Full wwPDB X-ray Structure Validation Report

Oct 17, 2017 – 03:17 AM EDT

PDB ID	:	3EKT
Title	:	Crystal Structure of the inhibitor Darunavir (DRV) in complex with a multi-drug resistant HIV-1 protease variant (L10F/G48V/I54V/V64I/V82A) (Refer: FLAP+ in citation.)
Authors	:	Prabu-Jeyabalan, M.; King, N.M.; Bandaranayake, R.M.
Deposited on	:	unknown
Resolution	:	1.97 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:

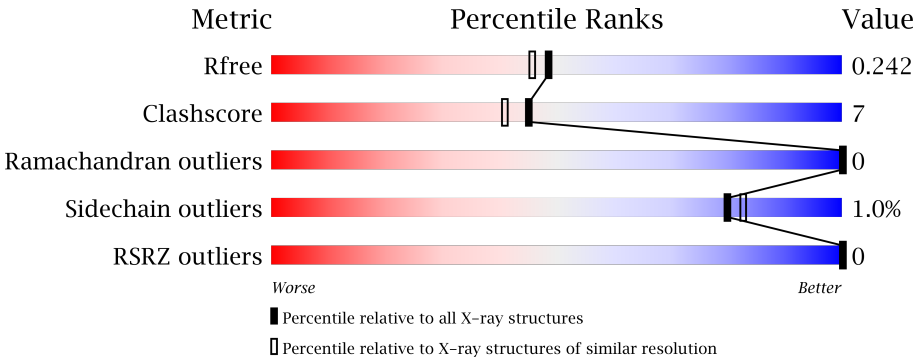
MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

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 1.97 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	9293 (2.00-1.96)
Clashscore	112137	10621 (2.00-1.96)
Ramachandran outliers	110173	10502 (2.00-1.96)
Sidechain outliers	110143	10501 (2.00-1.96)
RSRZ outliers	101464	9395 (2.00-1.96)

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	99	97%
1	B	99	94% 6%
1	C	99	87% 13%
1	D	99	79% 21%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	608	-	-	-	X
2	ACT	C	603	-	-	-	X
2	ACT	C	604	-	-	-	X
2	ACT	D	602	-	-	-	X
2	ACT	D	605	-	-	X	X
4	PO4	C	505	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3555 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 Protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	3	0
			761	492	128	137	4			
1	B	99	Total	C	N	O	S	0	3	0
			753	486	126	137	4			
1	C	99	Total	C	N	O	S	0	5	0
			767	493	130	138	6			
1	D	99	Total	C	N	O	S	0	6	0
			778	500	132	140	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	LYS	GLN	ENGINEERED	UNP P03369
A	10	ILE	LEU	ENGINEERED	UNP P03369
A	48	VAL	GLY	ENGINEERED	UNP P03369
A	54	VAL	ILE	ENGINEERED	UNP P03369
A	64	ILE	VAL	ENGINEERED	UNP P03369
A	82	ALA	VAL	ENGINEERED	UNP P03369
B	7	LYS	GLN	ENGINEERED	UNP P03369
B	10	ILE	LEU	ENGINEERED	UNP P03369
B	48	VAL	GLY	ENGINEERED	UNP P03369
B	54	VAL	ILE	ENGINEERED	UNP P03369
B	64	ILE	VAL	ENGINEERED	UNP P03369
B	82	ALA	VAL	ENGINEERED	UNP P03369
C	7	LYS	GLN	ENGINEERED	UNP P03369
C	10	ILE	LEU	ENGINEERED	UNP P03369
C	48	VAL	GLY	ENGINEERED	UNP P03369
C	54	VAL	ILE	ENGINEERED	UNP P03369
C	64	ILE	VAL	ENGINEERED	UNP P03369
C	82	ALA	VAL	ENGINEERED	UNP P03369
D	7	LYS	GLN	ENGINEERED	UNP P03369
D	10	ILE	LEU	ENGINEERED	UNP P03369
D	48	VAL	GLY	ENGINEERED	UNP P03369

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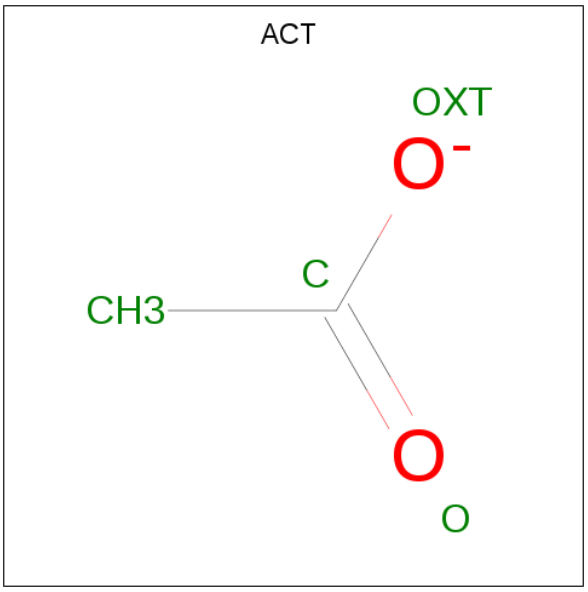
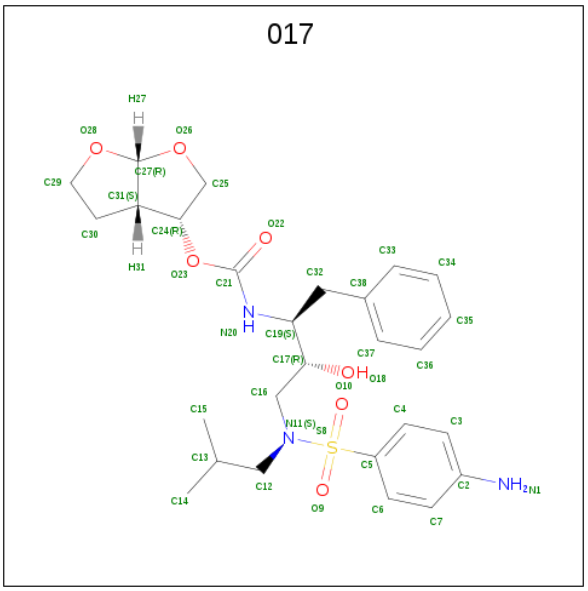
Chain	Residue	Modelled	Actual	Comment	Reference
D	54	VAL	ILE	ENGINEERED	UNP P03369
D	64	ILE	VAL	ENGINEERED	UNP P03369
D	82	ALA	VAL	ENGINEERED	UNP P03369

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



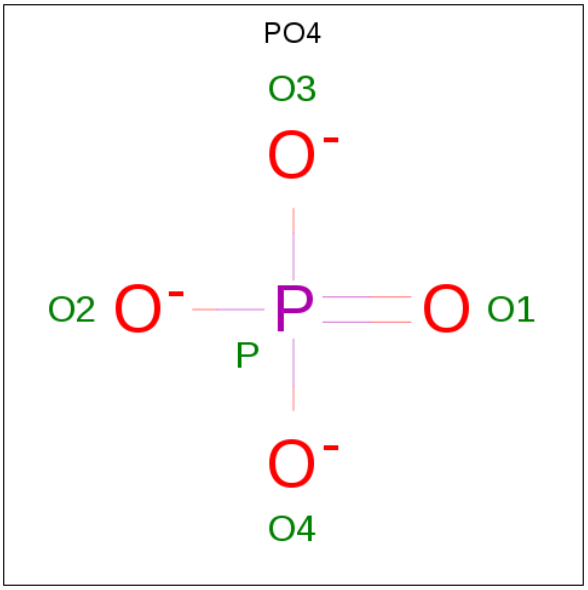
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0

- Molecule 3 is (3R,3AS,6AR)-HEXAHYDROFURO[2,3-B]FURAN-3-YL(1S,2R)-3-[[(4-AMINOPHENYL)SULFONYL](ISOBUTYL)AMINO]-1-BENZYL-2-HYDROXYPROPYLCARBAMATE (three-letter code: 017) (formula: $C_{27}H_{37}N_3O_7S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	S	0	0
			38	27	3	7	1		
3	D	1	Total	C	N	O	S	0	1
			76	54	6	14	2		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0

- Molecule 5 is water.

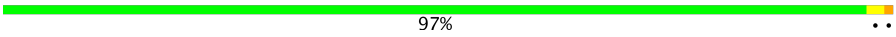
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	71	Total O 72 72	1	1
5	B	89	Total O 89 89	0	0
5	C	79	Total O 80 80	0	1
5	D	82	Total O 84 84	0	2

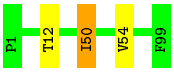
3 Residue-property plots (i)

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: Protease

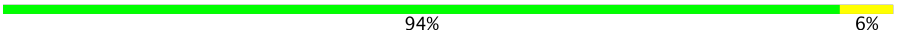
Chain A:





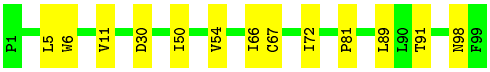
- Molecule 1: Protease

Chain B:



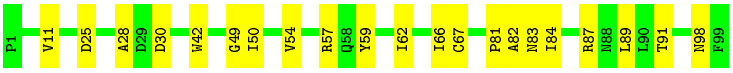
- Molecule 1: Protease

Chain C:



- Molecule 1: Protease

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α, β, γ	92.01Å 92.01Å 106.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 1.97 23.75 – 1.75	Depositor EDS
% Data completeness (in resolution range)	95.6 (40.00-1.97) 94.6 (23.75-1.75)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.55 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.209 , 0.258 0.223 , 0.242	Depositor DCC
R_{free} test set	1713 reflections (5.24%)	DCC
Wilson B-factor (Å^2)	17.5	Xtriage
Anisotropy	0.218	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 31.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.36$	Xtriage
Estimated twinning fraction	0.479 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3555	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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, 0.333 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: PO4, ACT, 017

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	0/775	0.51	0/1055
1	B	0.35	0/767	0.54	0/1047
1	C	0.39	0/781	0.53	0/1065
1	D	0.38	0/792	0.54	0/1079
All	All	0.37	0/3115	0.53	0/4246

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	761	0	787	5	0
1	B	753	0	765	12	0
1	C	767	0	778	13	0
1	D	778	0	795	27	0
2	A	8	0	6	1	0
2	B	4	0	3	1	0
2	C	8	0	6	2	0
2	D	12	0	9	5	0
3	B	38	0	37	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	76	0	74	14	0
4	C	15	0	0	0	0
4	D	10	0	0	0	0
5	A	72	0	0	0	0
5	B	89	0	0	0	0
5	C	80	0	0	1	0
5	D	84	0	0	1	0
All	All	3555	0	3260	48	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 7.

All (48) 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:D:83:ASN:H	2:D:605:ACT:H3	1.26	0.97
1:D:84[A]:ILE:CD1	3:D:200[A]:017:H6	2.12	0.80
1:D:83:ASN:H	2:D:605:ACT:CH3	1.98	0.77
1:B:28:ALA:CB	1:B:84[B]:ILE:HD11	2.21	0.71
1:D:84[A]:ILE:HD12	3:D:200[A]:017:H6	1.74	0.69
1:B:28:ALA:CB	1:B:84[B]:ILE:CD1	2.71	0.69
1:A:50[A]:ILE:HG23	1:B:54:VAL:HG21	1.76	0.67
1:D:87:ARG:O	1:D:91[B]:THR:HG23	1.97	0.65
1:C:72:ILE:H	2:C:604:ACT:H3	1.64	0.62
1:D:82:ALA:HA	2:D:605:ACT:H2	1.82	0.61
1:B:28:ALA:HB2	1:B:84[B]:ILE:CD1	2.30	0.61
1:B:84[A]:ILE:CD1	3:B:200:017:H6	2.31	0.59
1:B:28:ALA:HB1	1:B:84[B]:ILE:HD11	1.84	0.58
1:B:68:GLY:H	2:B:607:ACT:H1	1.70	0.55
1:D:81:PRO:HG2	3:D:200[A]:017:C35	2.38	0.54
1:D:83:ASN:N	2:D:605:ACT:H3	2.10	0.52
1:C:11:VAL:HB	1:C:67[A]:CYS:SG	2.50	0.52
1:C:81:PRO:HG2	3:D:200[B]:017:C35	2.40	0.51
1:C:50[B]:ILE:HG23	1:D:54:VAL:HG11	1.93	0.50
1:C:98:ASN:HD22	1:D:98:ASN:HD22	1.60	0.50
1:D:25:ASP:CG	3:D:200[A]:017:H18	2.14	0.49
1:D:84[B]:ILE:HD11	3:D:200[B]:017:H141	1.93	0.49
1:B:84[A]:ILE:HD12	3:B:200:017:H6	1.94	0.49
1:D:66:ILE:HD11	1:D:89:LEU:HD13	1.96	0.48
1:C:66:ILE:HD11	1:C:89:LEU:HD13	1.95	0.48
1:B:28:ALA:CB	1:B:84[B]:ILE:HD12	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:42:TRP:CE2	1:D:57:ARG:HD2	2.49	0.47
1:D:28:ALA:HB1	3:D:200[A]:017:C7	2.45	0.47
1:A:12:THR:HG22	2:A:606:ACT:H2	1.97	0.47
1:C:50[A]:ILE:HG22	1:D:49:GLY:HA3	1.97	0.47
1:D:84[A]:ILE:HD11	3:D:200[A]:017:H6	1.95	0.47
1:D:81:PRO:HG2	3:D:200[A]:017:H35	1.97	0.46
2:D:605:ACT:H1	5:D:661:HOH:O	2.15	0.46
1:D:59:TYR:HB3	1:D:62:ILE:HD11	1.98	0.45
1:C:5:LEU:C	1:D:91[B]:THR:HG22	2.37	0.45
1:C:6:TRP:HA	1:D:91[B]:THR:HG21	1.98	0.45
1:A:50[B]:ILE:HG23	1:B:54:VAL:HG21	1.99	0.43
1:A:54:VAL:HG21	1:B:50[B]:ILE:HG23	2.00	0.43
1:D:84[A]:ILE:HD11	3:D:200[A]:017:C6	2.49	0.43
1:D:11:VAL:HB	1:D:67[A]:CYS:SG	2.59	0.42
1:D:84[A]:ILE:CD1	3:D:200[A]:017:C6	2.92	0.42
1:C:30:ASP:OD2	5:C:656:HOH:O	2.21	0.42
1:C:91:THR:O	2:C:603:ACT:H1	2.19	0.42
1:C:50[B]:ILE:HG23	1:D:54:VAL:CG1	2.49	0.42
1:D:30:ASP:HB2	3:D:200[A]:017:H12	1.84	0.42
1:D:50[B]:ILE:HD12	3:D:200[B]:017:O10	2.20	0.41
1:A:54:VAL:HG21	1:B:50[B]:ILE:CG2	2.51	0.41
1:C:81:PRO:HG2	3:D:200[B]:017:C36	2.51	0.41

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	100/99 (101%)	100 (100%)	0	0	100	100
1	B	100/99 (101%)	100 (100%)	0	0	100	100
1	C	102/99 (103%)	102 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	103/99 (104%)	103 (100%)	0	0	100	100
All	All	405/396 (102%)	405 (100%)	0	0	100	100

There are no Ramachandran outliers to report.

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	81/83 (98%)	79 (98%)	2 (2%)	53	51
1	B	79/83 (95%)	78 (99%)	1 (1%)	73	75
1	C	81/83 (98%)	80 (99%)	1 (1%)	75	78
1	D	83/83 (100%)	83 (100%)	0	100	100
All	All	324/332 (98%)	320 (99%)	4 (1%)	80	78

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50[A]	ILE
1	A	50[B]	ILE
1	B	46	MET
1	C	54	VAL

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	37	ASN
1	A	69	HIS
1	B	2	GLN
1	B	37	ASN
1	B	98	ASN
1	D	2	GLN
1	D	69	HIS

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Mol	Chain	Res	Type
1	D	98	ASN

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

16 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	ACT	A	606	-	1,3,3	0.59	0	0,3,3	0.00	-
2	ACT	A	608	-	1,3,3	1.43	0	0,3,3	0.00	-
3	017	B	200	-	41,41,41	2.26	5 (12%)	52,58,58	2.07	15 (28%)
2	ACT	B	607	-	1,3,3	0.02	0	0,3,3	0.00	-
4	PO4	C	502	-	4,4,4	0.68	0	6,6,6	0.38	0
4	PO4	C	503	-	4,4,4	0.82	0	6,6,6	0.46	0
4	PO4	C	505	-	4,4,4	0.77	0	6,6,6	0.33	0
2	ACT	C	603	-	1,3,3	1.87	0	0,3,3	0.00	-
2	ACT	C	604	-	1,3,3	1.22	0	0,3,3	0.00	-
3	017	D	200[A]	-	41,41,41	2.26	4 (9%)	52,58,58	1.76	11 (21%)
3	017	D	200[B]	-	41,41,41	2.26	4 (9%)	52,58,58	1.73	12 (23%)
4	PO4	D	501	-	4,4,4	0.81	0	6,6,6	0.56	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	PO4	D	504	-	4,4,4	0.80	0	6,6,6	0.41	0
2	ACT	D	601	-	1,3,3	1.24	0	0,3,3	0.00	-
2	ACT	D	602	-	1,3,3	0.97	0	0,3,3	0.00	-
2	ACT	D	605	-	1,3,3	1.17	0	0,3,3	0.00	-

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	ACT	A	606	-	-	0/0/0/0	0/0/0/0
2	ACT	A	608	-	-	0/0/0/0	0/0/0/0
3	017	B	200	-	-	0/36/55/55	0/4/4/4
2	ACT	B	607	-	-	0/0/0/0	0/0/0/0
4	PO4	C	502	-	-	0/0/0/0	0/0/0/0
4	PO4	C	503	-	-	0/0/0/0	0/0/0/0
4	PO4	C	505	-	-	0/0/0/0	0/0/0/0
2	ACT	C	603	-	-	0/0/0/0	0/0/0/0
2	ACT	C	604	-	-	0/0/0/0	0/0/0/0
3	017	D	200[A]	-	-	0/36/55/55	0/4/4/4
3	017	D	200[B]	-	-	0/36/55/55	0/4/4/4
4	PO4	D	501	-	-	0/0/0/0	0/0/0/0
4	PO4	D	504	-	-	0/0/0/0	0/0/0/0
2	ACT	D	601	-	-	0/0/0/0	0/0/0/0
2	ACT	D	602	-	-	0/0/0/0	0/0/0/0
2	ACT	D	605	-	-	0/0/0/0	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	200[A]	017	C5-S8	-10.94	1.60	1.76
3	D	200[B]	017	C5-S8	-10.83	1.60	1.76
3	B	200	017	C5-S8	-10.02	1.62	1.76
3	B	200	017	C7-C6	-4.97	1.29	1.38
3	D	200[A]	017	C7-C6	-4.94	1.29	1.38
3	B	200	017	C7-C2	-4.87	1.29	1.39
3	D	200[B]	017	C7-C6	-4.68	1.30	1.38
3	D	200[B]	017	C7-C2	-4.68	1.29	1.39
3	D	200[A]	017	C7-C2	-4.58	1.29	1.39
3	B	200	017	O23-C24	-2.22	1.42	1.46
3	D	200[A]	017	O23-C21	5.64	1.45	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	200[B]	017	O23-C21	5.94	1.45	1.35
3	B	200	017	O23-C21	6.43	1.46	1.35

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	200	017	O10-S8-O9	-7.20	107.36	119.46
3	D	200[A]	017	O10-S8-O9	-5.95	109.45	119.46
3	D	200[B]	017	O10-S8-O9	-5.70	109.88	119.46
3	B	200	017	C32-C19-N20	-3.88	105.62	110.17
3	D	200[B]	017	C32-C19-N20	-3.71	105.83	110.17
3	D	200[A]	017	C32-C19-N20	-3.39	106.20	110.17
3	B	200	017	C12-N11-S8	-3.11	110.73	117.75
3	B	200	017	C16-N11-S8	-3.06	110.84	117.75
3	D	200[A]	017	C12-N11-S8	-2.77	111.51	117.75
3	D	200[A]	017	C16-N11-S8	-2.75	111.54	117.75
3	D	200[B]	017	C16-N11-S8	-2.74	111.56	117.75
3	D	200[B]	017	O23-C21-O22	-2.65	120.47	124.51
3	D	200[A]	017	O23-C21-O22	-2.57	120.58	124.51
3	D	200[A]	017	C17-C19-N20	-2.43	105.19	109.52
3	D	200[B]	017	C12-N11-S8	-2.36	112.43	117.75
3	D	200[B]	017	C38-C32-C19	-2.23	109.31	113.36
3	B	200	017	C38-C32-C19	-2.13	109.48	113.36
3	B	200	017	C6-C5-C4	-2.06	117.48	120.43
3	D	200[B]	017	C17-C19-N20	-2.00	105.95	109.52
3	B	200	017	C25-O26-C27	2.04	109.79	106.87
3	D	200[B]	017	C25-O26-C27	2.10	109.87	106.87
3	D	200[B]	017	C29-O28-C27	2.13	111.27	107.54
3	D	200[A]	017	C29-O28-C27	2.15	111.30	107.54
3	B	200	017	C29-O28-C27	2.16	111.32	107.54
3	B	200	017	C4-C5-S8	2.26	122.24	119.78
3	B	200	017	C7-C6-C5	2.33	121.98	119.46
3	B	200	017	C30-C31-C27	2.40	105.57	102.27
3	B	200	017	O23-C21-N20	2.44	113.55	110.36
3	D	200[A]	017	O10-S8-N11	2.57	109.12	106.69
3	D	200[B]	017	C30-C31-C27	2.64	105.90	102.27
3	D	200[A]	017	C30-C31-C27	2.69	105.97	102.27
3	B	200	017	C5-S8-N11	3.58	111.66	107.31
3	D	200[A]	017	O23-C21-N20	3.83	115.38	110.36
3	D	200[B]	017	O23-C21-N20	3.88	115.45	110.36
3	D	200[A]	017	C5-S8-N11	4.37	112.63	107.31
3	D	200[B]	017	C5-S8-N11	4.46	112.75	107.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	200	017	O9-S8-C5	4.64	114.05	108.00
3	B	200	017	C24-O23-C21	5.27	124.02	116.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	606	ACT	1	0
3	B	200	017	2	0
2	B	607	ACT	1	0
2	C	603	ACT	1	0
2	C	604	ACT	1	0
3	D	200[A]	017	10	0
3	D	200[B]	017	4	0
2	D	605	ACT	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, 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	99/99 (100%)	-0.18	0	100	100	3, 12, 22, 26	0
1	B	99/99 (100%)	-0.19	0	100	100	3, 11, 22, 25	0
1	C	99/99 (100%)	-0.18	0	100	100	5, 14, 28, 31	0
1	D	99/99 (100%)	-0.14	0	100	100	5, 13, 28, 32	0
All	All	396/396 (100%)	-0.17	0	100	100	3, 12, 23, 32	0

There are no RSRZ outliers to report.

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, 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
2	ACT	C	603	4/4	0.82	0.15	7.45	17,17,18,20	0
2	ACT	C	604	4/4	0.52	0.21	4.30	37,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
4	PO4	C	505	5/5	0.84	0.18	3.44	50,50,50,52	0
2	ACT	D	605	4/4	0.88	0.14	2.22	19,20,20,21	0
2	ACT	D	602	4/4	0.75	0.15	2.14	31,31,32,32	0
2	ACT	A	608	4/4	0.86	0.13	2.14	19,19,20,20	0
4	PO4	C	503	5/5	0.91	0.15	1.71	23,23,24,28	0
2	ACT	D	601	4/4	0.91	0.12	1.65	26,26,26,27	0
2	ACT	B	607	4/4	0.87	0.17	1.23	32,32,32,33	0
2	ACT	A	606	4/4	0.92	0.14	1.17	35,35,35,35	0
4	PO4	D	501	5/5	0.94	0.14	1.09	21,21,23,27	0
3	017	D	200[A]	38/38	0.95	0.13	0.86	2,2,2,3	38
3	017	D	200[B]	38/38	0.95	0.13	0.86	2,2,4,4	38
3	017	B	200	38/38	0.94	0.08	-0.70	2,4,8,8	0
4	PO4	C	502	5/5	0.88	0.18	-	34,38,39,39	0
4	PO4	D	504	5/5	0.82	0.20	-	51,51,52,53	0

6.5 Other polymers (i)

There are no such residues in this entry.