



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2024 – 05:36 PM EDT

PDB ID : 6G30
Title : Crystal structure of the p97 D2 domain in a helical split-washer conformation
Authors : Stach, L.; Morgan, R.M.L.; Freemont, P.S.
Deposited on : 2018-03-23
Resolution : 2.42 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

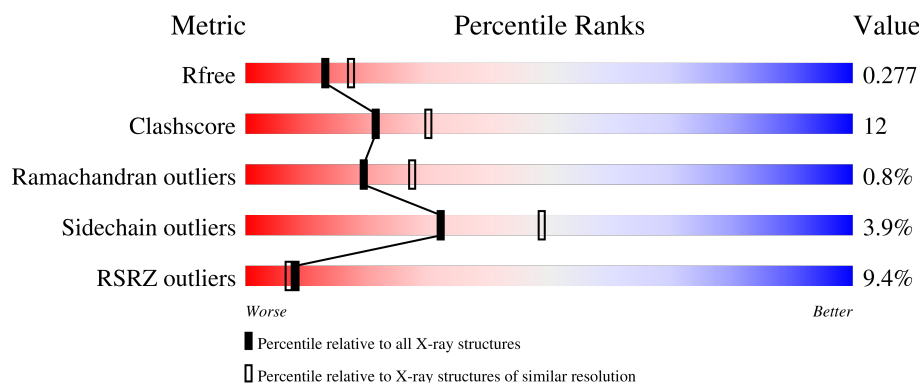
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	130704	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	306	

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	MPD	A	801	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NA	A	808	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 2193 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.

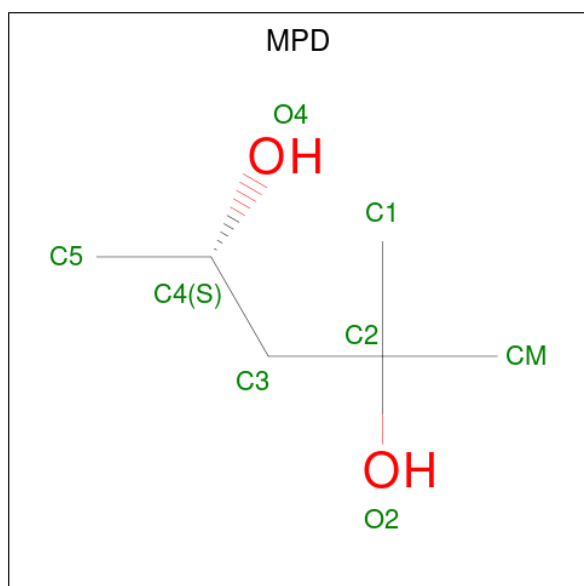
- Molecule 1 is a protein called Transitional endoplasmic reticulum ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	266	2112	1336	369	396	11	0	1	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	459	GLY	-	expression tag	UNP P55072
A	460	PRO	-	expression tag	UNP P55072
A	461	GLY	-	expression tag	UNP P55072

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



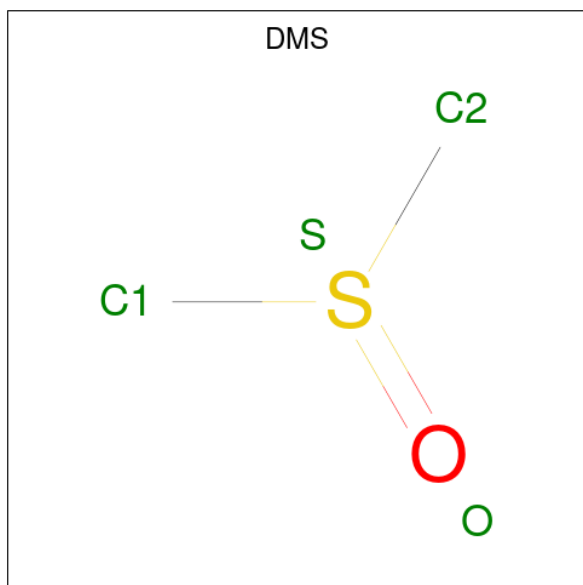
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	8	6	2	0	0

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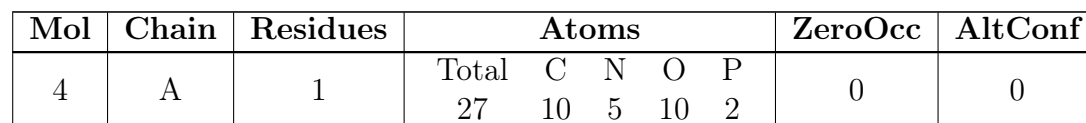
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 3 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 5 | A | 1 | Total Na
1 1 | 0 | 0 |

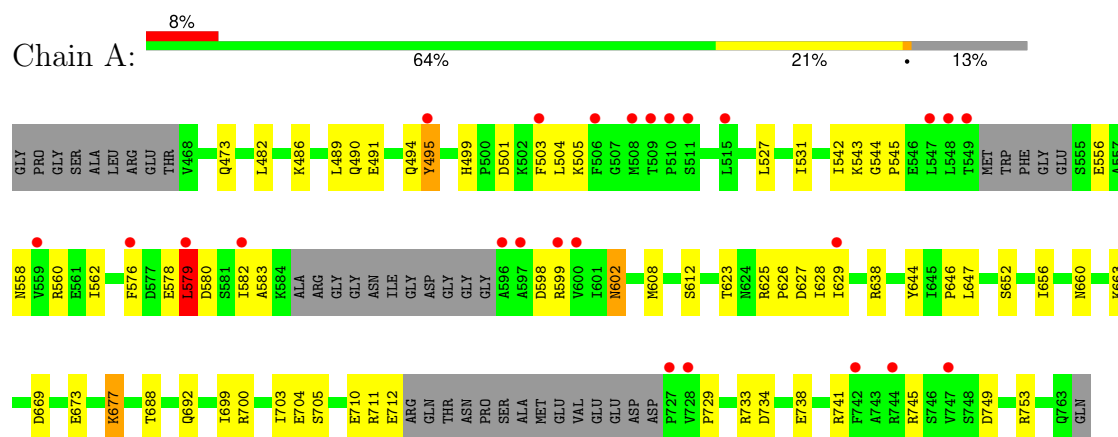
- ELN
-
- The chemical structure is 4-chlorobenzamide, 2-cyanoethyl ester. It features a central amide group (-C(=O)-NH-) connecting a 2-cyanoethyl group and a 4-chlorophenyl group. The 2-cyanoethyl group consists of a methylene group (C10) adjacent to a nitrile group (C11≡N12). The amide carbonyl is represented by C08=O09. The nitrogen of the amide is N07. The 4-chlorophenyl group is a benzene ring with carbons labeled C01 through C06 and a chlorine atom (Cl1) at the para position (C03).
- N#CCCC(=O)Nc1ccc(Cl)cc1

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	Cl	N	O	0	0
			13	9	1	2	1		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Transitional endoplasmic reticulum ATPase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	94.76Å 94.76Å 66.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.38 – 2.42 51.60 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.5 (47.38-2.42) 99.8 (51.60-2.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.42Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.227 , 0.273 0.236 , 0.277	Depositor DCC
R_{free} test set	669 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	68.8	Xtriage
Anisotropy	0.664	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 88.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.068 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2193	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, DMS, ELN, NA, MPD

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	0.48	0/2148	0.70	2/2893 (0.1%)

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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	579	LEU	CA-CB-CG	13.09	145.41	115.30
1	A	647	LEU	CA-CB-CG	5.32	127.54	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	579	LEU	Peptide

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	2112	0	2146	53	0
2	A	32	0	56	7	0
3	A	8	0	12	0	0
4	A	27	0	12	0	0
5	A	1	0	0	0	0
6	A	13	0	0	0	0
All	All	2193	0	2226	53	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 12.

All (53) 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:629:ILE:HA	2:A:802:MPD:H11	1.63	0.79
1:A:625:ARG:HH11	1:A:628:ILE:HD11	1.48	0.78
1:A:491:GLU:OE1	1:A:495:TYR:CE1	2.38	0.77
1:A:543:LYS:HG3	1:A:578:GLU:HG2	1.66	0.77
1:A:501:ASP:HA	1:A:504:LEU:CD2	2.18	0.74
1:A:738:GLU:OE1	1:A:741:ARG:NH1	2.30	0.65
1:A:700:ARG:O	1:A:704:GLU:HG3	1.99	0.63
1:A:501:ASP:HA	1:A:504:LEU:HD21	1.82	0.62
1:A:545:PRO:HD3	1:A:578:GLU:OE2	2.00	0.62
1:A:625:ARG:HH11	1:A:628:ILE:CD1	2.12	0.61
1:A:579:LEU:O	2:A:802:MPD:O2	2.17	0.61
1:A:556:GLU:OE2	1:A:560:ARG:NH1	2.35	0.60
1:A:626:PRO:O	1:A:629:ILE:HD12	2.03	0.57
1:A:652:SER:O	1:A:656:ILE:HG13	2.04	0.57
1:A:625:ARG:HD2	1:A:628:ILE:CD1	2.35	0.56
1:A:627:ASP:OD1	1:A:627:ASP:N	2.37	0.55
1:A:608:MET:SD	1:A:638:ARG:HB3	2.48	0.53
1:A:602:ASN:ND2	1:A:602:ASN:O	2.42	0.52
1:A:710:GLU:OE2	1:A:711:ARG:HG2	2.09	0.52
1:A:542:ILE:HG21	1:A:562:ILE:HG21	1.90	0.52
1:A:495:TYR:HB3	1:A:503:PHE:CZ	2.46	0.51
1:A:625:ARG:NH1	1:A:628:ILE:HD11	2.23	0.50
1:A:699:ILE:HG12	2:A:801:MPD:O2	2.12	0.50
1:A:580:ASP:HB3	1:A:623:THR:OG1	2.10	0.50
1:A:578:GLU:O	1:A:578:GLU:HG3	2.11	0.50
1:A:729:PRO:O	2:A:801:MPD:HM3	2.11	0.50
1:A:473:GLN:H	1:A:473:GLN:CD	2.15	0.49
1:A:482:LEU:O	1:A:486:LYS:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:GLN:O	1:A:494:GLN:HG3	2.13	0.49
1:A:745:ARG:NH2	1:A:749:ASP:OD1	2.38	0.48
1:A:583:ALA:HB2	2:A:802:MPD:H53	1.95	0.47
1:A:542:ILE:HD11	1:A:576:PHE:CE2	2.51	0.46
1:A:644:TYR:CZ	1:A:646:PRO:HB3	2.51	0.46
1:A:527:LEU:O	1:A:531:ILE:HG13	2.16	0.46
1:A:558:ASN:O	1:A:562:ILE:HG12	2.16	0.46
1:A:628:ILE:HG22	1:A:628:ILE:O	2.16	0.45
1:A:669:ASP:O	1:A:733:ARG:HD3	2.17	0.45
1:A:703:ILE:HG13	1:A:704:GLU:N	2.30	0.45
1:A:489:LEU:HD13	1:A:531:ILE:HD11	1.98	0.44
1:A:543:LYS:NZ	1:A:578:GLU:OE1	2.21	0.44
1:A:504:LEU:HG	1:A:505:LYS:N	2.33	0.43
1:A:710:GLU:O	1:A:712:GLU:OE2	2.35	0.43
1:A:580:ASP:HB2	1:A:625:ARG:HB2	1.99	0.43
1:A:733:ARG:HG3	2:A:803:MPD:O2	2.18	0.43
1:A:626:PRO:HA	1:A:629:ILE:HD11	2.00	0.43
1:A:629:ILE:HG23	2:A:802:MPD:H31	2.00	0.43
1:A:660:ASN:HD22	1:A:660:ASN:HA	1.63	0.43
1:A:579:LEU:O	1:A:582:ILE:HG22	2.19	0.43
1:A:688:THR:O	1:A:692:GLN:HG3	2.19	0.42
1:A:626:PRO:HA	1:A:629:ILE:CD1	2.50	0.42
1:A:580:ASP:CB	1:A:625:ARG:HB2	2.50	0.41
1:A:749:ASP:O	1:A:753:ARG:HG3	2.21	0.41
1:A:673:GLU:O	1:A:677:LYS:HE2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	259/306 (85%)	247 (95%)	10 (4%)	2 (1%)	19	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	598	ASP
1	A	544	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	231/258 (90%)	222 (96%)	9 (4%)	32	49

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	495	TYR
1	A	499	HIS
1	A	599	ARG
1	A	602	ASN
1	A	612	SER
1	A	663	LYS
1	A	677	LYS
1	A	705	SER
1	A	734	ASP

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

Mol	Chain	Res	Type
1	A	499	HIS
1	A	602	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	MPD	A	802	-	7,7,7	0.26	0	9,10,10	0.70	0
4	ADP	A	807	-	24,29,29	1.35	3 (12%)	29,45,45	1.29	3 (10%)
3	DMS	A	805	-	3,3,3	0.72	0	3,3,3	0.60	0
2	MPD	A	803	-	7,7,7	0.32	0	9,10,10	0.57	0
6	ELN	A	809	-	13,13,13	2.30	6 (46%)	16,16,16	1.57	3 (18%)
2	MPD	A	801	-	7,7,7	0.40	0	9,10,10	0.65	0
3	DMS	A	806	-	3,3,3	0.71	0	3,3,3	0.42	0
2	MPD	A	804	-	7,7,7	0.34	0	9,10,10	0.35	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	MPD	A	802	-	-	1/5/5/5	-
4	ADP	A	807	-	-	5/12/32/32	0/3/3/3
2	MPD	A	803	-	-	5/5/5/5	-
6	ELN	A	809	-	-	3/7/7/7	0/1/1/1
2	MPD	A	801	-	-	4/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MPD	A	804	-	-	4/5/5/5	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	809	ELN	C08-N07	5.70	1.47	1.35
6	A	809	ELN	C10-C11	3.69	1.52	1.45
4	A	807	ADP	C2-N3	3.18	1.37	1.32
4	A	807	ADP	O4'-C1'	3.10	1.45	1.40
4	A	807	ADP	PA-O3A	2.64	1.62	1.59
6	A	809	ELN	C06-N07	2.48	1.46	1.41
6	A	809	ELN	O09-C08	-2.36	1.18	1.23
6	A	809	ELN	C10-C08	2.14	1.54	1.52
6	A	809	ELN	C03-CL1	2.08	1.79	1.74

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	807	ADP	N3-C2-N1	-3.89	123.40	128.67
6	A	809	ELN	O09-C08-N07	-3.59	117.21	123.64
4	A	807	ADP	N6-C6-N1	3.07	124.89	118.33
6	A	809	ELN	C10-C08-N07	3.00	126.08	115.37
6	A	809	ELN	C11-C10-C08	2.54	116.68	111.36
4	A	807	ADP	O3B-PB-O2B	2.20	116.05	107.80

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	MPD	C2-C3-C4-O4
2	A	801	MPD	C2-C3-C4-C5
2	A	803	MPD	C2-C3-C4-O4
2	A	803	MPD	C2-C3-C4-C5
4	A	807	ADP	PA-O3A-PB-O2B
4	A	807	ADP	C5'-O5'-PA-O2A
6	A	809	ELN	O09-C08-N07-C06
6	A	809	ELN	C10-C08-N07-C06
2	A	801	MPD	C1-C2-C3-C4
2	A	803	MPD	C1-C2-C3-C4
2	A	803	MPD	CM-C2-C3-C4
2	A	804	MPD	C1-C2-C3-C4

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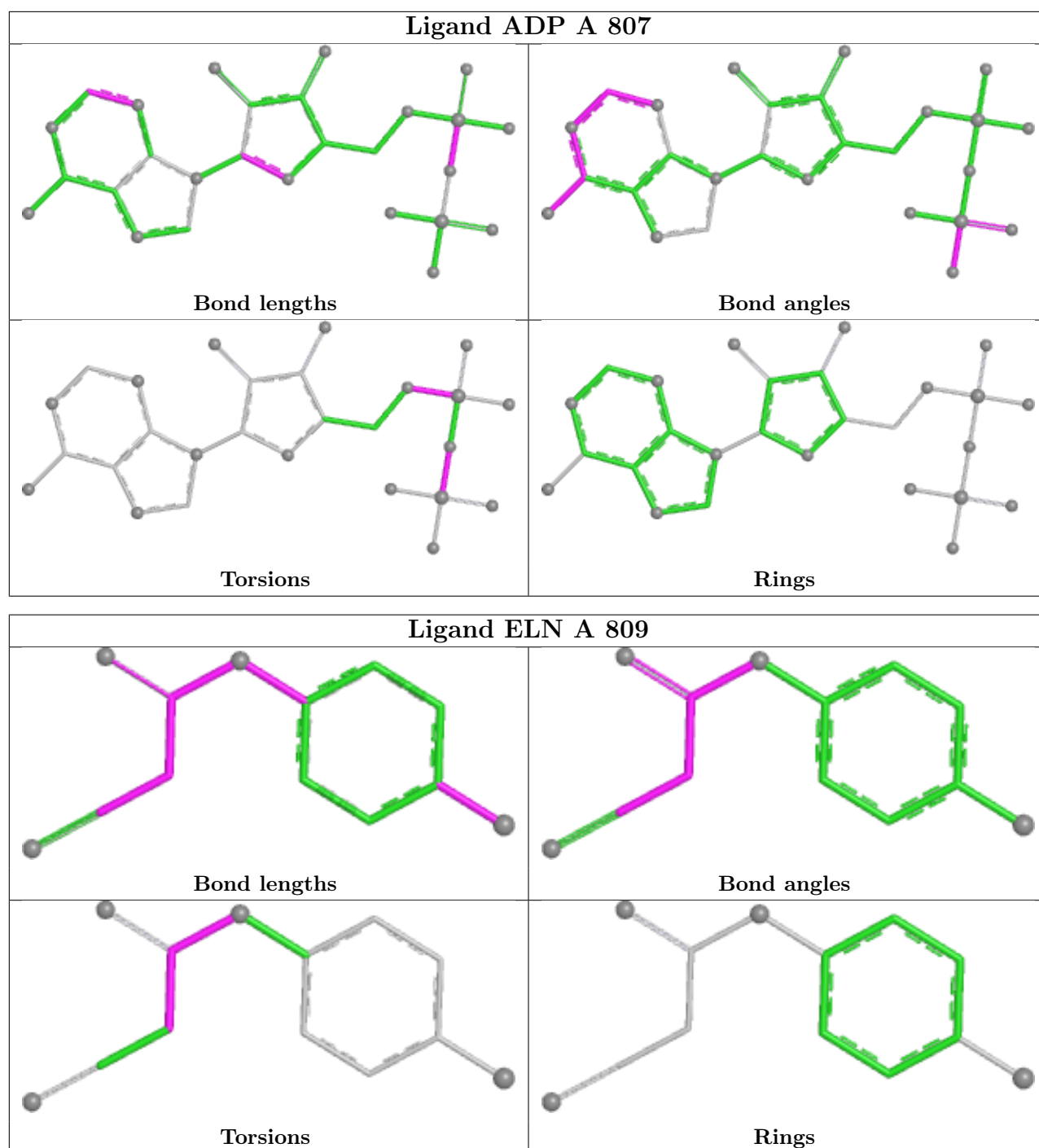
Mol	Chain	Res	Type	Atoms
2	A	804	MPD	CM-C2-C3-C4
2	A	802	MPD	C2-C3-C4-C5
4	A	807	ADP	C5'-O5'-PA-O1A
4	A	807	ADP	C5'-O5'-PA-O3A
4	A	807	ADP	PA-O3A-PB-O3B
6	A	809	ELN	N07-C08-C10-C11
2	A	803	MPD	O2-C2-C3-C4
2	A	804	MPD	O2-C2-C3-C4
2	A	804	MPD	C2-C3-C4-O4
2	A	801	MPD	CM-C2-C3-C4

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	802	MPD	4	0
2	A	803	MPD	1	0
2	A	801	MPD	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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, 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	266/306 (86%)	0.44	25 (9%) 8 7	67, 97, 137, 156	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	547	LEU	5.5
1	A	742	PHE	5.4
1	A	548	LEU	5.0
1	A	582	ILE	5.0
1	A	728	VAL	4.3
1	A	495	TYR	4.1
1	A	596	ALA	4.0
1	A	511	SER	4.0
1	A	597	ALA	3.9
1	A	579	LEU	3.9
1	A	506	PHE	3.7
1	A	503	PHE	3.7
1	A	549	THR	3.6
1	A	727	PRO	3.2
1	A	599	ARG	3.2
1	A	508	MET	2.7
1	A	559	VAL	2.4
1	A	576	PHE	2.4
1	A	509	THR	2.2
1	A	629	ILE	2.1
1	A	744	ARG	2.1
1	A	600	VAL	2.1
1	A	510	PRO	2.0
1	A	747	VAL	2.0
1	A	515	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

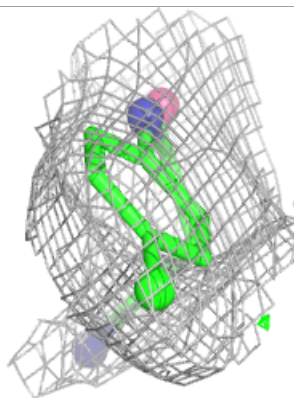
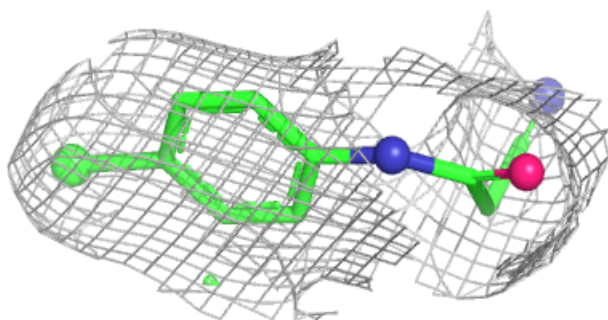
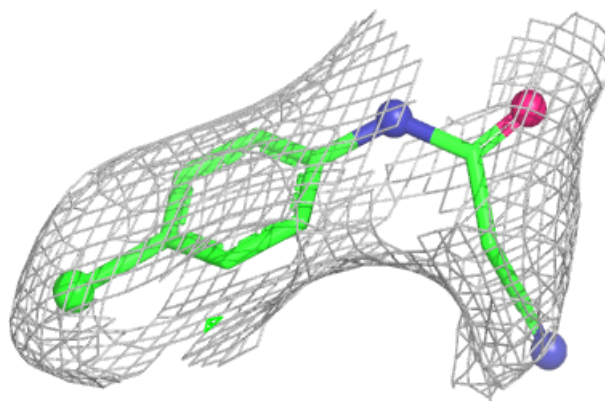
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NA	A	808	1/1	0.49	3.34	107,107,107,107	0
2	MPD	A	804	8/8	0.70	0.33	95,104,112,117	0
2	MPD	A	801	8/8	0.77	0.46	102,107,110,111	0
3	DMS	A	806	4/4	0.86	0.27	100,106,111,124	0
6	ELN	A	809	13/13	0.87	0.11	95,109,121,125	0
2	MPD	A	802	8/8	0.88	0.57	100,112,114,116	0
3	DMS	A	805	4/4	0.92	0.16	99,113,114,117	0
2	MPD	A	803	8/8	0.94	0.34	83,101,104,109	0
4	ADP	A	807	27/27	0.97	0.19	61,72,88,94	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

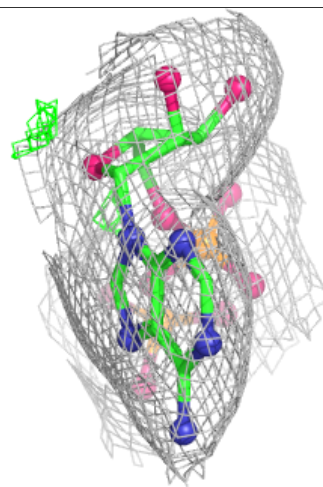
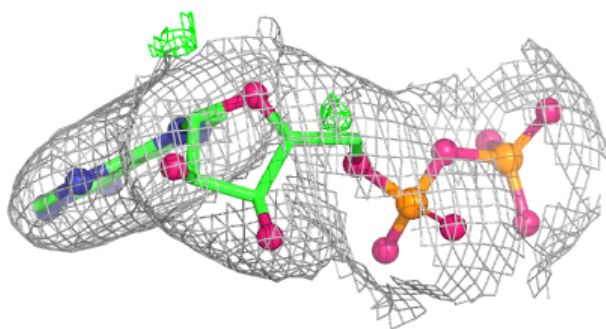
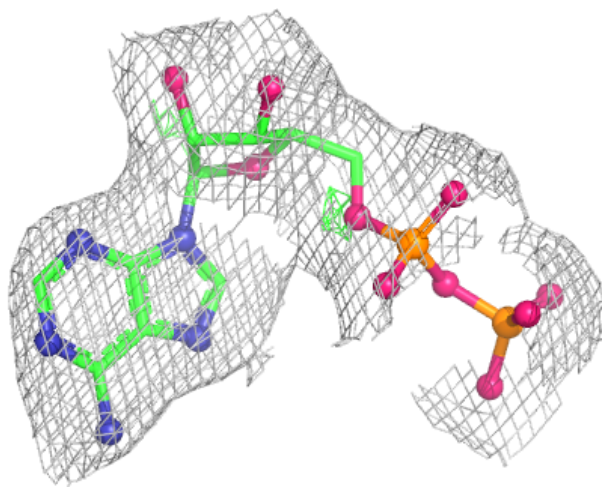
Electron density around ELN A 809:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ADP A 807:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.