Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:50 AM GMT

PDB ID	:	2J07
Title	:	THERMUS DNA PHOTOLYASE WITH 8-HDF ANTENNA CHROMOPHORE
Authors	:	Klar, T.; Kaiser, G.; Hennecke, U.; Carell, T.; Batschauer, A.; Essen, L.-O.
Deposited on	:	2006-08-01
Resolution	:	1.95 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:

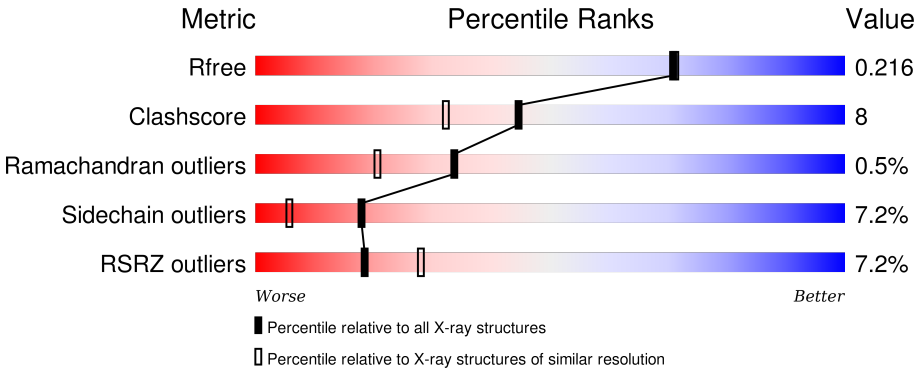
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	420	7% 83% 13% •

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
5	PO4	A	1426	-	-	-	X

2 Entry composition

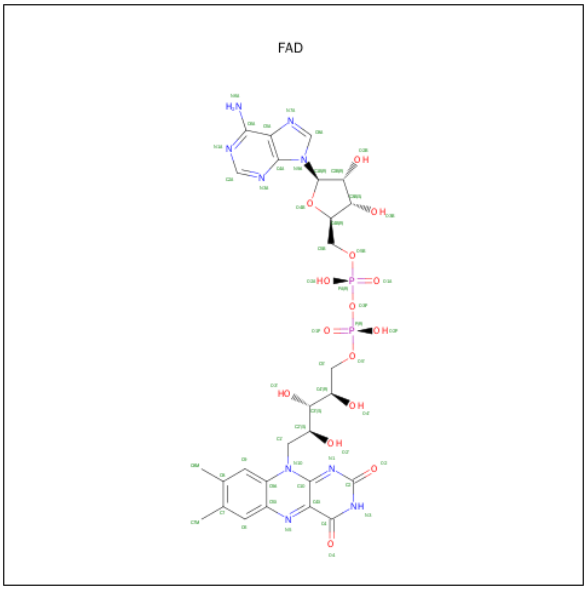
There are 6 unique types of molecules in this entry. The entry contains 3836 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DEOXYRIBODIPYRIMIDINE PHOTO-LYASE.

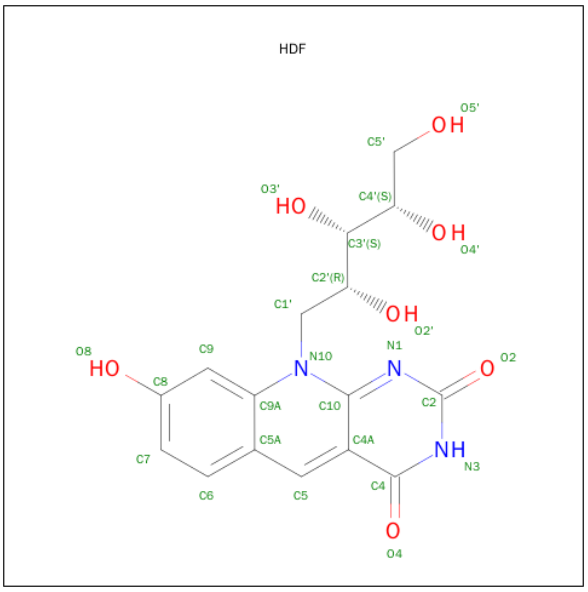
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	21	0	0
			3393	2186	631	572	4			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is 8-HYDROXY-10-(D-RIBO-2,3,4,5-TETRAHYDROXYPENTYL)-5-DEAZAI SOALLOXAZINE (three-letter code: HDF) (formula: $C_{16}H_{17}N_3O_7$).

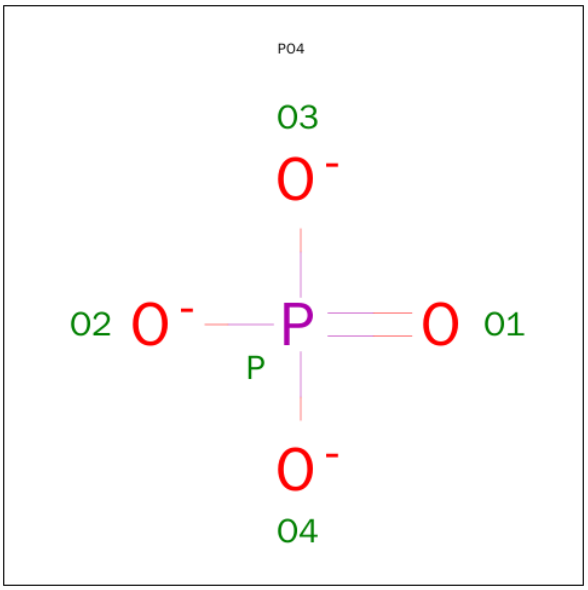


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			26	16	3	7		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Cl	0	0
			3	3		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



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

- Molecule 6 is water.

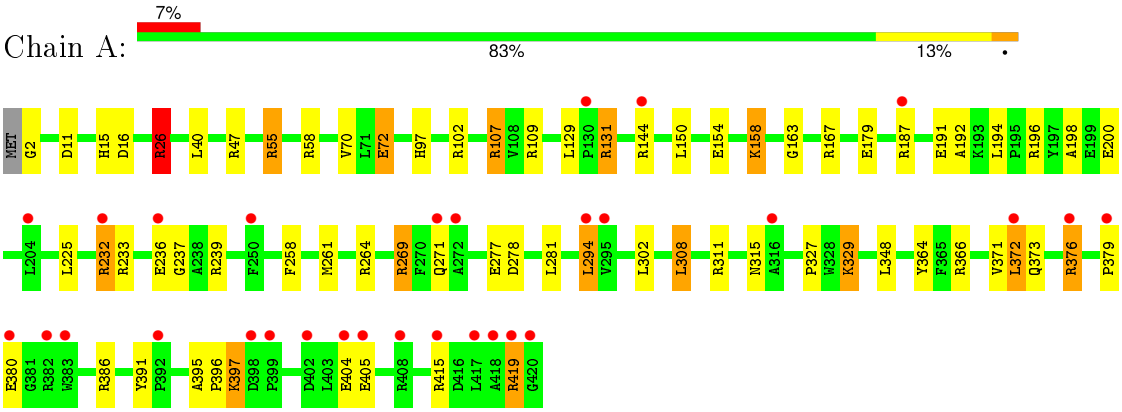
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	356	Total O 356 356	0	0

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: DEOXYRIBODIPYRIMIDINE PHOTO-LYASE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	112.62Å 112.62Å 140.26Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	18.86 – 1.95 18.79 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (18.86-1.95) 99.9 (18.79-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.63 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.169 , 0.212 0.174 , 0.216	Depositor DCC
R_{free} test set	1950 reflections (5.30%)	DCC
Wilson B-factor (Å^2)	42.1	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 60.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Outliers	1 of 38750 reflections (0.003%)	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	3836	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: HDF, PO4, FAD, CL

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	2.92	7/3508 (0.2%)	1.07	13/4783 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	419	ARG	CD-NE	164.41	4.25	1.46
1	A	419	ARG	CB-CG	-24.42	0.86	1.52
1	A	26	ARG	NE-CZ	15.01	1.52	1.33
1	A	158	LYS	CE-NZ	6.00	1.64	1.49
1	A	158	LYS	CG-CD	5.75	1.72	1.52
1	A	158	LYS	CD-CE	5.71	1.65	1.51
1	A	72	GLU	CG-CD	5.68	1.60	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	26	ARG	NE-CZ-NH2	-38.13	101.24	120.30
1	A	26	ARG	CD-NE-CZ	-26.72	86.19	123.60
1	A	419	ARG	CG-CD-NE	-16.37	77.42	111.80
1	A	419	ARG	CA-CB-CG	15.91	148.41	113.40
1	A	419	ARG	CB-CG-CD	13.92	147.80	111.60
1	A	26	ARG	NE-CZ-NH1	12.32	126.46	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	55	ARG	NE-CZ-NH2	-9.75	115.43	120.30
1	A	419	ARG	CD-NE-CZ	-7.45	113.17	123.60
1	A	58	ARG	NE-CZ-NH2	-6.39	117.11	120.30
1	A	55	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	A	269	ARG	CA-CB-CG	5.72	125.99	113.40
1	A	109	ARG	NE-CZ-NH1	5.67	123.13	120.30
1	A	109	ARG	NE-CZ-NH2	-5.13	117.73	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	154	GLU	Mainchain
1	A	26	ARG	Sidechain
1	A	55	ARG	Sidechain

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3393	0	3357	51	0
2	A	53	0	31	0	0
3	A	26	0	16	1	0
4	A	3	0	0	1	0
5	A	5	0	0	1	0
6	A	356	0	0	23	2
All	All	3836	0	3404	54	2

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 (54) 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:236:GLU:O	6:A:2261:HOH:O	1.65	1.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ARG:HH11	1:A:131:ARG:HG3	0.98	1.05
1:A:131:ARG:HH11	1:A:131:ARG:CG	1.71	1.03
1:A:131:ARG:HG3	1:A:131:ARG:NH1	1.65	0.97
1:A:72:GLU:OE2	6:A:2083:HOH:O	1.81	0.96
1:A:237:GLY:N	6:A:2264:HOH:O	1.97	0.96
4:A:1423:CL:CL	6:A:2312:HOH:O	2.21	0.95
1:A:167:ARG:NH1	6:A:2199:HOH:O	2.15	0.79
1:A:236:GLU:HB3	6:A:2263:HOH:O	1.85	0.74
1:A:376:ARG:HH11	1:A:376:ARG:HB2	1.52	0.74
1:A:236:GLU:HA	6:A:2260:HOH:O	1.88	0.73
1:A:107:ARG:HD2	6:A:2136:HOH:O	1.91	0.70
1:A:380:GLU:N	6:A:2332:HOH:O	2.08	0.68
1:A:72:GLU:HG2	6:A:2084:HOH:O	1.94	0.66
1:A:198:ALA:O	6:A:2228:HOH:O	2.14	0.65
1:A:232:ARG:HD2	6:A:2257:HOH:O	1.98	0.63
1:A:15:HIS:CD2	1:A:225:LEU:HD13	2.35	0.62
1:A:277:GLU:OE1	1:A:277:GLU:HA	2.00	0.61
1:A:15:HIS:ND1	1:A:179:GLU:OE2	2.27	0.59
1:A:311:ARG:HH11	1:A:373:GLN:HE22	1.53	0.57
1:A:364:TYR:OH	1:A:415:ARG:HG2	2.05	0.57
1:A:315:ASN:HD21	1:A:373:GLN:HE21	1.54	0.56
1:A:372:LEU:HD13	1:A:376:ARG:HH12	1.72	0.55
1:A:258:PHE:O	1:A:261:MET:HG2	2.07	0.55
1:A:2:GLY:N	6:A:2112:HOH:O	2.41	0.54
1:A:131:ARG:HD3	6:A:2160:HOH:O	2.08	0.53
1:A:232:ARG:HG2	6:A:2257:HOH:O	2.09	0.51
1:A:131:ARG:CD	6:A:2160:HOH:O	2.59	0.50
5:A:1426:PO4:O1	6:A:2356:HOH:O	2.19	0.50
1:A:192:ALA:HB2	6:A:2223:HOH:O	2.13	0.49
1:A:391:TYR:HD2	1:A:397:LYS:HD3	1.78	0.49
1:A:415:ARG:HD2	6:A:2340:HOH:O	2.12	0.48
1:A:194:LEU:O	1:A:237:GLY:HA3	2.14	0.48
1:A:327:PRO:HB2	6:A:2302:HOH:O	2.14	0.47
1:A:232:ARG:HB3	1:A:232:ARG:HE	1.25	0.47
1:A:302:LEU:HB2	1:A:308:LEU:HD22	1.98	0.46
1:A:329:LYS:HB2	6:A:2305:HOH:O	2.16	0.46
1:A:395:ALA:HA	1:A:396:PRO:HD3	1.81	0.46
1:A:70:VAL:O	1:A:163:GLY:HA3	2.16	0.46
1:A:40:LEU:O	1:A:47:ARG:NH1	2.48	0.45
1:A:97:HIS:O	1:A:102:ARG:NH1	2.44	0.45
3:A:1422:HDF:H4'	6:A:2347:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ASP:HA	1:A:150:LEU:HB2	1.99	0.45
1:A:405:GLU:O	1:A:405:GLU:OE2	2.34	0.44
1:A:294:LEU:O	1:A:294:LEU:HD12	2.17	0.44
1:A:131:ARG:NH1	1:A:131:ARG:CG	2.41	0.44
1:A:376:ARG:CB	1:A:376:ARG:HH11	2.26	0.44
1:A:236:GLU:O	1:A:237:GLY:C	2.56	0.42
1:A:311:ARG:HH11	1:A:373:GLN:NE2	2.15	0.42
1:A:281:LEU:HD23	1:A:281:LEU:HA	1.90	0.42
1:A:294:LEU:C	1:A:294:LEU:HD12	2.40	0.42
1:A:366:ARG:NH1	6:A:2323:HOH:O	2.33	0.42
1:A:264:ARG:HD2	1:A:271:GLN:NE2	2.35	0.41
1:A:196:ARG:NH2	1:A:200:GLU:OE2	2.53	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:2039:HOH:O	6:A:2084:HOH:O[11_655]	1.93	0.27
6:A:2041:HOH:O	6:A:2041:HOH:O[11_655]	2.19	0.01

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	417/420 (99%)	409 (98%)	6 (1%)	2 (0%)	34 21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	278	ASP
1	A	379	PRO

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	334/335 (100%)	310 (93%)	24 (7%)	18 6

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	26	ARG
1	A	107	ARG
1	A	129	LEU
1	A	131	ARG
1	A	144	ARG
1	A	158	LYS
1	A	187	ARG
1	A	191	GLU
1	A	232	ARG
1	A	233	ARG
1	A	239	ARG
1	A	269	ARG
1	A	294	LEU
1	A	308	LEU
1	A	329	LYS
1	A	348	LEU
1	A	371	VAL
1	A	372	LEU
1	A	376	ARG
1	A	386	ARG
1	A	397	LYS
1	A	404	GLU
1	A	419	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	39	ASN

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Mol	Chain	Res	Type
1	A	373	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

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 for Mogul analysis.

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	FAD	A	1421	-	48,58,58	1.28	7 (14%)	54,89,89	2.04	9 (16%)
3	HDF	A	1422	-	25,28,28	3.62	12 (48%)	28,41,41	4.52	20 (71%)
5	PO4	A	1426	-	4,4,4	0.40	0	6,6,6	0.31	0

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	FAD	A	1421	-	-	0/30/50/50	0/6/6/6
3	HDF	A	1422	-	-	0/14/14/14	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PO4	A	1426	-	-	0/0/0/0	0/0/0/0

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1422	HDF	C5-C4A	-8.87	1.40	1.53
3	A	1422	HDF	C4A-C10	-6.92	1.45	1.54
3	A	1422	HDF	C6-C5A	-5.73	1.40	1.53
3	A	1422	HDF	C9-C9A	-5.53	1.40	1.53
3	A	1422	HDF	C9-C8	-5.24	1.42	1.51
3	A	1422	HDF	C5-C5A	-4.82	1.41	1.53
3	A	1422	HDF	C7-C8	-4.62	1.38	1.51
3	A	1422	HDF	C7-C6	-4.56	1.41	1.52
3	A	1422	HDF	C5A-C9A	-3.02	1.47	1.53
3	A	1422	HDF	C9A-N10	-2.89	1.42	1.48
3	A	1422	HDF	C4'-C3'	-2.76	1.47	1.53
3	A	1422	HDF	O8-C8	-2.39	1.36	1.43
2	A	1421	FAD	C1'-N10	2.23	1.50	1.48
2	A	1421	FAD	C10-N1	2.27	1.39	1.35
2	A	1421	FAD	C4-N3	2.63	1.38	1.33
2	A	1421	FAD	C5X-N5	2.95	1.40	1.35
2	A	1421	FAD	C2A-N1A	3.32	1.40	1.33
2	A	1421	FAD	C4X-N5	3.46	1.38	1.33
2	A	1421	FAD	C2A-N3A	4.03	1.39	1.32

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1421	FAD	N3A-C2A-N1A	-9.46	121.65	128.89
2	A	1421	FAD	C4X-C4-N3	-4.01	118.10	123.59
3	A	1422	HDF	O4'-C4'-C3'	-3.81	99.44	109.02
2	A	1421	FAD	C2B-C1B-N9A	-2.85	109.94	114.29
2	A	1421	FAD	C1B-N9A-C4A	-2.83	122.67	126.94
2	A	1421	FAD	C4A-C5A-N7A	-2.82	106.89	109.48
3	A	1422	HDF	C1'-C2'-C3'	2.32	116.45	109.82
2	A	1421	FAD	C1'-C2'-C3'	2.32	116.46	109.82
3	A	1422	HDF	C8-C9-C9A	2.84	118.75	112.96
3	A	1422	HDF	O8-C8-C7	2.85	118.11	110.05
2	A	1421	FAD	C4X-N5-C5X	2.85	120.04	116.76
2	A	1421	FAD	C5X-C9A-N10	3.00	119.90	117.62
3	A	1422	HDF	C4A-C5-C5A	3.10	120.55	110.08
3	A	1422	HDF	C7-C6-C5A	3.46	120.42	112.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1422	HDF	C5A-C9A-N10	4.16	119.16	110.12
3	A	1422	HDF	C9-C9A-N10	4.20	121.61	113.03
3	A	1422	HDF	C6-C7-C8	4.21	119.38	111.59
3	A	1422	HDF	C5-C5A-C6	4.21	121.03	112.68
3	A	1422	HDF	O8-C8-C9	4.68	119.17	109.86
3	A	1422	HDF	N3-C2-N1	5.21	121.85	116.14
3	A	1422	HDF	O3'-C3'-C2'	5.43	122.43	108.75
3	A	1422	HDF	O4'-C4'-C5'	5.61	122.30	109.22
3	A	1422	HDF	O2'-C2'-C3'	5.73	123.41	109.02
3	A	1422	HDF	C6-C5A-C9A	6.04	119.44	110.23
2	A	1421	FAD	C4-N3-C2	6.25	120.65	115.25
3	A	1422	HDF	C5-C5A-C9A	6.48	119.30	109.94
3	A	1422	HDF	C5-C4A-C4	6.70	124.05	110.89
3	A	1422	HDF	C5-C4A-C10	8.94	120.62	107.89
3	A	1422	HDF	C9-C8-C7	9.53	122.67	110.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1422	HDF	1	0
5	A	1426	PO4	1	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	419/420 (99%)	0.25	30 (7%) 18 28	30, 41, 57, 79	5 (1%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	419	ARG	8.2
1	A	420	GLY	5.8
1	A	236	GLU	5.1
1	A	418	ALA	5.1
1	A	379	PRO	4.9
1	A	417	LEU	4.7
1	A	372	LEU	4.6
1	A	398	ASP	4.6
1	A	376	ARG	3.9
1	A	272	ALA	3.8
1	A	405	GLU	3.4
1	A	415	ARG	3.3
1	A	295	VAL	3.1
1	A	404	GLU	3.1
1	A	144	ARG	2.8
1	A	294	LEU	2.7
1	A	382	ARG	2.7
1	A	380	GLU	2.7
1	A	130	PRO	2.7
1	A	383	TRP	2.5
1	A	399	PRO	2.5
1	A	402	ASP	2.5
1	A	187	ARG	2.4
1	A	250	PHE	2.4
1	A	408	ARG	2.4
1	A	232	ARG	2.3
1	A	392	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	204	LEU	2.2
1	A	316	ALA	2.1
1	A	271	GLN	2.1

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
5	PO4	A	1426	5/5	0.97	0.18	3.33	47,48,56,58	0
4	CL	A	1425	1/1	0.93	0.11	0.61	74,74,74,74	0
2	FAD	A	1421	53/53	0.96	0.09	-0.51	39,42,45,46	0
3	HDF	A	1422	26/26	0.93	0.08	-0.63	35,39,53,56	0
4	CL	A	1423	1/1	0.96	0.10	-0.81	86,86,86,86	0
4	CL	A	1424	1/1	0.92	0.12	-	85,85,85,85	0

6.5 Other polymers (i)

There are no such residues in this entry.