



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:38 pm GMT

PDB ID : 1M52
Title : Crystal Structure of the c-Abl Kinase domain in complex with PD173955
Authors : Nagar, B.; Bornmann, W.; Pellicena, P.; Schindler, T.; Veach, D.; Miller, W.T.; Clarkson, B.; Kuriyan, J.
Deposited on : 2002-07-08
Resolution : 2.60 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

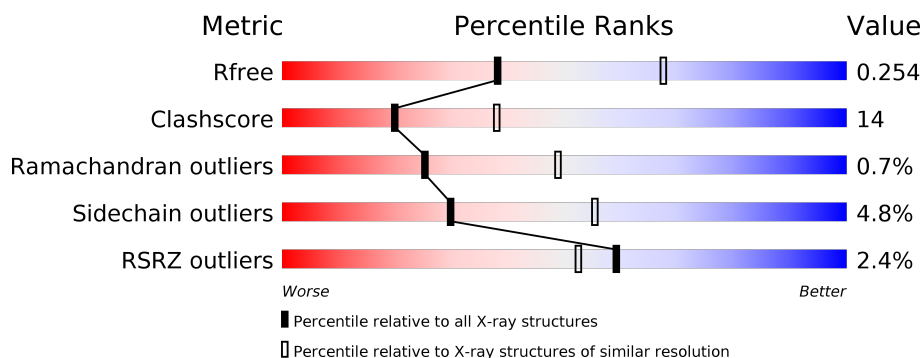
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	293	<div> <div>4%</div> <div> <div></div> <div>65%</div> <div>26%</div> <div>• 8%</div> </div> </div>
1	B	293	<div> <div>%</div> <div> <div></div> <div>66%</div> <div>24%</div> <div>• 7%</div> </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
2	MES	B	117	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4620 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 PROTO-ONCOGENE TYROSINE-PROTEIN KINASE ABL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2208	1423	358	410	17			
1	B	272	Total	C	N	O	S	0	0	0
			2214	1426	359	412	17			

There are 12 discrepancies between the modelled and reference sequences:

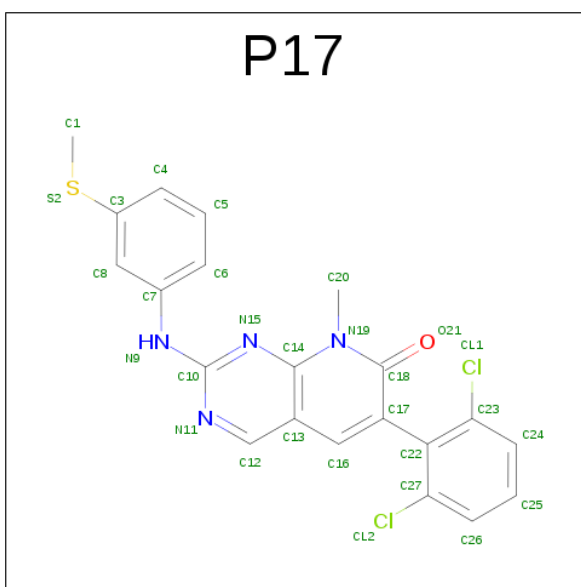
Chain	Residue	Modelled	Actual	Comment	Reference
A	223	GLY	-	CLONING ARTIFACT	UNP P00520
A	224	ALA	-	CLONING ARTIFACT	UNP P00520
A	225	MET	-	CLONING ARTIFACT	UNP P00520
A	226	ASP	-	CLONING ARTIFACT	UNP P00520
A	227	PRO	-	CLONING ARTIFACT	UNP P00520
A	228	SER	-	CLONING ARTIFACT	UNP P00520
B	223	GLY	-	CLONING ARTIFACT	UNP P00520
B	224	ALA	-	CLONING ARTIFACT	UNP P00520
B	225	MET	-	CLONING ARTIFACT	UNP P00520
B	226	ASP	-	CLONING ARTIFACT	UNP P00520
B	227	PRO	-	CLONING ARTIFACT	UNP P00520
B	228	SER	-	CLONING ARTIFACT	UNP P00520

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 3 is 6-(2,6-DICHLORO-PHENYL)-8-METHYL-2-(3-METHYLSULFANYL-PHENYLAMINO)-8H-PYRIDO[2,3-D]PYRIMIDIN-7-ONE (three-letter code: P17) (formula: C₂₁H₁₆Cl₂N₄OS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	S	0	0
			29	21	2	4	1	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	Cl	N	O	S	0	0
			29	21	2	4	1	1		

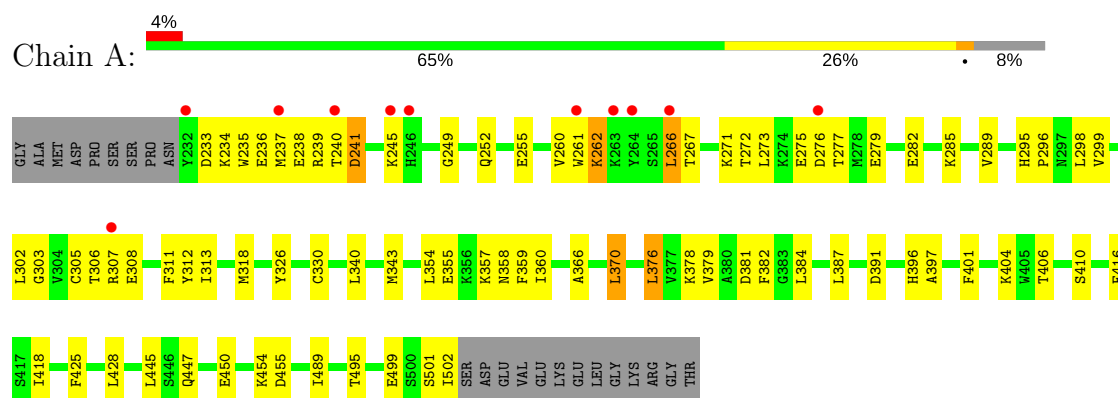
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	47	Total	O	0	0
			47	47		
4	B	69	Total	O	0	0
			69	69		

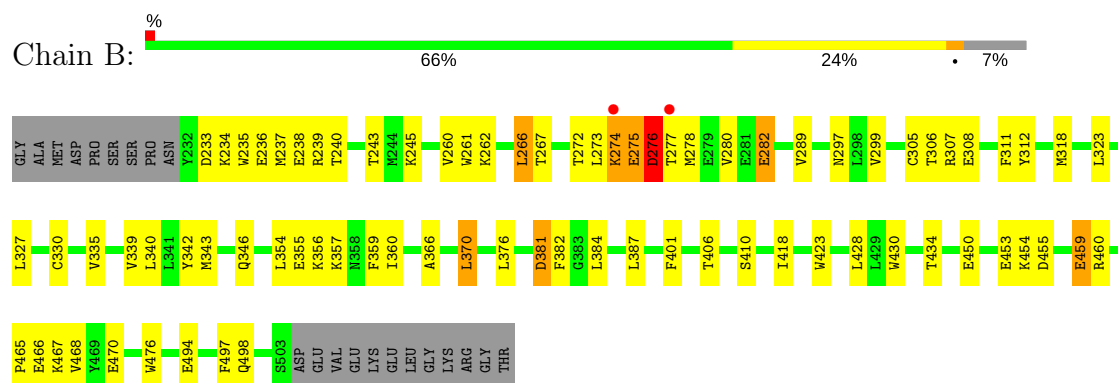
3 Residue-property plots [i](#)

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: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE ABL1



• Molecule 1: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE ABL1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	115.85Å 125.71Å 56.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	55.25 – 2.60 55.25 – 2.49	Depositor EDS
% Data completeness (in resolution range)	94.7 (55.25-2.60) 93.4 (55.25-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.211 , 0.254 0.211 , 0.254	Depositor DCC
R_{free} test set	2469 reflections (9.94%)	DCC
Wilson B-factor (Å ²)	30.7	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.0	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	4620	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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: MES, P17

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	0.37	0/2265	0.62	0/3065
1	B	0.41	0/2271	0.65	0/3073
All	All	0.39	0/4536	0.63	0/6138

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

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	2208	0	2162	63	0
1	B	2214	0	2167	59	0
2	B	24	0	26	0	0
3	A	29	0	16	2	0
3	B	29	0	16	2	0
4	A	47	0	0	0	0
4	B	69	0	0	1	0
All	All	4620	0	4387	126	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 14.

All (126) 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:305:CYS:HB2	1:A:312:TYR:HB2	1.58	0.84
1:B:299:VAL:HB	1:B:381:ASP:OD1	1.80	0.81
1:B:305:CYS:HB2	1:B:312:TYR:HB2	1.64	0.79
1:A:340:LEU:HA	1:A:343:MET:HE2	1.67	0.75
1:B:276:ASP:O	1:B:276:ASP:CG	2.25	0.74
1:B:340:LEU:HA	1:B:343:MET:HE2	1.71	0.73
1:B:357:LYS:HB3	1:B:359:PHE:HD1	1.54	0.72
1:A:495:THR:O	1:A:499:GLU:HG2	1.93	0.69
3:A:119:P17:H16	3:A:119:P17:N15	2.08	0.68
1:A:360:ILE:HG23	1:A:387:LEU:HD21	1.74	0.66
1:A:235:TRP:O	1:A:237:MET:HG3	1.96	0.66
1:A:318:MET:HG3	1:A:370:LEU:HB3	1.78	0.66
1:B:459:GLU:CD	1:B:459:GLU:H	2.00	0.65
1:B:239:ARG:HG3	1:B:312:TYR:CE1	2.32	0.64
1:B:355:GLU:HG3	1:B:418:ILE:HG12	1.80	0.64
1:B:360:ILE:HG23	1:B:387:LEU:HD21	1.81	0.63
1:B:357:LYS:HB3	1:B:359:PHE:CD1	2.33	0.63
1:A:238:GLU:HB3	1:A:241:ASP:OD1	1.99	0.63
1:A:239:ARG:HG3	1:A:312:TYR:CE1	2.34	0.62
1:B:306:THR:HA	1:B:311:PHE:HD2	1.65	0.62
1:A:276:ASP:O	1:A:277:THR:HG23	2.00	0.61
1:A:238:GLU:HG3	1:A:240:THR:H	1.65	0.61
3:B:120:P17:H16	3:B:120:P17:N15	2.17	0.60
1:B:318:MET:HG3	1:B:370:LEU:HB3	1.84	0.60
1:B:381:ASP:O	1:B:384:LEU:HG	2.01	0.60
1:A:273:LEU:HD11	1:A:279:GLU:HB2	1.83	0.60
1:B:282:GLU:HA	1:B:282:GLU:OE1	2.03	0.58
1:A:306:THR:HA	1:A:311:PHE:HD2	1.68	0.58
1:A:275:GLU:HG3	1:A:276:ASP:N	2.19	0.58
1:A:406:THR:CG2	1:A:410:SER:HB2	2.34	0.57
1:A:357:LYS:HB3	1:A:359:PHE:HD1	1.69	0.57
1:A:299:VAL:HB	1:A:381:ASP:OD1	2.05	0.56
1:B:245:LYS:HD2	1:B:260:VAL:CG2	2.36	0.56
1:B:245:LYS:HD2	1:B:260:VAL:HG23	1.88	0.55
1:B:498:GLN:NE2	1:B:498:GLN:HA	2.22	0.54
1:A:255:GLU:HB2	1:A:272:THR:O	2.09	0.53
1:A:454:LYS:O	1:A:455:ASP:HB2	2.09	0.53
1:B:498:GLN:HE21	1:B:498:GLN:HA	1.73	0.53
1:B:307:ARG:HH11	1:B:307:ARG:HG2	1.73	0.53
3:A:119:P17:C6	3:A:119:P17:N15	2.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:PRO:O	1:A:378:LYS:HE2	2.09	0.53
1:A:340:LEU:HD23	1:A:343:MET:HE1	1.91	0.53
1:B:467:LYS:HB2	4:B:34:HOH:O	2.07	0.53
1:B:465:PRO:HG2	1:B:468:VAL:CG2	2.40	0.52
1:B:238:GLU:HG3	1:B:240:THR:H	1.74	0.52
1:A:376:LEU:HD11	1:A:378:LYS:HE3	1.91	0.52
1:A:318:MET:HG3	1:A:370:LEU:CB	2.39	0.51
1:B:238:GLU:HG3	1:B:240:THR:HB	1.92	0.51
1:B:237:MET:SD	1:B:261:TRP:CH2	3.04	0.51
1:B:274:LYS:HG3	1:B:275:GLU:HG2	1.91	0.51
1:A:237:MET:CE	1:A:303:GLY:HA3	2.40	0.51
1:B:342:TYR:HB2	1:B:497:PHE:CE1	2.46	0.50
1:A:366:ALA:N	1:A:428:LEU:HD13	2.27	0.50
1:B:274:LYS:CG	1:B:275:GLU:H	2.24	0.50
1:A:298:LEU:HD23	1:A:379:VAL:HB	1.92	0.50
1:B:340:LEU:HD23	1:B:343:MET:HE1	1.92	0.50
1:A:245:LYS:HE3	1:A:260:VAL:CG2	2.43	0.49
3:B:120:P17:C6	3:B:120:P17:N15	2.75	0.49
1:B:450:GLU:O	1:B:454:LYS:HG2	2.12	0.49
1:A:340:LEU:HD23	1:A:343:MET:CE	2.43	0.49
1:A:305:CYS:CB	1:A:312:TYR:HB2	2.36	0.49
1:A:233:ASP:HB3	1:A:236:GLU:HG2	1.95	0.48
1:B:318:MET:HG3	1:B:370:LEU:CB	2.43	0.48
1:B:406:THR:CG2	1:B:410:SER:HB2	2.43	0.48
1:B:233:ASP:OD1	1:B:235:TRP:HB2	2.14	0.48
1:A:295:HIS:CG	1:A:296:PRO:HD2	2.49	0.47
1:B:274:LYS:HG3	1:B:275:GLU:H	1.79	0.47
1:B:494:GLU:O	1:B:498:GLN:HG2	2.13	0.47
1:B:243:THR:OG1	1:B:262:LYS:HE2	2.14	0.47
1:B:453:GLU:C	1:B:455:ASP:H	2.17	0.47
1:A:326:TYR:O	1:A:330:CYS:HB3	2.14	0.47
1:A:355:GLU:HG3	1:A:418:ILE:HG12	1.96	0.47
1:B:274:LYS:CG	1:B:275:GLU:N	2.78	0.47
1:B:274:LYS:HG3	1:B:275:GLU:N	2.30	0.47
1:B:307:ARG:NH1	1:B:307:ARG:HG2	2.30	0.46
1:A:307:ARG:HG2	1:A:307:ARG:HH11	1.79	0.46
1:A:357:LYS:HB3	1:A:359:PHE:CD1	2.48	0.46
1:A:307:ARG:HG2	1:A:307:ARG:NH1	2.31	0.46
1:A:266:LEU:HD23	1:A:267:THR:N	2.30	0.46
1:A:501:SER:O	1:A:502:ILE:HG13	2.16	0.46
1:A:404:LYS:HE2	1:A:445:LEU:HD21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:ASP:HB3	1:B:236:GLU:HG2	1.98	0.45
1:A:236:GLU:OE2	1:A:306:THR:OG1	2.34	0.45
1:B:273:LEU:HD23	1:B:280:VAL:HG22	1.98	0.45
1:B:339:VAL:O	1:B:343:MET:HG3	2.15	0.45
1:A:382:PHE:C	1:A:382:PHE:CD1	2.90	0.45
1:A:396:HIS:O	1:A:397:ALA:C	2.55	0.45
1:A:450:GLU:O	1:A:454:LYS:HG2	2.16	0.45
1:A:245:LYS:HE3	1:A:260:VAL:HG23	1.99	0.44
1:B:340:LEU:HD23	1:B:343:MET:CE	2.47	0.44
1:A:425:PHE:CE2	1:A:489:ILE:HG21	2.52	0.44
1:A:416:PHE:CD1	1:A:416:PHE:N	2.85	0.44
1:A:237:MET:SD	1:A:261:TRP:CH2	3.11	0.44
1:B:266:LEU:HD23	1:B:267:THR:N	2.32	0.44
1:B:455:ASP:CG	1:B:455:ASP:O	2.56	0.44
1:B:330:CYS:HB2	1:B:335:VAL:HG21	1.99	0.44
1:A:273:LEU:HD11	1:A:279:GLU:CB	2.46	0.43
1:A:381:ASP:O	1:A:384:LEU:HG	2.17	0.43
1:B:366:ALA:N	1:B:428:LEU:HD13	2.34	0.43
1:A:357:LYS:O	1:A:358:ASN:HB2	2.18	0.43
1:B:278:MET:HG3	1:B:278:MET:O	2.19	0.43
1:A:285:LYS:O	1:A:289:VAL:HG23	2.19	0.42
1:A:406:THR:HG22	1:A:410:SER:HB2	2.01	0.42
1:A:275:GLU:HG3	1:A:276:ASP:H	1.82	0.42
1:B:423:TRP:CE3	1:B:476:TRP:HA	2.55	0.42
1:A:391:ASP:OD2	1:A:391:ASP:N	2.51	0.42
1:B:382:PHE:C	1:B:382:PHE:CD1	2.93	0.42
1:B:237:MET:SD	1:B:261:TRP:CZ2	3.12	0.42
1:B:289:VAL:HG13	1:B:359:PHE:CE2	2.54	0.42
1:A:450:GLU:OE2	1:A:454:LYS:HE3	2.19	0.42
1:A:360:ILE:HG23	1:A:387:LEU:CD2	2.47	0.42
1:A:249:GLY:O	1:A:252:GLN:HG2	2.20	0.41
1:A:273:LEU:HA	1:A:273:LEU:HD12	1.87	0.41
1:A:261:TRP:O	1:A:262:LYS:C	2.58	0.41
1:A:261:TRP:HZ3	1:A:302:LEU:HB3	1.86	0.41
1:A:233:ASP:OD1	1:A:235:TRP:HB2	2.21	0.41
1:B:466:GLU:O	1:B:470:GLU:HG3	2.21	0.41
1:A:271:LYS:HB3	1:A:313:ILE:HB	2.03	0.40
1:B:272:THR:HG22	1:B:273:LEU:N	2.37	0.40
1:A:239:ARG:CB	1:A:239:ARG:NH1	2.84	0.40
1:A:360:ILE:O	1:A:360:ILE:HG13	2.19	0.40
1:B:238:GLU:HG3	1:B:240:THR:CB	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:297:ASN:ND2	1:B:346:GLN:HB3	2.36	0.40
1:B:323:LEU:O	1:B:327:LEU:HG	2.22	0.40
1:B:356:LYS:HE3	1:B:356:LYS:HB2	1.85	0.40
1:B:430:TRP:O	1:B:434:THR:HG23	2.22	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	269/293 (92%)	253 (94%)	15 (6%)	1 (0%)	38	63
1	B	270/293 (92%)	255 (94%)	12 (4%)	3 (1%)	17	35
All	All	539/586 (92%)	508 (94%)	27 (5%)	4 (1%)	25	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	276	ASP
1	B	277	THR
1	A	262	LYS
1	B	274	LYS

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	237/255 (93%)	227 (96%)	10 (4%)	34	62
1	B	238/255 (93%)	225 (94%)	13 (6%)	25	49
All	All	475/510 (93%)	452 (95%)	23 (5%)	30	55

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	234	LYS
1	A	241	ASP
1	A	266	LEU
1	A	282	GLU
1	A	308	GLU
1	A	354	LEU
1	A	370	LEU
1	A	376	LEU
1	A	401	PHE
1	A	447	GLN
1	B	234	LYS
1	B	266	LEU
1	B	275	GLU
1	B	276	ASP
1	B	282	GLU
1	B	308	GLU
1	B	354	LEU
1	B	370	LEU
1	B	376	LEU
1	B	381	ASP
1	B	401	PHE
1	B	459	GLU
1	B	460	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	447	GLN
1	B	498	GLN

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 ⓘ

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	$\# Z > 2$
3	P17	A	119	-	31,32,32	2.15	11 (35%)	40,46,46	2.40	12 (30%)
2	MES	B	117	-	12,12,12	3.27	6 (50%)	14,16,16	2.35	4 (28%)
2	MES	B	118	-	12,12,12	3.31	6 (50%)	14,16,16	2.40	6 (42%)
3	P17	B	120	-	31,32,32	2.11	9 (29%)	40,46,46	2.38	14 (35%)

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
3	P17	A	119	-	-	0/10/10/10	0/4/4/4
2	MES	B	117	-	-	0/6/14/14	0/1/1/1
2	MES	B	118	-	-	0/6/14/14	0/1/1/1
3	P17	B	120	-	-	0/10/10/10	0/4/4/4

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	120	P17	C18-N19	-2.90	1.34	1.38
3	A	119	P17	C18-N19	-2.75	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	120	P17	C4-C3	2.03	1.43	1.39
3	A	119	P17	C5-C6	2.04	1.42	1.38
3	B	120	P17	C25-C26	2.07	1.42	1.38
3	A	119	P17	C8-C3	2.09	1.43	1.39
3	A	119	P17	C12-N11	2.16	1.37	1.32
3	A	119	P17	C4-C3	2.20	1.43	1.39
3	A	119	P17	C10-N11	2.21	1.37	1.34
3	B	120	P17	C5-C6	2.55	1.43	1.38
3	B	120	P17	C18-C17	2.81	1.50	1.44
3	A	119	P17	C18-C17	2.86	1.50	1.44
2	B	118	MES	C7-N4	3.04	1.54	1.47
3	A	119	P17	C6-C7	3.31	1.44	1.39
2	B	117	MES	C5-N4	3.36	1.56	1.47
2	B	117	MES	C7-N4	3.41	1.55	1.47
2	B	118	MES	O3S-S	3.43	1.59	1.47
2	B	118	MES	C3-N4	3.57	1.56	1.47
2	B	118	MES	C5-N4	3.67	1.56	1.47
2	B	117	MES	O3S-S	3.72	1.60	1.47
3	B	120	P17	C6-C7	3.74	1.45	1.39
2	B	117	MES	C3-N4	3.84	1.57	1.47
2	B	117	MES	O1S-S	4.02	1.56	1.45
3	B	120	P17	C22-C27	4.07	1.44	1.40
3	B	120	P17	C16-C13	4.08	1.51	1.42
2	B	118	MES	O1S-S	4.14	1.57	1.45
3	A	119	P17	C22-C27	4.38	1.44	1.40
3	A	119	P17	C16-C13	4.46	1.52	1.42
3	A	119	P17	C22-C23	5.15	1.45	1.40
3	B	120	P17	C22-C23	5.76	1.46	1.40
2	B	117	MES	O2S-S	7.17	1.66	1.45
2	B	118	MES	O2S-S	7.44	1.66	1.45

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	119	P17	N11-C10-N15	-7.02	119.65	126.65
3	B	120	P17	N11-C10-N15	-6.23	120.43	126.65
3	B	120	P17	C13-C14-N15	-3.87	117.83	123.30
3	A	119	P17	C13-C14-N15	-3.78	117.96	123.30
3	B	120	P17	C17-C22-C23	-3.60	117.98	121.93
3	B	120	P17	C26-C27-C22	-2.98	119.33	122.38
3	A	119	P17	C24-C23-C22	-2.78	119.53	122.38
3	B	120	P17	C24-C23-C22	-2.71	119.61	122.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	119	P17	C17-C22-C23	-2.56	119.12	121.93
3	B	120	P17	C17-C16-C13	-2.53	118.23	121.92
3	A	119	P17	C26-C27-C22	-2.51	119.81	122.38
3	A	119	P17	C17-C16-C13	-2.50	118.28	121.92
3	B	120	P17	C13-C12-N11	-2.05	119.63	123.86
2	B	118	MES	O2S-S-O1S	-2.01	106.90	113.86
3	B	120	P17	C7-C8-C3	2.01	121.25	119.57
3	A	119	P17	C16-C17-C18	2.15	119.78	117.67
2	B	117	MES	C6-O1-C2	2.17	117.23	109.89
2	B	118	MES	C6-O1-C2	2.25	117.51	109.89
2	B	118	MES	C5-N4-C3	2.35	114.20	108.87
3	B	120	P17	C22-C23-CL1	2.41	123.08	119.72
3	A	119	P17	C22-C23-CL1	2.49	123.20	119.72
3	B	120	P17	C16-C17-C18	2.52	120.14	117.67
3	B	120	P17	C12-N11-C10	2.74	120.34	115.89
3	A	119	P17	C12-N11-C10	2.88	120.57	115.89
2	B	117	MES	O1S-S-C8	2.93	109.31	106.79
3	A	119	P17	C23-C22-C27	2.99	119.38	115.93
3	B	120	P17	C23-C22-C27	3.29	119.73	115.93
2	B	118	MES	O1S-S-C8	3.63	109.91	106.79
2	B	118	MES	O3S-S-C8	3.78	110.70	106.06
2	B	117	MES	O3S-S-C8	4.05	111.04	106.06
2	B	118	MES	O2S-S-C8	5.96	111.91	106.79
2	B	117	MES	O2S-S-C8	5.96	111.91	106.79
3	B	120	P17	C17-C18-N19	6.03	120.29	116.11
3	A	119	P17	C17-C18-N19	6.53	120.64	116.11
3	B	120	P17	C10-N15-C14	6.56	122.59	115.11
3	A	119	P17	C10-N15-C14	6.78	122.84	115.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	119	P17	2	0
3	B	120	P17	2	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	271/293 (92%)	-0.03	11 (4%) 38 30	15, 39, 84, 115	0
1	B	272/293 (92%)	-0.36	2 (0%) 87 85	14, 29, 69, 100	0
All	All	543/586 (92%)	-0.20	13 (2%) 59 52	14, 34, 76, 115	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	237	MET	4.7
1	B	274	LYS	3.4
1	B	277	THR	3.3
1	A	261	TRP	3.3
1	A	264	TYR	3.2
1	A	263	LYS	3.1
1	A	232	TYR	2.8
1	A	245	LYS	2.5
1	A	246	HIS	2.5
1	A	240	THR	2.4
1	A	266	LEU	2.4
1	A	276	ASP	2.1
1	A	307	ARG	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(Å ²)	Q<0.9
2	MES	B	117	12/12	0.95	0.27	2.57	52,53,55,56	0
2	MES	B	118	12/12	0.92	0.20	1.33	50,52,52,53	0
3	P17	A	119	29/29	0.91	0.18	0.64	38,41,46,50	0
3	P17	B	120	29/29	0.94	0.14	-0.23	25,28,32,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.