	0.44
1:A:38:VAL:HG22	1:A:53:LYS:HB2	2.00	0.44
1:A:238:LEU:HD11	1:A:270:PHE:CE1	2.53	0.44
1:A:4:GLU:HG3	1:A:5:ARG:H	1.80	0.44
1:A:27:LEU:CD1	1:A:41:ALA:HB2	2.43	0.44
1:A:297:ILE:HG13	1:A:301:GLN:OE1	2.18	0.44
1:A:129:PHE:O	1:A:132:TYR:HB3	2.18	0.44
1:A:109:MET:SD	1:A:165:LYS:HE3	2.58	0.43
1:A:297:ILE:CD1	1:A:302:ALA:N	2.80	0.43
1:A:312:HIS:CG	1:A:313:ASP:H	2.36	0.43
1:A:117:VAL:HG22	1:A:122:LEU:HD21	2.00	0.43
1:A:186:ARG:O	1:A:187:TRP:CB	2.65	0.43
1:A:280:LEU:HD13	1:A:306:ALA:HB3	2.00	0.43
1:A:225:GLY:HA2	1:A:234:LEU:HD12	1.96	0.43
1:A:49:ARG:HB2	3:A:388:HOH:O	2.18	0.43
1:A:280:LEU:HD23	1:A:280:LEU:HA	1.68	0.43
1:A:290:VAL:CG1	1:A:296:ARG:HG2	2.48	0.43
1:A:341:THR:O	1:A:345:VAL:HG23	2.18	0.43
1:A:131:ILE:HG22	1:A:308:PHE:CZ	2.54	0.43
1:A:93:ALA:CB	1:A:98:GLU:HB3	2.48	0.43
1:A:283:ASP:O	1:A:287:LYS:HG3	2.18	0.43
1:A:45:LYS:HB2	1:A:45:LYS:HE3	1.42	0.43
1:A:258:TYR:CG	2:A:361:P41:H3	2.53	0.43
1:A:187:TRP:HA	1:A:231:GLN:HE22	1.84	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:18:TRP:CH2	1:A:39:CYS:HB2	2.54	0.43
1:A:74:LEU:HD21	1:A:146:ILE:CD1	2.49	0.43
1:A:159:ASN:OD1	1:A:161:ASP:HB2	2.18	0.42
1:A:152:LYS:HB2	1:A:153:PRO:CD	2.50	0.42
1:A:232:LEU:HD22	2:A:361:P41:C15	2.49	0.42
1:A:59:PHE:CE2	1:A:338:LYS:HD2	2.54	0.42
1:A:117:VAL:HG12	1:A:118:LYS:N	2.34	0.42
1:A:202:GLN:C	1:A:204:VAL:H	2.23	0.42
1:A:75:LEU:HD23	1:A:75:LEU:H	1.81	0.42
1:A:13:LEU:HB2	1:A:18:TRP:CD1	2.55	0.42
1:A:209:VAL:HG12	1:A:288:MET:HE1	2.01	0.42
1:A:31:GLY:C	1:A:33:GLY:N	2.74	0.42
1:A:351:PRO:HA	1:A:352:PRO:HD3	1.77	0.42
1:A:193:ILE:HD13	1:A:200:TYR:CE1	2.55	0.42
1:A:255:ALA:O	1:A:259:ILE:HG13	2.20	0.41
1:A:258:TYR:C	1:A:258:TYR:CD1	2.93	0.41
1:A:76:LYS:HD3	1:A:344:GLU:OE1	2.19	0.41
1:A:123:THR:O	1:A:126:HIS:N	2.53	0.41
1:A:158:VAL:HA	1:A:163:GLU:O	2.21	0.41
1:A:54:LYS:HG3	1:A:103:TYR:CZ	2.55	0.41
1:A:209:VAL:HG12	1:A:210:GLY:N	2.36	0.41
1:A:152:LYS:HG2	1:A:155:ASN:OD1	2.21	0.41
1:A:297:ILE:CD1	1:A:302:ALA:CA	2.99	0.41
1:A:269:ASN:HB3	1:A:272:ASN:ND2	2.35	0.41
1:A:99:PHE:O	1:A:338:LYS:HE3	2.21	0.40
1:A:267:LYS:HD3	1:A:268:MET:O	2.21	0.40
1:A:330:ARG:HG2	1:A:332:LEU:HD21	2.03	0.40
1:A:118:LYS:HG3	1:A:118:LYS:O	2.21	0.40
1:A:158:VAL:CG2	1:A:162:CYS:HA	2.51	0.40
1:A:269:ASN:HB3	1:A:272:ASN:HD22	1.86	0.40
1:A:123:THR:HG23	1:A:126:HIS:HB2	2.03	0.40
1:A:280:LEU:CD1	1:A:306:ALA:CB	3.00	0.40
1:A:144:ALA:HB1	1:A:323:TYR:HB2	2.04	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	334/366 (91%)	274 (82%)	38 (11%)	22 (7%)	2 5

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	32	SER
1	A	34	ALA
1	A	36	GLY
1	A	117	VAL
1	A	160	GLU
1	A	161	ASP
1	A	307	TYR
1	A	61	SER
1	A	116	ILE
1	A	118	LYS
1	A	187	TRP
1	A	198	MET
1	A	312	HIS
1	A	93	ALA
1	A	142	HIS
1	A	169	PHE
1	A	313	ASP
1	A	350	PRO
1	A	186	ARG
1	A	334	ILE
1	A	6	PRO
1	A	170	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	300/325 (92%)	237 (79%)	63 (21%)	1 5

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	5	ARG
1	A	10	ARG
1	A	11	GLN
1	A	12	GLU
1	A	15	LYS
1	A	22	GLU
1	A	23	ARG
1	A	28	SER
1	A	35	TYR
1	A	37	SER
1	A	39	CYS
1	A	57	ARG
1	A	61	SER
1	A	74	LEU
1	A	75	LEU
1	A	78	MET
1	A	100	ASN
1	A	108	LEU
1	A	114	ASN
1	A	115	ASN
1	A	118	LYS
1	A	121	LYS
1	A	122	LEU
1	A	123	THR
1	A	125	ASP
1	A	126	HIS
1	A	131	ILE
1	A	136	ARG
1	A	138	LEU
1	A	139	LYS
1	A	141	ILE
1	A	145	ASP
1	A	149	ARG
1	A	151	LEU
1	A	152	LYS
1	A	154	SER
1	A	155	ASN
1	A	161	ASP
1	A	164	LEU
1	A	169	PHE
1	A	171	LEU
1	A	183	VAL

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Mol	Chain	Res	Type
1	A	209	VAL
1	A	212	ILE
1	A	213	MET
1	A	218	THR
1	A	220	ARG
1	A	222	LEU
1	A	227	ASP
1	A	237	ARG
1	A	238	LEU
1	A	251	SER
1	A	254	SER
1	A	259	ILE
1	A	261	SER
1	A	263	THR
1	A	267	LYS
1	A	283	ASP
1	A	290	VAL
1	A	326	SER
1	A	330	ARG
1	A	334	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	60	GLN
1	A	114	ASN
1	A	133	GLN
1	A	199	HIS
1	A	257	ASN
1	A	264	GLN
1	A	272	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

2 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	$\# Z > 2$
2	P41	A	361	-	27,28,28	1.77	3 (11%)	38,40,40	3.90	16 (42%)
2	P41	A	362	-	27,28,28	1.28	2 (7%)	38,40,40	2.66	15 (39%)

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	P41	A	361	-	-	0/8/12/12	0/2/4/4
2	P41	A	362	-	-	0/8/12/12	0/2/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	361	P41	C10-C14	-7.30	1.42	1.49
2	A	362	P41	C10-C14	-4.67	1.44	1.49
2	A	361	P41	C5-C4	2.97	1.42	1.36
2	A	361	P41	C16-N21	-2.18	1.31	1.34
2	A	362	P41	C13-S25	-2.17	1.71	1.74

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	361	P41	C8-O23-C15	17.77	107.52	103.54
2	A	362	P41	C9-C15-C14	7.66	142.61	127.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	361	P41	C9-C15-C14	7.62	142.53	127.53
2	A	362	P41	S25-C16-N22	6.19	125.75	121.80
2	A	361	P41	C16-S25-C13	5.64	91.16	88.89
2	A	361	P41	C3-C10-C12	5.12	121.10	115.78
2	A	362	P41	C18-C19-N22	-4.96	98.50	109.90
2	A	361	P41	C16-N22-C19	4.76	129.01	124.06
2	A	362	P41	S25-C16-N21	-4.05	113.00	120.99
2	A	362	P41	O23-C15-C9	-4.04	110.22	117.12
2	A	361	P41	C7-C13-C11	3.78	123.48	120.82
2	A	362	P41	C17-C19-N22	3.69	118.40	109.90
2	A	362	P41	C3-C10-C12	3.67	119.59	115.78
2	A	361	P41	C17-C19-C18	3.58	119.26	111.76
2	A	362	P41	C16-S25-C13	3.51	90.30	88.89
2	A	361	P41	C11-C13-S25	-3.42	107.73	112.03
2	A	361	P41	C13-C11-N21	3.34	116.19	108.37
2	A	362	P41	C17-C19-C18	-3.24	104.98	111.76
2	A	361	P41	S25-C16-N21	-3.21	114.66	120.99
2	A	361	P41	C17-C19-N22	2.83	116.41	109.90
2	A	362	P41	C8-O23-C15	2.79	104.17	103.54
2	A	361	P41	C18-C19-N22	2.76	116.26	109.90
2	A	362	P41	C13-C11-N21	2.60	114.47	108.37
2	A	362	P41	C4-C5-C11	-2.52	117.97	120.92
2	A	361	P41	C6-C12-C10	-2.44	120.56	123.58
2	A	362	P41	C16-N22-C19	2.28	126.43	124.06
2	A	361	P41	C15-C14-N20	2.26	112.01	107.97
2	A	362	P41	C7-C13-C11	2.23	122.39	120.82
2	A	361	P41	C4-C5-C11	-2.07	118.50	120.92
2	A	361	P41	F24-C12-C10	2.06	121.75	118.79
2	A	362	P41	C9-C7-C13	-2.02	117.52	121.24

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	338/366 (92%)	0.89	46 (13%) 4 3	44, 70, 105, 137	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	119	CYS	13.5
1	A	118	LYS	7.1
1	A	271	ALA	5.7
1	A	125	ASP	4.9
1	A	185	THR	4.9
1	A	220	ARG	4.6
1	A	64	HIS	4.6
1	A	8	PHE	4.1
1	A	120	GLN	4.1
1	A	122	LEU	4.1
1	A	261	SER	4.0
1	A	15	LYS	3.9
1	A	310	GLN	3.8
1	A	275	ILE	3.8
1	A	218	THR	3.7
1	A	115	ASN	3.7
1	A	272	ASN	3.6
1	A	42	PHE	3.6
1	A	254	SER	3.5
1	A	14	ASN	3.2
1	A	12	GLU	3.2
1	A	280	LEU	3.1
1	A	124	ASP	3.1
1	A	253	GLU	3.1
1	A	311	TYR	3.0
1	A	9	TYR	3.0
1	A	38	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	7	THR	2.8
1	A	171	LEU	2.7
1	A	282	VAL	2.7
1	A	22	GLU	2.6
1	A	121	LYS	2.6
1	A	123	THR	2.5
1	A	159	ASN	2.5
1	A	296	ARG	2.5
1	A	315	ASP	2.4
1	A	65	ALA	2.4
1	A	234	LEU	2.3
1	A	117	VAL	2.2
1	A	160	GLU	2.2
1	A	334	ILE	2.2
1	A	60	GLN	2.2
1	A	186	ARG	2.1
1	A	281	ALA	2.1
1	A	98	GLU	2.1
1	A	19	GLU	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. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	P41	A	361	25/25	0.25	0.14	47,63,72,73	0
2	P41	A	362	25/25	0.19	-1.01	62,74,79,80	0

6.5 Other polymers ⓘ

There are no such residues in this entry.