



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:02 AM UTC

PDB ID : 2QYN / pdb_00002qyn
Title : Crystal structure of PDE4D2 in complex with inhibitor NPV
Authors : Ke, H.
Deposited on : 2007-08-15
Resolution : 1.57 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

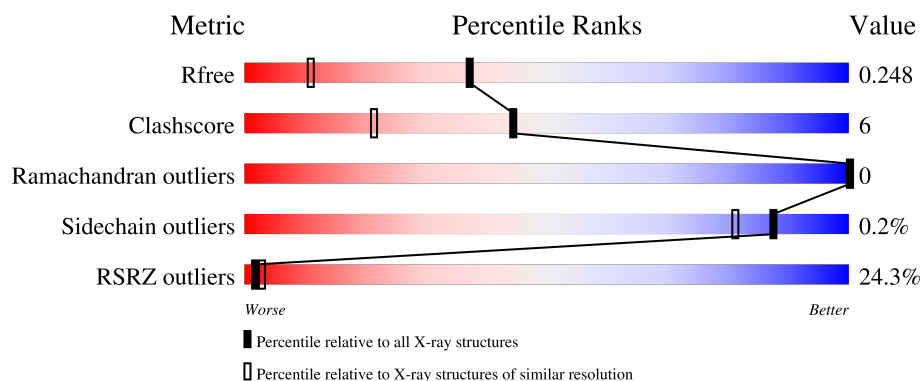
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.57 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}	180053	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	328	11% 87% 12% ..
1	B	328	37% 81% 18% .

2 Entry composition [i](#)

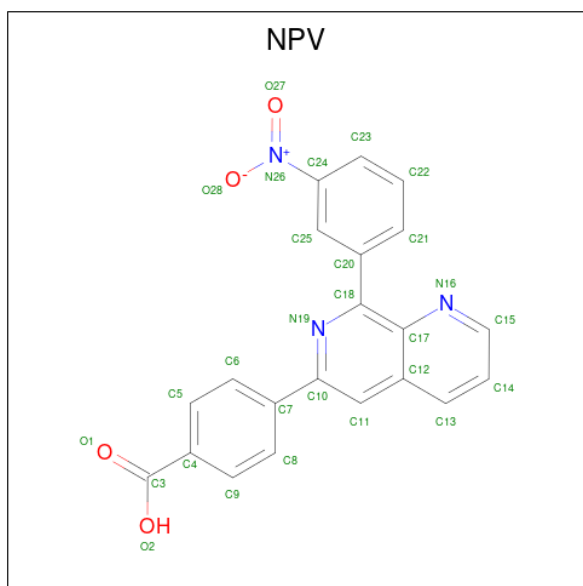
There are 5 unique types of molecules in this entry. The entry contains 5694 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 cAMP-specific 3',5'-cyclic phosphodiesterase 4D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2638	1668	450	506	14			
1	B	325	Total	C	N	O	S	0	0	0
			2631	1664	450	503	14			

- Molecule 2 is 4-[8-(3-nitrophenyl)-1,7-naphthyridin-6-yl]benzoic acid (CCD ID: NPV) (formula: C₂₁H₁₃N₃O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	21	3	4		
2	B	1	Total	C	N	O	0	0
			28	21	3	4		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Zn 1	0	0
3	B	1	Total 1	Zn 1	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Mg 1	0	0
4	B	1	Total 1	Mg 1	0	0

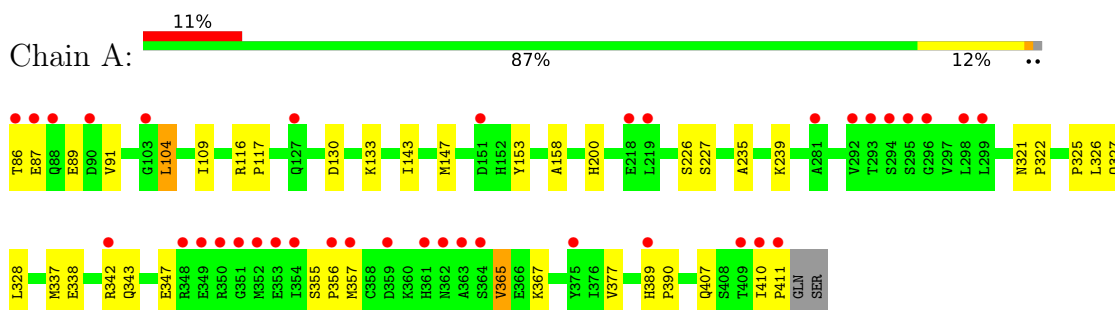
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	225	Total 225	O 225	0	0
5	B	140	Total 140	O 140	0	0

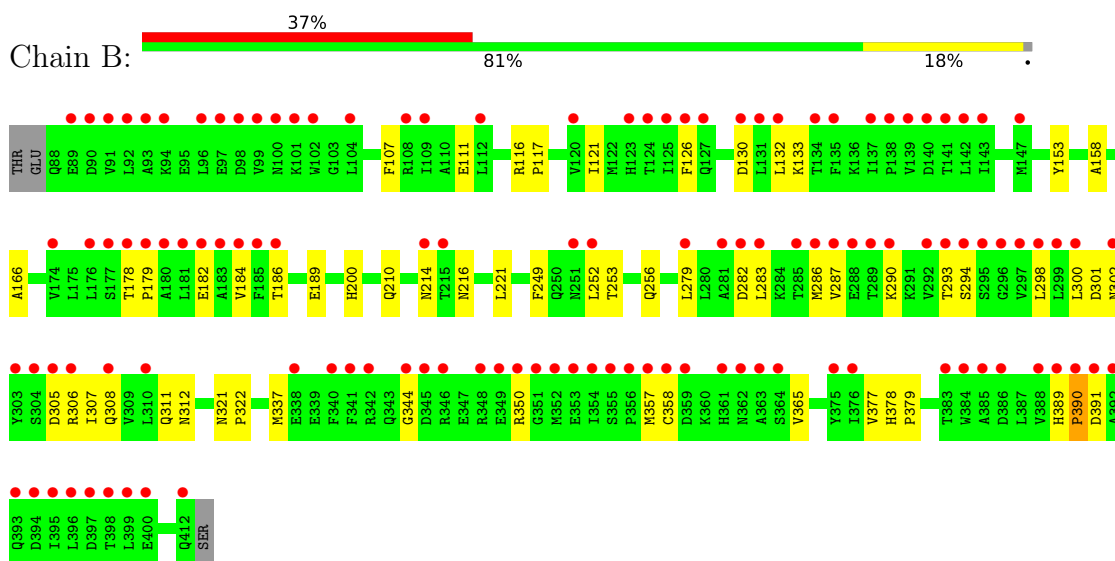
3 Residue-property plots

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: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.10Å 80.63Å 163.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.57 30.00 – 1.57	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.57) 92.6 (30.00-1.57)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 1.57Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.225 , 0.248 0.224 , 0.248	Depositor DCC
R_{free} test set	10439 reflections (10.05%)	wwPDB-VP
Wilson B-factor (Å ²)	19.3	Xtriage
Anisotropy	0.325	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5694	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NPV, ZN, MG

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	0/2692	0.82	7/3658 (0.2%)
1	B	0.33	0/2685	0.82	5/3648 (0.1%)
All	All	0.34	0/5377	0.82	12/7306 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	377	VAL	N-CA-C	7.43	117.52	110.53
1	B	200	HIS	N-CA-C	6.72	121.08	112.34
1	B	377	VAL	N-CA-C	6.58	116.71	110.53
1	A	200	HIS	N-CA-C	6.53	120.83	112.34
1	A	158	ALA	N-CA-C	6.32	118.16	111.28
1	B	158	ALA	N-CA-C	6.26	118.11	111.28
1	A	153	TYR	N-CA-C	-6.00	100.78	110.32
1	A	326	LEU	N-CA-C	5.39	117.85	111.33
1	A	365	VAL	N-CA-C	5.38	116.01	110.36
1	B	182	GLU	N-CA-C	5.20	117.62	111.33
1	B	153	TYR	N-CA-C	-5.03	101.52	109.76
1	A	104	LEU	N-CA-C	-5.02	103.71	110.24

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	2638	0	2591	24	0
1	B	2631	0	2586	41	0
2	A	28	0	12	1	0
2	B	28	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	225	0	0	1	0
5	B	140	0	0	3	0
All	All	5694	0	5201	63	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 6.

All (63) 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:B:293:THR:HG22	1:B:294:SER:H	1.30	0.93
1:A:226:SER:HB2	5:A:661:HOH:O	1.75	0.87
1:A:239:LYS:HG2	1:B:221:LEU:HD21	1.55	0.86
1:B:184:VAL:HG11	1:B:300:LEU:HD12	1.69	0.73
1:B:210:GLN:HG3	1:B:214:ASN:HD21	1.55	0.71
1:A:239:LYS:HG2	1:B:221:LEU:CD2	2.21	0.69
1:B:282:ASP:HB3	1:B:308:GLN:NE2	2.07	0.69
1:A:104:LEU:HD11	1:A:109:ILE:HD11	1.75	0.68
1:B:179:PRO:HD2	1:B:391:ASP:CG	2.19	0.68
1:B:321:ASN:HB2	1:B:322:PRO:HD3	1.79	0.64
1:B:293:THR:HG22	1:B:294:SER:N	2.10	0.63
1:B:302:ASN:O	1:B:306:ARG:HG3	1.98	0.62
1:A:389:HIS:CE1	1:A:390:PRO:HB3	2.37	0.60
1:B:282:ASP:HB3	1:B:308:GLN:CD	2.28	0.59
1:B:178:THR:HG23	1:B:391:ASP:OD2	2.02	0.59
1:B:210:GLN:HG3	1:B:214:ASN:ND2	2.18	0.58
1:B:283:LEU:O	1:B:287:VAL:HG23	2.04	0.58
1:B:301:ASP:OD1	5:B:606:HOH:O	2.17	0.58
1:B:107:PHE:O	1:B:111:GLU:HG3	2.02	0.57
1:A:356:PRO:HB2	1:A:357:MET:HE2	1.87	0.56
1:B:253:THR:OG1	1:B:256:GLN:HG3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:GLN:O	1:A:347:GLU:HG3	2.06	0.54
1:B:184:VAL:CG1	1:B:300:LEU:HD12	2.37	0.54
1:A:235:ALA:O	1:A:239:LYS:HG3	2.08	0.53
1:B:189:GLU:OE1	1:B:306:ARG:NH2	2.41	0.53
1:B:186:THR:HG23	1:B:306:ARG:HH22	1.73	0.52
1:B:121:ILE:HD12	1:B:166:ALA:HB1	1.93	0.51
1:B:186:THR:OG1	1:B:189:GLU:HG3	2.10	0.50
1:A:130:ASP:OD1	1:A:133:LYS:HD2	2.11	0.50
1:B:357:MET:SD	2:B:1:NPV:C23	3.01	0.49
1:A:321:ASN:HB2	1:A:322:PRO:HD3	1.93	0.49
1:B:130:ASP:OD1	1:B:133:LYS:HD2	2.13	0.49
1:B:389:HIS:CE1	1:B:390:PRO:HB3	2.47	0.49
1:A:87:GLU:O	1:A:91:VAL:HG23	2.13	0.49
1:B:344:GLY:HA3	1:B:358:CYS:O	2.12	0.49
1:A:86:THR:HA	1:A:89:GLU:OE2	2.14	0.48
1:B:307:ILE:O	1:B:311:GLN:HG3	2.13	0.47
1:A:410:ILE:HB	1:A:411:PRO:HD3	1.95	0.47
1:B:293:THR:CG2	1:B:294:SER:H	2.11	0.47
1:B:179:PRO:HD2	1:B:391:ASP:OD2	2.15	0.47
1:B:290:LYS:HG3	1:B:298:LEU:CD1	2.44	0.46
1:B:378:HIS:HB3	1:B:379:PRO:HD3	1.97	0.46
1:A:355:SER:HB3	1:A:356:PRO:HD2	1.98	0.46
1:A:357:MET:SD	2:A:1:NPV:C23	3.04	0.46
1:B:126:PHE:HB3	1:B:132:LEU:CD1	2.46	0.45
1:A:116:ARG:N	1:A:117:PRO:CD	2.80	0.45
1:B:249:PHE:HA	1:B:252:LEU:HD13	1.98	0.45
1:B:286:MET:HE2	1:B:305:ASP:HB3	1.99	0.45
1:A:407:GLN:O	1:A:411:PRO:HD3	2.17	0.45
1:B:389:HIS:HA	1:B:390:PRO:HA	1.71	0.45
1:A:337:MET:HE2	1:A:365:VAL:HG22	1.99	0.45
1:A:143:ILE:O	1:A:147:MET:HG3	2.16	0.44
1:B:116:ARG:N	1:B:117:PRO:CD	2.80	0.44
1:B:350:ARG:HD3	5:B:577:HOH:O	2.18	0.44
1:A:327:GLN:NE2	1:A:328:LEU:HG	2.33	0.43
1:B:216:ASN:HA	5:B:637:HOH:O	2.19	0.42
1:B:337:MET:HE2	1:B:365:VAL:HG22	2.01	0.42
1:B:279:LEU:HG	1:B:312:ASN:OD1	2.20	0.41
1:A:367:LYS:HG2	1:A:410:ILE:HD13	2.03	0.41
1:A:226:SER:O	1:A:227:SER:C	2.64	0.41
1:A:338:GLU:OE2	1:A:342:ARG:NH2	2.55	0.40
1:A:325:PRO:HB2	1:A:327:GLN:HE22	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:LEU:H	1:B:132:LEU:HD12	1.87	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	324/328 (99%)	316 (98%)	8 (2%)	0	100	100
1	B	323/328 (98%)	312 (97%)	11 (3%)	0	100	100
All	All	647/656 (99%)	628 (97%)	19 (3%)	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	298/300 (99%)	298 (100%)	0	100	100
1	B	297/300 (99%)	296 (100%)	1 (0%)	86	79
All	All	595/600 (99%)	594 (100%)	1 (0%)	87	81

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	390	PRO

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

Mol	Chain	Res	Type
1	A	127	GLN
1	A	210	GLN
1	A	242	GLN
1	A	321	ASN
1	A	327	GLN
1	A	361	HIS
1	A	362	ASN
1	B	210	GLN
1	B	214	ASN
1	B	242	GLN
1	B	245	ASN
1	B	258	GLN
1	B	308	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	NPV	A	1	-	31,31,31	2.28	7 (22%)	39,44,44	1.36	7 (17%)
2	NPV	B	1	-	31,31,31	2.29	6 (19%)	39,44,44	1.38	5 (12%)

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	NPV	A	1	-	-	0/14/16/16	0/4/4/4
2	NPV	B	1	-	-	0/14/16/16	0/4/4/4

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NPV	C20-C18	-6.60	1.39	1.49
2	A	1	NPV	C20-C18	-6.54	1.39	1.49
2	B	1	NPV	C7-C10	-6.08	1.39	1.49
2	A	1	NPV	C7-C10	-6.02	1.39	1.49
2	B	1	NPV	C4-C3	-5.11	1.38	1.49
2	A	1	NPV	C4-C3	-5.10	1.38	1.49
2	B	1	NPV	C24-N26	-4.90	1.33	1.45
2	A	1	NPV	C24-N26	-4.89	1.33	1.45
2	A	1	NPV	C11-C10	2.35	1.39	1.37
2	B	1	NPV	C11-C10	2.27	1.39	1.37
2	B	1	NPV	O28-N26	-2.06	1.21	1.35
2	A	1	NPV	O28-N26	-2.06	1.21	1.35
2	A	1	NPV	C12-C17	-2.04	1.39	1.42

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	NPV	C15-N16-C17	3.68	121.65	117.31
2	B	1	NPV	C15-N16-C17	3.63	121.59	117.31
2	B	1	NPV	C20-C18-N19	3.05	118.96	115.29
2	A	1	NPV	C20-C18-N19	2.85	118.72	115.29
2	A	1	NPV	C18-N19-C10	2.72	122.43	118.07
2	B	1	NPV	C18-N19-C10	2.60	122.25	118.07
2	A	1	NPV	C14-C15-N16	-2.21	120.72	123.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NPV	C14-C15-N16	-2.17	120.78	123.97
2	B	1	NPV	C12-C17-N16	-2.11	120.41	122.63
2	A	1	NPV	C11-C10-N19	-2.09	119.82	121.47
2	A	1	NPV	C12-C17-N16	-2.07	120.45	122.63
2	A	1	NPV	C10-C11-C12	-2.03	119.05	120.55

There are no chirality outliers.

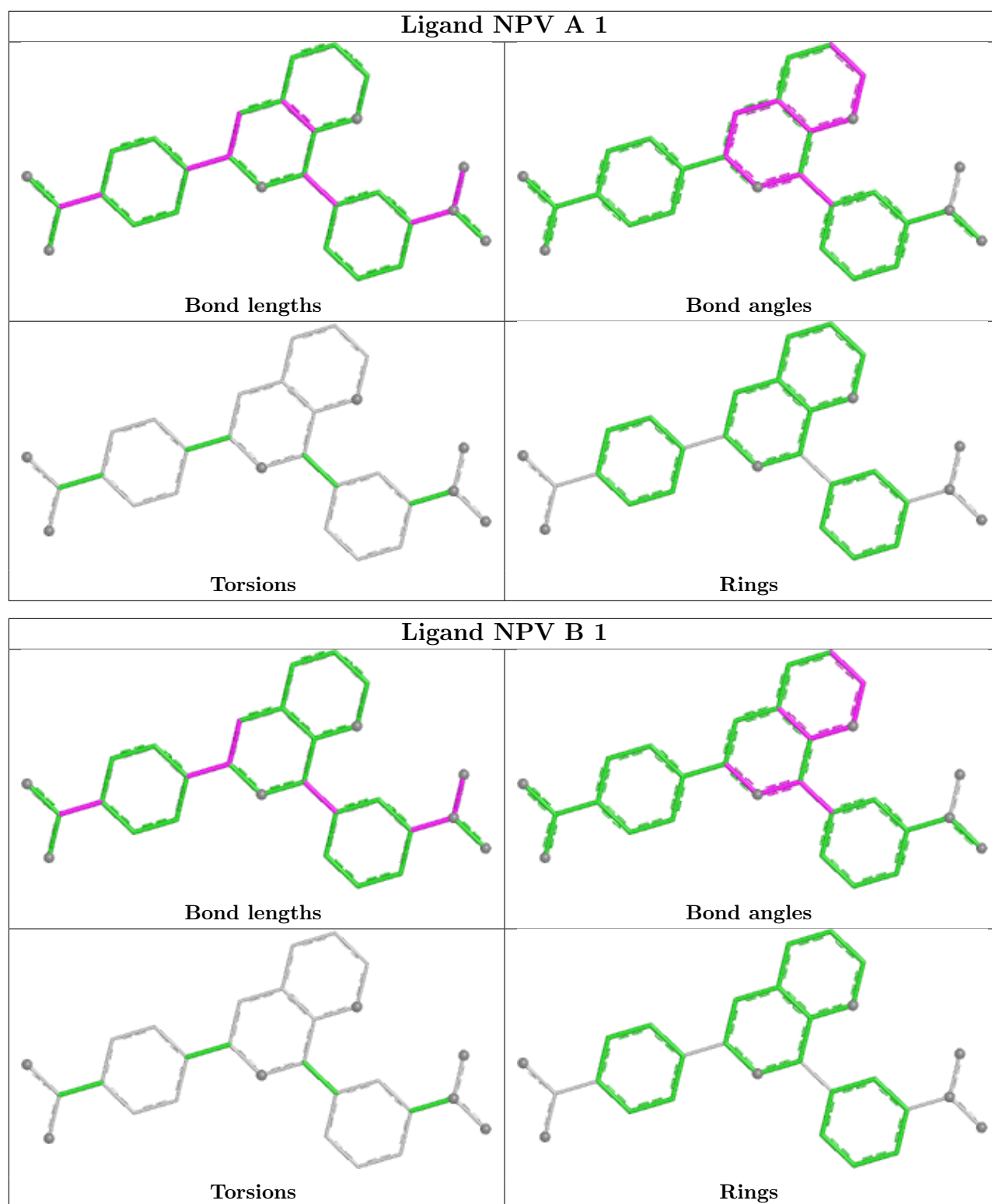
There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	NPV	1	0
2	B	1	NPV	1	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 ⓘ

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	326/328 (99%)	0.57	37 (11%)	10 15	11, 21, 41, 50	0
1	B	325/328 (99%)	1.63	121 (37%)	1 1	13, 29, 51, 58	0
All	All	651/656 (99%)	1.10	158 (24%)	2 3	11, 24, 46, 58	0

All (158) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	297	VAL	6.3
1	A	86	THR	5.7
1	B	354	ILE	5.7
1	B	93	ALA	5.5
1	B	293	THR	5.4
1	B	183	ALA	5.3
1	B	92	LEU	5.2
1	B	349	GLU	5.0
1	B	292	VAL	4.8
1	B	296	GLY	4.8
1	B	353	GLU	4.8
1	B	392	ALA	4.8
1	B	391	ASP	4.8
1	B	350	ARG	4.4
1	A	294	SER	4.4
1	B	279	LEU	4.2
1	B	184	VAL	4.1
1	A	410	ILE	4.1
1	B	286	MET	4.1
1	B	298	LEU	4.0
1	B	396	LEU	4.0
1	B	289	THR	4.0
1	B	99	VAL	4.0
1	B	294	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	348	ARG	4.0
1	B	178	THR	4.0
1	B	91	VAL	3.9
1	A	292	VAL	3.9
1	B	181	LEU	3.8
1	B	179	PRO	3.8
1	B	386	ASP	3.7
1	B	135	PHE	3.7
1	B	385	ALA	3.7
1	A	351	GLY	3.7
1	B	185	PHE	3.7
1	B	287	VAL	3.7
1	B	363	ALA	3.7
1	A	90	ASP	3.6
1	B	125	ILE	3.6
1	B	137	ILE	3.6
1	B	412	GLN	3.6
1	B	132	LEU	3.6
1	B	344	GLY	3.5
1	A	363	ALA	3.5
1	B	295	SER	3.5
1	B	395	ILE	3.5
1	A	342	ARG	3.4
1	B	302	ASN	3.4
1	B	351	GLY	3.4
1	B	176	LEU	3.4
1	A	103	GLY	3.3
1	B	390	PRO	3.3
1	B	358	CYS	3.3
1	B	138	PRO	3.3
1	B	96	LEU	3.2
1	A	411	PRO	3.2
1	B	139	VAL	3.2
1	B	388	VAL	3.2
1	B	104	LEU	3.2
1	B	131	LEU	3.2
1	A	353	GLU	3.2
1	B	109	ILE	3.2
1	B	340	PHE	3.1
1	B	177	SER	3.1
1	B	180	ALA	3.1
1	B	394	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	281	ALA	3.1
1	B	389	HIS	3.1
1	B	127	GLN	3.1
1	B	130	ASP	3.0
1	B	342	ARG	3.0
1	B	98	ASP	3.0
1	B	282	ASP	3.0
1	B	285	THR	3.0
1	A	350	ARG	2.9
1	B	214	ASN	2.9
1	B	120	VAL	2.9
1	A	88	GLN	2.9
1	B	375	TYR	2.9
1	B	186	THR	2.9
1	B	123	HIS	2.9
1	A	357	MET	2.9
1	A	87	GLU	2.9
1	A	218	GLU	2.9
1	A	362	ASN	2.9
1	B	352	MET	2.9
1	B	345	ASP	2.8
1	B	94	LYS	2.8
1	B	306	ARG	2.8
1	B	299	LEU	2.8
1	A	375	TYR	2.8
1	A	409	THR	2.8
1	B	397	ASP	2.7
1	B	141	THR	2.7
1	B	308	GLN	2.7
1	B	357	MET	2.7
1	A	293	THR	2.7
1	A	389	HIS	2.7
1	A	354	ILE	2.7
1	B	356	PRO	2.7
1	B	288	GLU	2.6
1	B	364	SER	2.6
1	B	362	ASN	2.6
1	A	356	PRO	2.6
1	B	140	ASP	2.6
1	B	384	TRP	2.5
1	B	147	MET	2.5
1	B	174	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	295	SER	2.5
1	A	299	LEU	2.5
1	B	303	TYR	2.5
1	B	310	LEU	2.5
1	A	127	GLN	2.5
1	B	142	LEU	2.5
1	B	126	PHE	2.5
1	B	359	ASP	2.5
1	B	361	HIS	2.5
1	B	283	LEU	2.4
1	B	341	PHE	2.4
1	B	300	LEU	2.4
1	B	398	THR	2.4
1	B	355	SER	2.4
1	B	100	ASN	2.4
1	B	90	ASP	2.4
1	B	252	LEU	2.4
1	B	134	THR	2.3
1	B	338	GLU	2.3
1	B	346	ARG	2.3
1	B	215	THR	2.3
1	B	112	LEU	2.3
1	B	143	ILE	2.3
1	A	349	GLU	2.3
1	B	251	ASN	2.2
1	A	361	HIS	2.2
1	B	89	GLU	2.2
1	B	393	GLN	2.2
1	A	298	LEU	2.2
1	B	182	GLU	2.2
1	B	102	TRP	2.2
1	A	151	ASP	2.2
1	A	352	MET	2.2
1	A	296	GLY	2.1
1	A	364	SER	2.1
1	A	359	ASP	2.1
1	B	101	LYS	2.1
1	B	124	THR	2.1
1	B	376	ILE	2.1
1	B	304	SER	2.1
1	A	219	LEU	2.1
1	B	290	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	348	ARG	2.0
1	B	305	ASP	2.0
1	B	108	ARG	2.0
1	B	383	THR	2.0
1	B	97	GLU	2.0
1	B	400	GLU	2.0
1	A	281	ALA	2.0
1	B	399	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

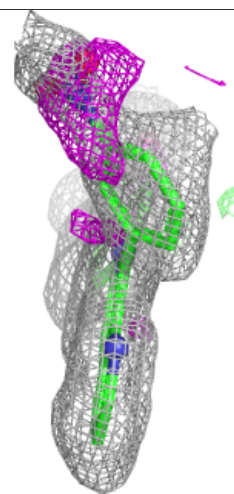
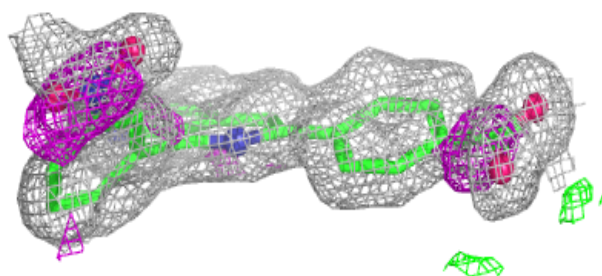
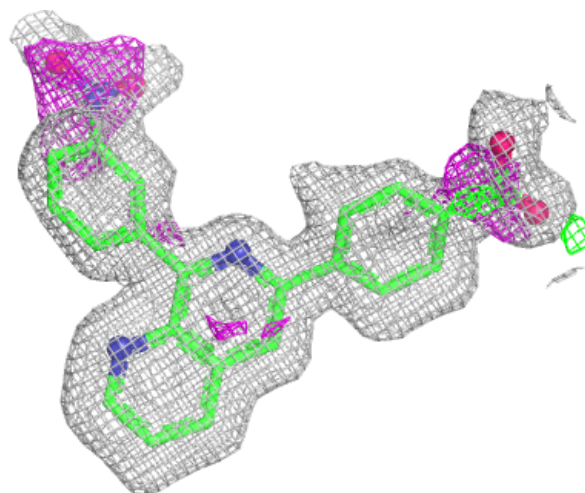
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

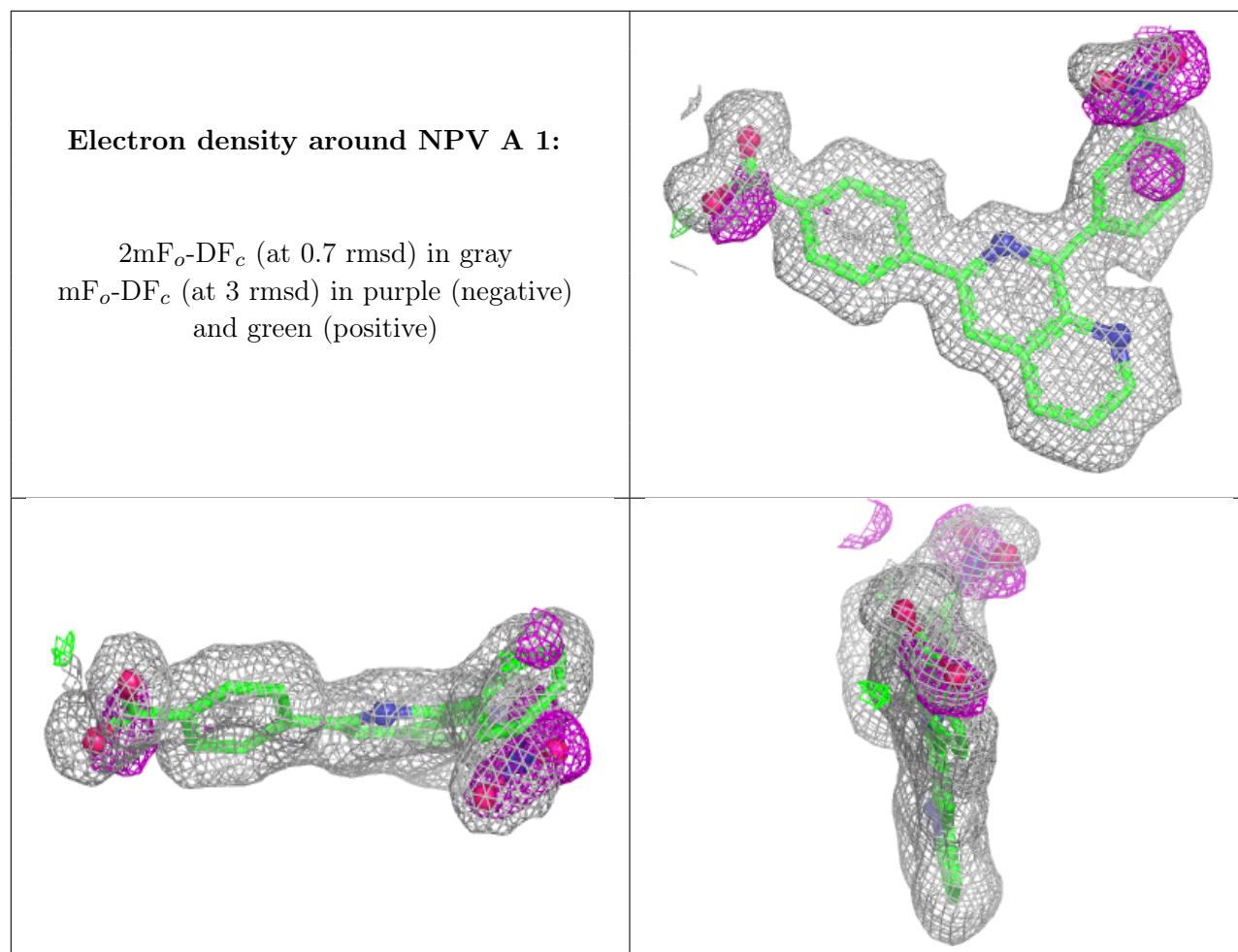
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NPV	B	1	28/28	0.84	0.13	22,25,40,41	0
2	NPV	A	1	28/28	0.89	0.10	17,21,38,41	0
3	ZN	A	501	1/1	0.99	0.01	14,14,14,14	0
3	ZN	B	501	1/1	0.99	0.03	18,18,18,18	0
4	MG	A	502	1/1	0.99	0.02	12,12,12,12	0
4	MG	B	502	1/1	0.99	0.03	14,14,14,14	0

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 NPV B 1:

$2mF_o-DF_c$ (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.