Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:15 AM GMT

PDB ID	:	3DYA
Title	:	HIV-1 RT with non-nucleoside inhibitor annulated Pyrazole 1
Authors	:	Harris, S.F.; Villasenor, A.
Deposited on	:	2008-07-25
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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

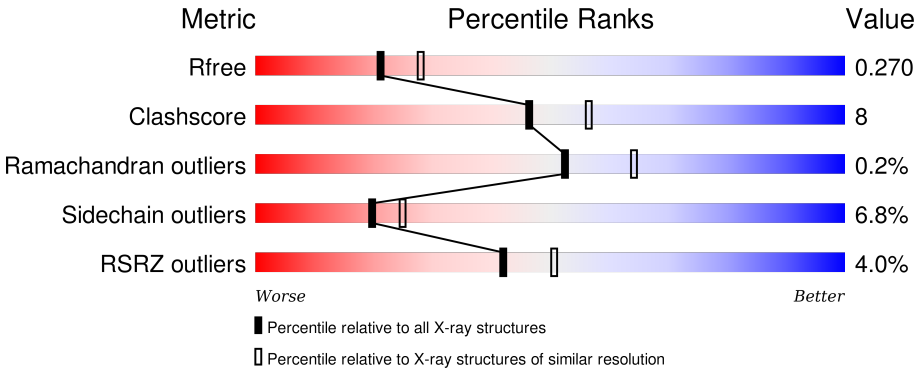
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 <=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	561	3%, 80%, 15%
2	B	440	5%, 73%, 16%, 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8161 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.

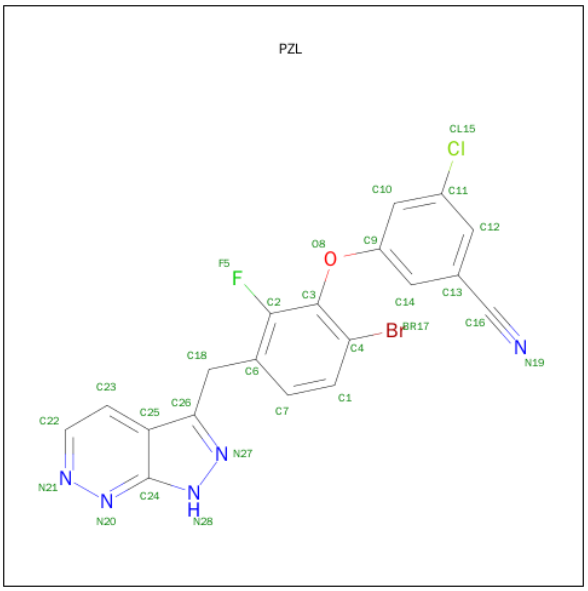
- Molecule 1 is a protein called REVERSE TRANSCRIPTASE/RIBONUCLEASE H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	551	4495	2906	751	830	8	0	1	0

- Molecule 2 is a protein called p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	402	3327	2170	546	604	7	0	0	0

- Molecule 3 is 3-[6-BROMO-2-FLUORO-3-(1H-PYRAZOLO[3,4-C]PYRIDAZIN-3-YL METHYL)PHENOXY]-5-CHLOROBENZONITRILE (three-letter code: PZL) (formula: $C_{19}H_{10}BrClFN_5O$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	Br	C	Cl	F	N	O		
3	A	1	28	1	19	1	1	5	1	0	0

- Molecule 4 is water.

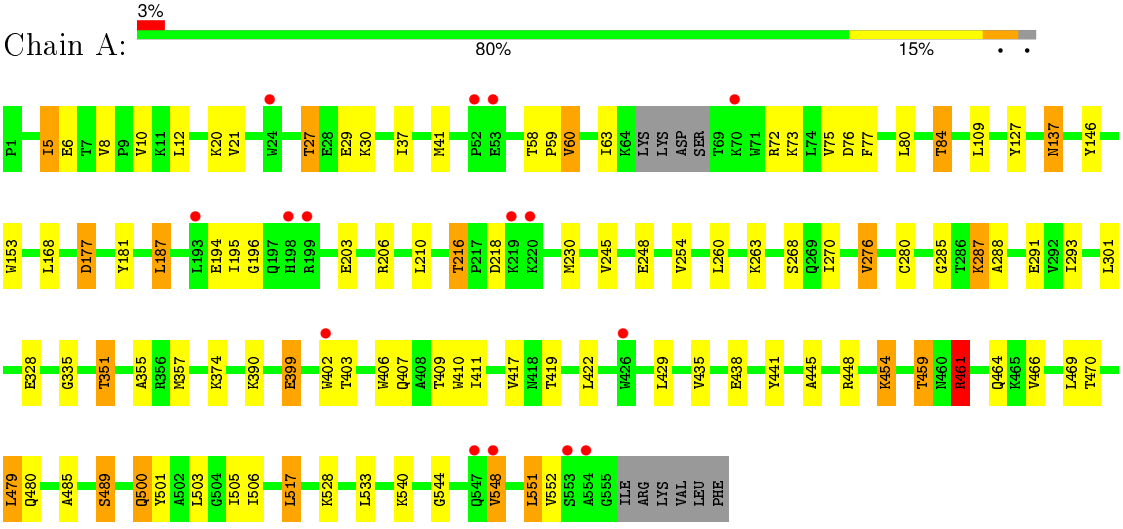
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	183	Total O 183 183	0	0
4	B	128	Total O 128 128	0	0

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 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.

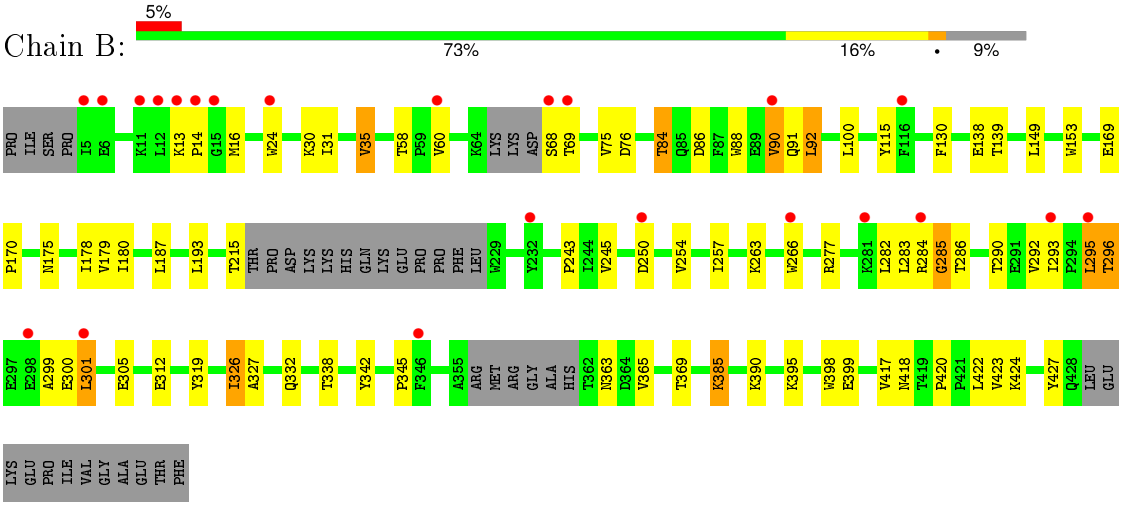
- Molecule 1: REVERSE TRANSCRIPTASE/RIBONUCLEASE H

Chain A:



- Molecule 2: p51 RT

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	117.66Å 153.67Å 155.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.98 – 2.30 46.97 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.6 (46.98-2.30) 96.6 (46.97-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	1.39 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.218 , 0.266 0.222 , 0.270	Depositor DCC
R_{free} test set	3093 reflections (5.37%)	DCC
Wilson B-factor (Å^2)	52.5	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 46.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Outliers	0 of 60666 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8161	wwPDB-VP
Average B, all atoms (Å^2)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: PZL

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	0.59	0/4612	0.65	1/6267 (0.0%)
2	B	0.60	0/3421	0.63	0/4650
All	All	0.59	0/8033	0.64	1/10917 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	461	ARG	NE-CZ-NH1	7.66	124.13	120.30

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	4495	0	4536	69	0
2	B	3327	0	3348	61	0
3	A	28	0	10	0	0
4	A	183	0	0	5	0
4	B	128	0	0	4	0
All	All	8161	0	7894	120	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 8.

All (120) 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:500:GLN:NE2	2:B:422:LEU:HD12	1.66	1.08
1:A:268:SER:O	1:A:351:THR:HG22	1.71	0.89
2:B:84:THR:HG21	2:B:153:TRP:HE1	1.38	0.88
1:A:27:THR:HG21	4:A:660:HOH:O	1.75	0.85
1:A:80:LEU:O	1:A:84:THR:HG22	1.76	0.85
2:B:326:ILE:HD13	2:B:390:LYS:HD2	1.58	0.85
1:A:37:ILE:HG22	1:A:41:MET:CE	2.11	0.80
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.49	0.77
1:A:37:ILE:HG22	1:A:41:MET:HE1	1.66	0.76
1:A:461:ARG:HH11	1:A:461:ARG:HG2	1.50	0.76
2:B:254:VAL:HG13	2:B:283:LEU:HD22	1.68	0.75
2:B:138:GLU:HG2	2:B:139:THR:HG23	1.71	0.73
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.24	0.72
1:A:406:TRP:CZ3	1:A:407:GLN:HG2	2.25	0.71
2:B:365:VAL:O	2:B:369:THR:HG23	1.92	0.69
1:A:216:THR:HG22	4:A:563:HOH:O	1.92	0.68
1:A:466:VAL:CG2	1:A:551:LEU:HD12	2.24	0.68
2:B:24:TRP:HB2	4:B:443:HOH:O	1.94	0.67
1:A:485:ALA:O	1:A:489:SER:HB2	1.96	0.66
1:A:417:VAL:HG22	1:A:419:THR:HG23	1.77	0.66
1:A:435:VAL:HG23	2:B:290:THR:HG21	1.78	0.65
1:A:73:LYS:NZ	1:A:146:TYR:OH	2.30	0.65
2:B:296:THR:HG23	2:B:299:ALA:CB	2.27	0.64
1:A:27:THR:HG22	1:A:30:LYS:H	1.62	0.63
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.35	0.61
2:B:91:GLN:HG2	2:B:92:LEU:HD13	1.82	0.61
2:B:295:LEU:HD12	2:B:300:GLU:OE1	1.99	0.61
2:B:326:ILE:HD12	2:B:327:ALA:N	2.15	0.60
1:A:409:THR:HG21	4:A:567:HOH:O	2.01	0.60
1:A:254:VAL:HG22	1:A:293:ILE:HD11	1.82	0.60
2:B:424:LYS:NZ	4:B:485:HOH:O	2.34	0.60
1:A:454:LYS:HE3	1:A:466:VAL:HG11	1.82	0.59
1:A:37:ILE:HG22	1:A:41:MET:HE2	1.83	0.59
1:A:466:VAL:HG23	1:A:551:LEU:HD12	1.85	0.58
2:B:254:VAL:HG13	2:B:283:LEU:CD2	2.33	0.58
2:B:86:ASP:HA	2:B:90:VAL:HG13	1.85	0.58
2:B:301:LEU:HD22	4:B:482:HOH:O	2.02	0.58
2:B:58:THR:HG23	2:B:76:ASP:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:266:TRP:HZ2	2:B:427:TYR:HH	1.53	0.56
2:B:332:GLN:HG3	2:B:338:THR:HG23	1.88	0.56
1:A:12:LEU:HD11	1:A:127:TYR:CE1	2.41	0.56
1:A:503:LEU:HD12	1:A:533:LEU:CD2	2.36	0.55
1:A:406:TRP:CZ3	1:A:407:GLN:CG	2.90	0.54
2:B:84:THR:HG23	2:B:88:TRP:HD1	1.73	0.54
2:B:243:PRO:O	2:B:245:VAL:HG13	2.07	0.54
2:B:296:THR:HG23	2:B:299:ALA:HB2	1.89	0.53
2:B:266:TRP:HZ2	2:B:427:TYR:OH	1.92	0.53
2:B:420:PRO:HB2	2:B:423:VAL:HG13	1.90	0.53
1:A:60:VAL:HG12	1:A:75:VAL:HG22	1.90	0.52
1:A:276:VAL:HG13	1:A:280:CYS:SG	2.50	0.52
1:A:8:VAL:O	1:A:10:VAL:HG23	2.10	0.52
1:A:459:THR:O	2:B:286:THR:HG21	2.09	0.52
1:A:37:ILE:CG2	1:A:41:MET:HE1	2.39	0.52
2:B:266:TRP:CZ2	2:B:427:TYR:CZ	2.97	0.52
1:A:429:LEU:HD11	1:A:506:ILE:HG22	1.91	0.51
2:B:84:THR:HG23	2:B:88:TRP:CD1	2.46	0.51
2:B:301:LEU:HD12	2:B:305:GLU:OE2	2.11	0.51
2:B:296:THR:HG23	2:B:299:ALA:HB3	1.93	0.50
1:A:84:THR:HG21	1:A:153:TRP:NE1	2.24	0.50
1:A:245:VAL:O	1:A:263:LYS:NZ	2.44	0.50
2:B:68:SER:OG	2:B:69:THR:N	2.44	0.50
1:A:438:GLU:HG3	1:A:461:ARG:HD3	1.94	0.48
2:B:395:LYS:NZ	2:B:399:GLU:OE1	2.46	0.48
1:A:435:VAL:CG2	2:B:290:THR:HG21	2.43	0.48
1:A:469:LEU:HD21	1:A:480:GLN:HG2	1.95	0.48
1:A:75:VAL:HG11	1:A:77:PHE:CZ	2.49	0.48
1:A:27:THR:CG2	1:A:30:LYS:H	2.25	0.47
1:A:137:ASN:ND2	4:A:585:HOH:O	2.46	0.47
2:B:326:ILE:CD1	2:B:390:LYS:HD2	2.38	0.47
1:A:445:ALA:HB3	1:A:552:VAL:HG12	1.96	0.47
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.97	0.46
1:A:503:LEU:HD12	1:A:533:LEU:HD23	1.98	0.45
1:A:410:TRP:CE3	2:B:363:ASN:HB2	2.51	0.45
1:A:287:LYS:HB3	1:A:291:GLU:OE1	2.17	0.45
1:A:177:ASP:OD1	1:A:177:ASP:C	2.55	0.45
1:A:181:TYR:CD1	2:B:138:GLU:HA	2.52	0.45
1:A:168:LEU:HD11	1:A:187:LEU:HD11	1.99	0.45
1:A:441:TYR:O	1:A:548:VAL:HG21	2.17	0.44
2:B:282:LEU:HD11	2:B:296:THR:CG2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:GLY:O	1:A:355:ALA:HA	2.17	0.44
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.99	0.44
1:A:328:GLU:HG2	1:A:390:LYS:HB2	2.00	0.44
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.99	0.44
1:A:195:ILE:CG2	1:A:196:GLY:N	2.80	0.44
1:A:406:TRP:HH2	2:B:418:ASN:HA	1.80	0.43
1:A:270:ILE:HA	1:A:351:THR:HB	2.00	0.43
1:A:399:GLU:O	1:A:403:THR:HG23	2.18	0.43
1:A:500:GLN:HE21	2:B:422:LEU:HD12	1.71	0.43
1:A:58:THR:HG23	1:A:76:ASP:O	2.19	0.43
1:A:5:ILE:HD13	1:A:6:GLU:O	2.18	0.43
1:A:206:ARG:HG2	1:A:216:THR:HG21	2.01	0.43
1:A:459:THR:O	2:B:286:THR:CG2	2.67	0.43
2:B:257:ILE:HD12	2:B:293:ILE:HG21	2.01	0.42
2:B:31:ILE:O	2:B:35:VAL:HG13	2.18	0.42
1:A:109:LEU:HD22	1:A:216:THR:OG1	2.19	0.42
2:B:284:ARG:CG	2:B:285:GLY:N	2.83	0.42
2:B:342:TYR:CE1	2:B:345:PRO:HA	2.55	0.42
2:B:169:GLU:HB3	2:B:170:PRO:HD3	2.01	0.42
2:B:60:VAL:HG11	2:B:130:PHE:CD2	2.55	0.42
2:B:180:ILE:CG2	2:B:187:LEU:HD22	2.49	0.42
1:A:288:ALA:HB3	1:A:291:GLU:HB2	2.01	0.42
1:A:501:TYR:CZ	1:A:505:ILE:HD11	2.54	0.42
2:B:13:LYS:HG2	2:B:14:PRO:CD	2.48	0.42
1:A:27:THR:HG23	1:A:29:GLU:H	1.85	0.42
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.55	0.41
2:B:100:LEU:CD1	2:B:179:VAL:HG23	2.50	0.41
2:B:30:LYS:NZ	4:B:536:HOH:O	2.52	0.41
2:B:295:LEU:HD23	2:B:295:LEU:H	1.85	0.41
1:A:479:LEU:HB3	1:A:517:LEU:HD13	2.02	0.41
2:B:282:LEU:HB3	2:B:293:ILE:HD13	2.03	0.41
2:B:266:TRP:CH2	2:B:427:TYR:CZ	3.08	0.41
2:B:369:THR:HG22	2:B:398:TRP:CZ3	2.55	0.40
1:A:407:GLN:HB3	1:A:407:GLN:HE21	1.69	0.40
1:A:544:GLY:O	1:A:548:VAL:HG23	2.21	0.40
1:A:407:GLN:NE2	4:A:623:HOH:O	2.54	0.40
1:A:489:SER:HB3	1:A:528:LYS:NZ	2.37	0.40
2:B:282:LEU:HD11	2:B:296:THR:HG22	2.04	0.40
2:B:266:TRP:HH2	2:B:427:TYR:CE1	2.40	0.40
2:B:319:TYR:OH	2:B:385:LYS:HE3	2.22	0.40
2:B:175:ASN:HB3	2:B:178:ILE:HD12	2.04	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	548/561 (98%)	528 (96%)	19 (4%)	1 (0%)	52 64
2	B	394/440 (90%)	381 (97%)	12 (3%)	1 (0%)	46 57
All	All	942/1001 (94%)	909 (96%)	31 (3%)	2 (0%)	52 64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	285	GLY
2	B	285	GLY

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	492/501 (98%)	452 (92%)	40 (8%)	15 18
2	B	366/400 (92%)	348 (95%)	18 (5%)	31 41
All	All	858/901 (95%)	800 (93%)	58 (7%)	20 25

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE

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Mol	Chain	Res	Type
1	A	20	LYS
1	A	27	THR
1	A	60	VAL
1	A	63	ILE
1	A	72	ARG
1	A	84	THR
1	A	137	ASN
1	A	177	ASP
1	A	187	LEU
1	A	194	GLU
1	A	203	GLU
1	A	210	LEU
1	A	216	THR
1	A	218	ASP
1	A	230	MET
1	A	248	GLU
1	A	260	LEU
1	A	276	VAL
1	A	287	LYS
1	A	301	LEU
1	A	351	THR
1	A	357	MET
1	A	374	LYS
1	A	399	GLU
1	A	411	ILE
1	A	422	LEU
1	A	448	ARG
1	A	454	LYS
1	A	459	THR
1	A	461	ARG
1	A	464	GLN
1	A	470	THR
1	A	479	LEU
1	A	489	SER
1	A	500	GLN
1	A	517	LEU
1	A	540	LYS
1	A	548	VAL
1	A	551	LEU
2	B	16	MET
2	B	35	VAL
2	B	84	THR

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Mol	Chain	Res	Type
2	B	90	VAL
2	B	92	LEU
2	B	193	LEU
2	B	215	THR
2	B	250	ASP
2	B	263	LYS
2	B	277	ARG
2	B	292	VAL
2	B	295	LEU
2	B	296	THR
2	B	301	LEU
2	B	312	GLU
2	B	326	ILE
2	B	385	LYS
2	B	417	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	198	HIS
1	A	221	HIS
1	A	278	GLN
1	A	336	GLN
1	A	407	GLN
1	A	428	GLN
1	A	500	GLN
1	A	507	GLN
1	A	524	GLN
2	B	151	GLN
2	B	208	HIS
2	B	278	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 (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	$\#\|Z\| > 2$
3	PZL	A	562	-	30,31,31	2.00	4 (13%)	31,44,44	2.53	10 (32%)

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	PZL	A	562	-	-	0/10/10/10	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	562	PZL	C13-C16	-6.61	1.27	1.44
3	A	562	PZL	C3-C4	-2.41	1.35	1.40
3	A	562	PZL	C23-C22	2.87	1.40	1.36
3	A	562	PZL	N21-N20	6.04	1.42	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	562	PZL	C6-C18-C26	-8.11	101.37	114.85
3	A	562	PZL	O8-C3-C4	-2.95	115.83	120.59
3	A	562	PZL	C13-C16-N19	-2.82	171.17	177.99
3	A	562	PZL	BR17-C4-C3	-2.82	114.19	118.53
3	A	562	PZL	C2-C3-C4	2.06	120.89	116.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	562	PZL	C14-C13-C16	2.49	122.66	119.51
3	A	562	PZL	C9-O8-C3	2.92	123.55	118.46
3	A	562	PZL	BR17-C4-C1	3.34	124.46	117.81
3	A	562	PZL	C7-C6-C2	5.30	119.56	116.42
3	A	562	PZL	F5-C2-C6	5.51	123.11	117.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	551/561 (98%)	0.32	15 (2%)	58	67	39, 56, 73, 95	0
2	B	402/440 (91%)	0.49	23 (5%)	27	36	39, 54, 85, 90	0
All	All	953/1001 (95%)	0.39	38 (3%)	42	51	39, 55, 81, 95	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	14	PRO	7.2
2	B	232	TYR	5.4
2	B	5	ILE	4.6
2	B	301	LEU	4.2
2	B	13	LYS	4.1
2	B	15	GLY	3.9
2	B	12	LEU	3.7
2	B	69	THR	3.7
2	B	284	ARG	3.6
2	B	24	TRP	3.3
2	B	6	GLU	3.3
2	B	295	LEU	3.3
2	B	116	PHE	3.1
2	B	250	ASP	3.1
1	A	52	PRO	2.9
1	A	426	TRP	2.8
2	B	346	PHE	2.8
1	A	547	GLN	2.8
2	B	266	TRP	2.8
2	B	11	LYS	2.7
1	A	193	LEU	2.7
1	A	198	HIS	2.6
1	A	548	VAL	2.5
1	A	24	TRP	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	293	ILE	2.5
2	B	68	SER	2.4
2	B	90	VAL	2.4
1	A	402	TRP	2.4
2	B	298	GLU	2.3
1	A	554	ALA	2.3
2	B	281	LYS	2.2
1	A	553	SER	2.2
1	A	219	LYS	2.2
2	B	60	VAL	2.1
1	A	53	GLU	2.1
1	A	220	LYS	2.1
1	A	70	LYS	2.0
1	A	199	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, 95^{th} 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
3	PZL	A	562	28/28	0.97	0.13	-0.54	46,48,52,53	0

6.5 Other polymers (i)

There are no such residues in this entry.