



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:25 PM UTC

PDB ID : 3RHH / pdb_00003rhh
Title : Crystal structure of NADP-dependent glyceraldehyde-3-phosphate dehydrogenase from *Bacillus halodurans* C-125 complexed with NADP
Authors : Malashkevich, V.N.; Toro, R.; Seidel, R.; Garrett, S.; Foti, R.; Almo, S.C.; New York Structural Genomics Research Consortium (NYSGRG)
Deposited on : 2011-04-11
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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

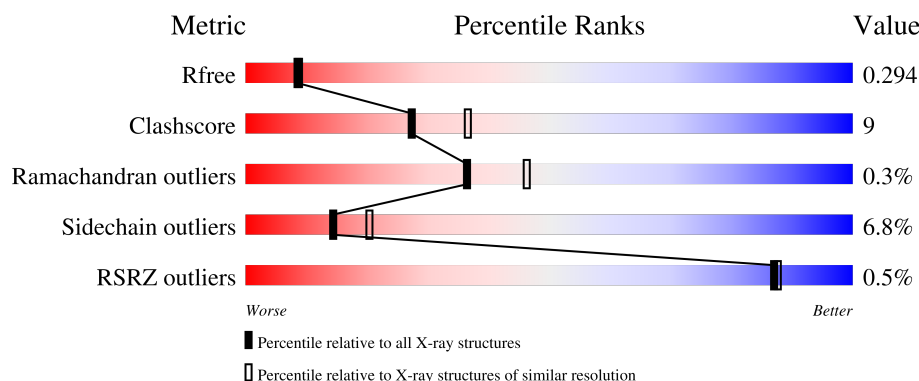
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	505	 71% 21% • 6%
1	B	505	 69% 23% • 6%
1	C	505	 78% 15% • 6%
1	D	505	 2% 70% 22% • 5%

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 14718 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 NADP-dependent glyceraldehyde-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	475	Total	C	N	O	S	0	0	0
			3573	2251	621	694	7			
1	B	476	Total	C	N	O	S	0	0	0
			3578	2254	622	695	7			
1	C	476	Total	C	N	O	S	0	2	0
			3594	2265	623	699	7			
1	D	480	Total	C	N	O	S	0	1	0
			3624	2286	627	704	7			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q9KAQ0
A	2	VAL	-	expression tag	UNP Q9KAQ0
A	484	ALA	-	expression tag	UNP Q9KAQ0
A	485	GLU	-	expression tag	UNP Q9KAQ0
A	486	ASN	-	expression tag	UNP Q9KAQ0
A	487	LEU	-	expression tag	UNP Q9KAQ0
A	488	TYR	-	expression tag	UNP Q9KAQ0
A	489	PHE	-	expression tag	UNP Q9KAQ0
A	490	GLN	-	expression tag	UNP Q9KAQ0
A	491	SER	-	expression tag	UNP Q9KAQ0
A	492	HIS	-	expression tag	UNP Q9KAQ0
A	493	HIS	-	expression tag	UNP Q9KAQ0
A	494	HIS	-	expression tag	UNP Q9KAQ0
A	495	HIS	-	expression tag	UNP Q9KAQ0
A	496	HIS	-	expression tag	UNP Q9KAQ0
A	497	HIS	-	expression tag	UNP Q9KAQ0
A	498	TRP	-	expression tag	UNP Q9KAQ0
A	499	SER	-	expression tag	UNP Q9KAQ0
A	500	HIS	-	expression tag	UNP Q9KAQ0
A	501	PRO	-	expression tag	UNP Q9KAQ0
A	502	GLN	-	expression tag	UNP Q9KAQ0

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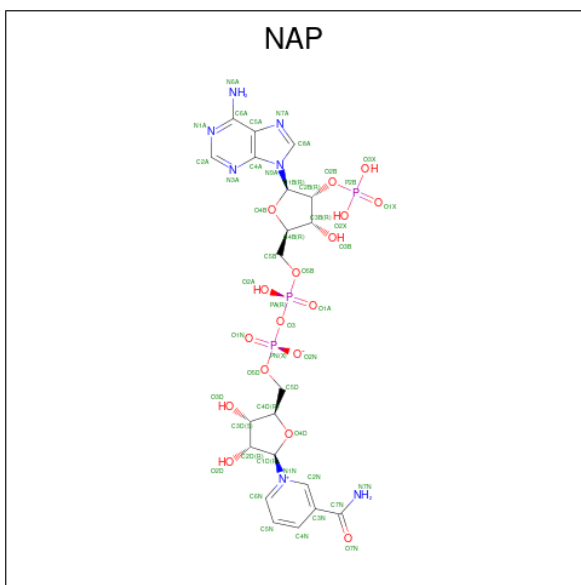
Chain	Residue	Modelled	Actual	Comment	Reference
A	503	PHE	-	expression tag	UNP Q9KAQ0
A	504	GLU	-	expression tag	UNP Q9KAQ0
A	505	LYS	-	expression tag	UNP Q9KAQ0
B	1	MET	-	expression tag	UNP Q9KAQ0
B	2	VAL	-	expression tag	UNP Q9KAQ0
B	484	ALA	-	expression tag	UNP Q9KAQ0
B	485	GLU	-	expression tag	UNP Q9KAQ0
B	486	ASN	-	expression tag	UNP Q9KAQ0
B	487	LEU	-	expression tag	UNP Q9KAQ0
B	488	TYR	-	expression tag	UNP Q9KAQ0
B	489	PHE	-	expression tag	UNP Q9KAQ0
B	490	GLN	-	expression tag	UNP Q9KAQ0
B	491	SER	-	expression tag	UNP Q9KAQ0
B	492	HIS	-	expression tag	UNP Q9KAQ0
B	493	HIS	-	expression tag	UNP Q9KAQ0
B	494	HIS	-	expression tag	UNP Q9KAQ0
B	495	HIS	-	expression tag	UNP Q9KAQ0
B	496	HIS	-	expression tag	UNP Q9KAQ0
B	497	HIS	-	expression tag	UNP Q9KAQ0
B	498	TRP	-	expression tag	UNP Q9KAQ0
B	499	SER	-	expression tag	UNP Q9KAQ0
B	500	HIS	-	expression tag	UNP Q9KAQ0
B	501	PRO	-	expression tag	UNP Q9KAQ0
B	502	GLN	-	expression tag	UNP Q9KAQ0
B	503	PHE	-	expression tag	UNP Q9KAQ0
B	504	GLU	-	expression tag	UNP Q9KAQ0
B	505	LYS	-	expression tag	UNP Q9KAQ0
C	1	MET	-	expression tag	UNP Q9KAQ0
C	2	VAL	-	expression tag	UNP Q9KAQ0
C	484	ALA	-	expression tag	UNP Q9KAQ0
C	485	GLU	-	expression tag	UNP Q9KAQ0
C	486	ASN	-	expression tag	UNP Q9KAQ0
C	487	LEU	-	expression tag	UNP Q9KAQ0
C	488	TYR	-	expression tag	UNP Q9KAQ0
C	489	PHE	-	expression tag	UNP Q9KAQ0
C	490	GLN	-	expression tag	UNP Q9KAQ0
C	491	SER	-	expression tag	UNP Q9KAQ0
C	492	HIS	-	expression tag	UNP Q9KAQ0
C	493	HIS	-	expression tag	UNP Q9KAQ0
C	494	HIS	-	expression tag	UNP Q9KAQ0
C	495	HIS	-	expression tag	UNP Q9KAQ0
C	496	HIS	-	expression tag	UNP Q9KAQ0

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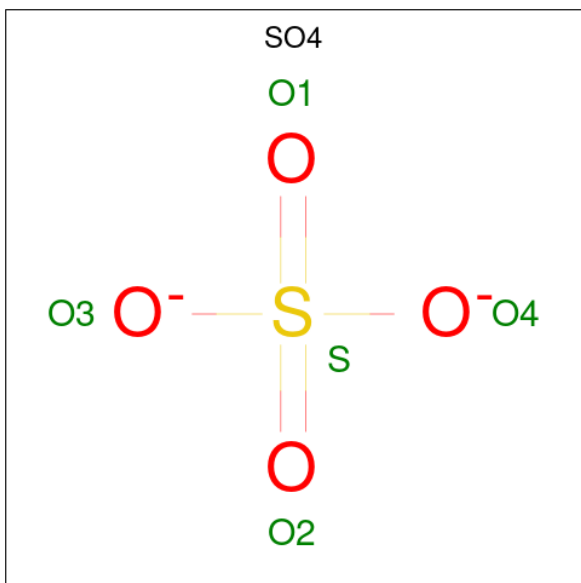
Chain	Residue	Modelled	Actual	Comment	Reference
C	497	HIS	-	expression tag	UNP Q9KAQ0
C	498	TRP	-	expression tag	UNP Q9KAQ0
C	499	SER	-	expression tag	UNP Q9KAQ0
C	500	HIS	-	expression tag	UNP Q9KAQ0
C	501	PRO	-	expression tag	UNP Q9KAQ0
C	502	GLN	-	expression tag	UNP Q9KAQ0
C	503	PHE	-	expression tag	UNP Q9KAQ0
C	504	GLU	-	expression tag	UNP Q9KAQ0
C	505	LYS	-	expression tag	UNP Q9KAQ0
D	1	MET	-	expression tag	UNP Q9KAQ0
D	2	VAL	-	expression tag	UNP Q9KAQ0
D	484	ALA	-	expression tag	UNP Q9KAQ0
D	485	GLU	-	expression tag	UNP Q9KAQ0
D	486	ASN	-	expression tag	UNP Q9KAQ0
D	487	LEU	-	expression tag	UNP Q9KAQ0
D	488	TYR	-	expression tag	UNP Q9KAQ0
D	489	PHE	-	expression tag	UNP Q9KAQ0
D	490	GLN	-	expression tag	UNP Q9KAQ0
D	491	SER	-	expression tag	UNP Q9KAQ0
D	492	HIS	-	expression tag	UNP Q9KAQ0
D	493	HIS	-	expression tag	UNP Q9KAQ0
D	494	HIS	-	expression tag	UNP Q9KAQ0
D	495	HIS	-	expression tag	UNP Q9KAQ0
D	496	HIS	-	expression tag	UNP Q9KAQ0
D	497	HIS	-	expression tag	UNP Q9KAQ0
D	498	TRP	-	expression tag	UNP Q9KAQ0
D	499	SER	-	expression tag	UNP Q9KAQ0
D	500	HIS	-	expression tag	UNP Q9KAQ0
D	501	PRO	-	expression tag	UNP Q9KAQ0
D	502	GLN	-	expression tag	UNP Q9KAQ0
D	503	PHE	-	expression tag	UNP Q9KAQ0
D	504	GLU	-	expression tag	UNP Q9KAQ0
D	505	LYS	-	expression tag	UNP Q9KAQ0

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	B	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	C	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	D	1	Total 31	C 10	N 5	O 13	P 3	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

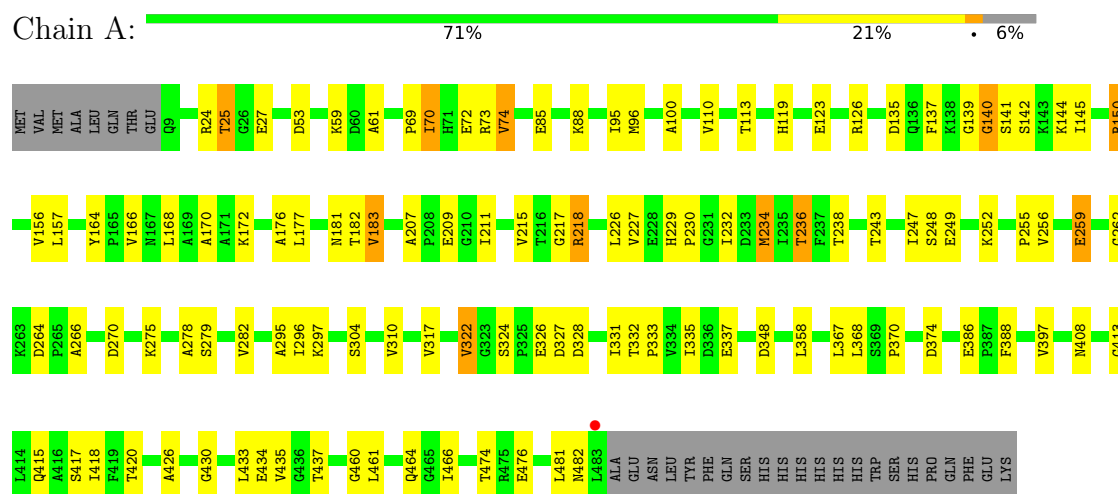
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	60	Total O 60 60	0	0
4	B	65	Total O 65 65	0	0
4	C	37	Total O 37 37	0	0
4	D	43	Total O 43 43	0	0

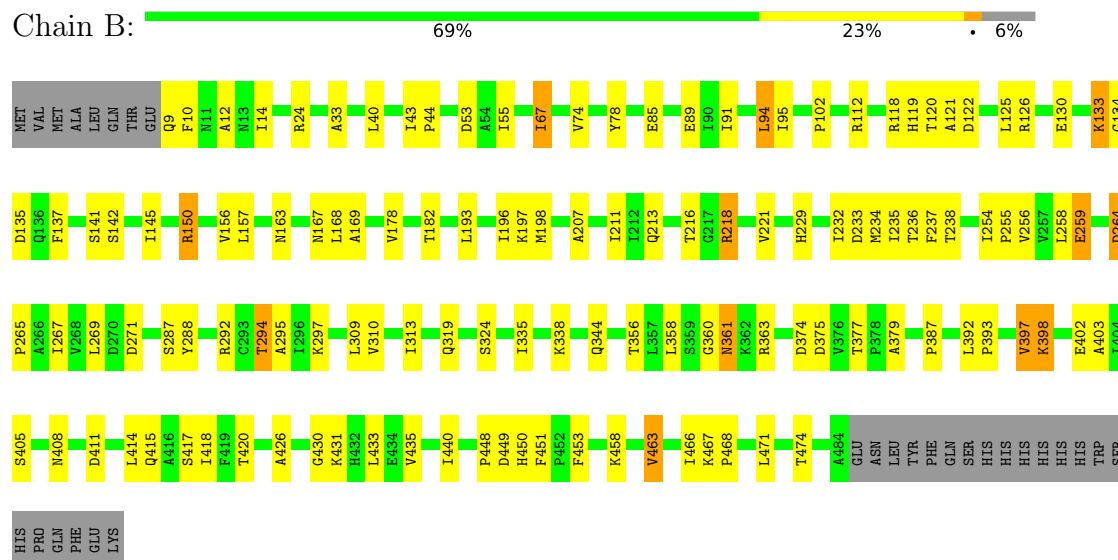
3 Residue-property plots

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: NADP-dependent glyceraldehyde-3-phosphate dehydrogenase

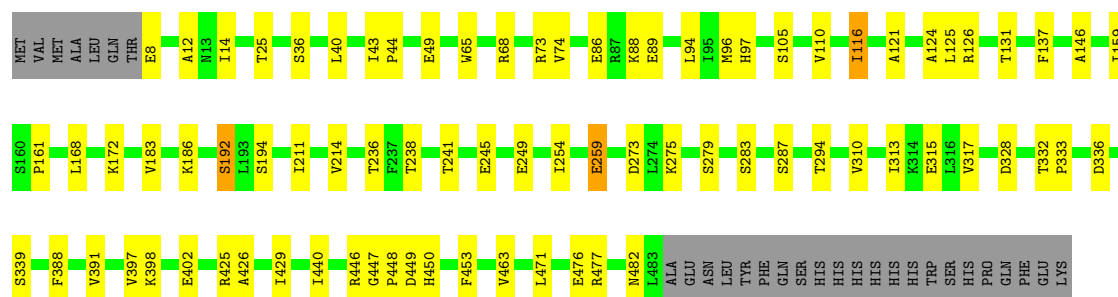


- Molecule 1: NADP-dependent glyceraldehyde-3-phosphate dehydrogenase



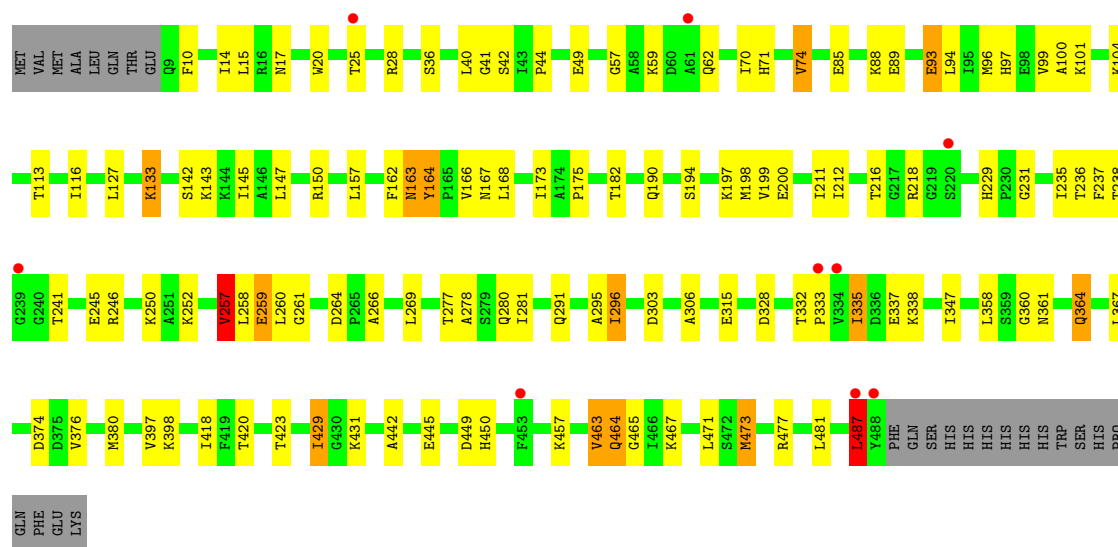
- Molecule 1: NADP-dependent glyceraldehyde-3-phosphate dehydrogenase

Chain C:



- Molecule 1: NADP-dependent glyceraldehyde-3-phosphate dehydrogenase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.65Å 130.75Å 86.92Å 90.00° 102.69° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.3 (20.00-2.30) 95.1 (20.00-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.30Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.232 , 0.297 0.233 , 0.294	Depositor DCC
R_{free} test set	3880 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	45.8	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.030 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14718	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.59% 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: NAP, SO4

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	0.62	0/3627	0.94	3/4919 (0.1%)
1	B	0.70	0/3632	0.98	8/4926 (0.2%)
1	C	0.59	0/3654	0.92	2/4955 (0.0%)
1	D	0.63	2/3683 (0.1%)	0.94	9/4996 (0.2%)
All	All	0.63	2/14596 (0.0%)	0.95	22/19796 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	464	GLN	C-O	-5.79	1.17	1.24
1	D	257	VAL	CA-CB	-5.66	1.47	1.54

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	ILE	N-CA-C	7.92	118.69	110.62
1	D	465	GLY	N-CA-C	-7.46	102.90	111.85
1	B	33	ALA	CA-C-N	6.67	126.36	119.56
1	B	33	ALA	C-N-CA	6.67	126.36	119.56
1	D	261	GLY	N-CA-C	6.57	121.05	112.25
1	D	164	TYR	CA-C-N	6.49	126.72	119.32
1	D	164	TYR	C-N-CA	6.49	126.72	119.32
1	D	487	LEU	N-CA-C	-6.19	105.63	113.43
1	B	335	ILE	N-CA-C	6.07	116.81	110.62
1	A	460	GLY	N-CA-C	5.95	119.44	110.75
1	C	254	ILE	N-CA-C	5.82	113.96	109.19
1	B	324	SER	CA-C-N	5.79	125.65	119.28
1	B	324	SER	C-N-CA	5.79	125.65	119.28
1	B	264	ASP	CA-C-N	5.78	126.23	119.99
1	B	264	ASP	C-N-CA	5.78	126.23	119.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	335	ILE	N-CA-C	5.70	116.35	110.36
1	D	457	LYS	CB-CA-C	-5.37	110.37	116.54
1	A	183	VAL	N-CA-C	5.26	115.48	108.11
1	B	435	VAL	N-CA-C	5.26	114.94	108.53
1	D	464	GLN	N-CA-C	5.11	115.03	108.34
1	C	183	VAL	N-CA-C	5.04	115.38	108.12
1	D	173	ILE	N-CA-C	5.02	115.23	110.42

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	3573	0	3656	69	0
1	B	3578	0	3661	87	0
1	C	3594	0	3681	45	0
1	D	3624	0	3702	78	0
2	A	31	0	11	1	0
2	B	31	0	11	0	0
2	C	31	0	11	0	0
2	D	31	0	11	0	0
3	A	5	0	0	0	0
3	B	5	0	0	1	0
3	C	5	0	0	0	0
3	D	5	0	0	1	0
4	A	60	0	0	3	0
4	B	65	0	0	8	0
4	C	37	0	0	2	0
4	D	43	0	0	4	0
All	All	14718	0	14744	265	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 9.

All (265) 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:238:THR:HG23	1:B:259:GLU:HG3	1.48	0.95
1:D:259:GLU:CD	1:D:464:GLN:HE22	1.75	0.94
1:D:257:VAL:HG12	1:D:257:VAL:O	1.64	0.94
1:B:218:ARG:HG3	1:B:221:VAL:HB	1.52	0.91
1:B:294:THR:HG22	1:B:414:LEU:HD21	1.54	0.88
1:D:266:ALA:HB2	1:D:296:ILE:HD13	1.57	0.86
1:B:207:ALA:HB1	1:B:211:ILE:HD11	1.58	0.85
1:A:259:GLU:HB3	1:A:464:GLN:HE22	1.46	0.81
1:B:397:VAL:HG22	1:B:402:GLU:HB3	1.61	0.80
1:B:238:THR:HG23	1:B:259:GLU:CG	2.12	0.80
1:B:236:THR:HG21	4:B:559:HOH:O	1.82	0.77
1:C:447:GLY:HA3	4:C:523:HOH:O	1.86	0.75
1:A:156:VAL:HG13	1:A:234:MET:CE	2.19	0.72
1:D:229:HIS:HD2	1:D:231:GLY:H	1.35	0.72
1:D:259:GLU:CD	1:D:464:GLN:NE2	2.46	0.72
1:D:277:THR:O	1:D:281:ILE:HG12	1.90	0.71
1:C:74:VAL:HG21	1:C:125:LEU:HD13	1.72	0.71
1:D:113:THR:OG1	1:D:167:ASN:HA	1.90	0.71
1:D:163:ASN:ND2	1:D:291:GLN:O	2.23	0.71
1:D:182:THR:HG22	1:D:211:ILE:HA	1.75	0.69
1:A:211:ILE:HB	4:A:517:HOH:O	1.93	0.68
1:D:96:MET:O	1:D:100:ALA:HA	1.92	0.68
1:B:430:GLY:O	1:C:477:ARG:NH2	2.26	0.68
1:B:156:VAL:HG13	1:B:234:MET:CE	2.24	0.68
1:D:10:PHE:O	1:D:41:GLY:HA3	1.94	0.68
1:A:418:ILE:HD11	1:A:433:LEU:HD11	1.75	0.68
1:A:141:SER:HB3	1:A:144:LYS:HE2	1.76	0.67
1:D:40:LEU:HD11	1:D:97:HIS:HB3	1.76	0.67
1:A:332:THR:HB	1:A:333:PRO:HD2	1.76	0.67
1:C:238:THR:HG23	1:C:259:GLU:CG	2.24	0.67
1:A:156:VAL:HG13	1:A:234:MET:HE2	1.76	0.66
1:D:264:ASP:HB2	1:D:295:ALA:O	1.95	0.66
1:B:218:ARG:HB3	4:B:509:HOH:O	1.95	0.66
1:A:168:LEU:HD13	1:A:238:THR:HG21	1.79	0.65
1:D:259:GLU:CG	1:D:464:GLN:HE22	2.10	0.65
1:A:358:LEU:HD11	1:A:374:ASP:HB2	1.77	0.65
1:B:156:VAL:HG13	1:B:234:MET:HE2	1.79	0.65
1:B:24:ARG:NH2	1:B:53:ASP:OD2	2.29	0.64
1:B:130:GLU:HB3	4:D:527:HOH:O	1.98	0.64
1:A:70:ILE:HD12	1:A:73:ARG:NH2	2.14	0.63
1:A:264:ASP:HB2	1:A:295:ALA:O	2.00	0.62
1:C:172[B]:LYS:NZ	1:C:238:THR:OG1	2.25	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:175:PRO:HG2	1:D:473:MET:HE3	1.81	0.62
1:A:207:ALA:HB1	4:A:517:HOH:O	1.99	0.61
1:B:358:LEU:HD11	1:B:374:ASP:HB2	1.82	0.61
1:C:65:TRP:HB2	1:C:211:ILE:HD11	1.82	0.61
1:D:164:TYR:HB3	1:D:167:ASN:HB3	1.80	0.61
1:B:218:ARG:HG3	1:B:221:VAL:CB	2.26	0.61
1:A:96:MET:HE1	1:A:326:GLU:HG3	1.82	0.61
1:B:264:ASP:HB2	1:B:295:ALA:O	2.01	0.60
1:D:303:ASP:HA	1:D:306:ALA:HB2	1.84	0.60
1:A:420:THR:CG2	1:A:426:ALA:HB2	2.31	0.60
1:C:168:LEU:HD13	1:C:238:THR:HG21	1.84	0.60
1:A:156:VAL:HB	1:A:183:VAL:HG22	1.83	0.60
1:D:347:ILE:HG21	1:D:361:ASN:OD1	2.03	0.59
1:D:237:PHE:HB3	1:D:258:LEU:HD23	1.84	0.59
1:B:271:ASP:OD1	1:B:420:THR:OG1	2.20	0.58
1:D:229:HIS:CD2	1:D:231:GLY:H	2.19	0.58
1:D:133:LYS:HD3	1:D:145:ILE:HD12	1.84	0.58
1:A:182:THR:HG22	1:A:211:ILE:HA	1.86	0.57
1:D:376:VAL:HA	1:D:380:MET:SD	2.44	0.57
1:D:236:THR:HA	1:D:257:VAL:HG12	1.86	0.57
1:B:137:PHE:HB2	1:C:450:HIS:ND1	2.19	0.57
1:D:257:VAL:O	1:D:257:VAL:CG1	2.30	0.57
1:B:78:TYR:OH	1:B:122:ASP:OD2	2.23	0.57
1:C:259:GLU:HG3	1:C:259:GLU:O	2.04	0.57
1:A:61:ALA:HB1	1:A:211:ILE:HD13	1.87	0.56
1:C:131:THR:HA	1:C:146:ALA:O	2.05	0.56
1:B:411:ASP:HB3	1:B:458:LYS:HG3	1.86	0.56
1:D:59:LYS:HE3	4:D:514:HOH:O	2.05	0.56
1:B:267:ILE:HG12	1:B:403:ALA:HB1	1.88	0.55
1:B:463:VAL:HG21	1:C:476:GLU:OE1	2.07	0.55
1:C:43:ILE:HG21	1:C:192:SER:HB2	1.89	0.55
1:C:172[A]:LYS:HD3	1:C:236:THR:HB	1.89	0.55
1:B:134:GLY:HA3	1:B:141:SER:O	2.07	0.55
1:D:418:ILE:CD1	1:D:429:ILE:HG22	2.36	0.55
1:C:259:GLU:OE1	1:C:446:ARG:NH2	2.39	0.55
1:A:317:VAL:HG13	1:A:370:PRO:HB2	1.88	0.55
1:C:238:THR:HG23	1:C:259:GLU:HG2	1.88	0.55
1:B:157:LEU:HD23	1:B:235:ILE:HG12	1.89	0.55
1:B:287:SER:HB2	1:B:292:ARG:HH11	1.72	0.55
1:B:310:VAL:CG1	1:B:358:LEU:HD13	2.37	0.55
1:A:415:GLN:HG2	1:A:437:THR:HB	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:157:LEU:HD23	1:D:235:ILE:CD1	2.38	0.54
1:B:408:ASN:HB3	4:B:532:HOH:O	2.07	0.54
1:B:119:HIS:HD2	1:B:466:ILE:HG21	1.73	0.54
1:B:12:ALA:HB1	1:B:196:ILE:HD13	1.90	0.54
1:A:145:ILE:HD11	1:A:481:LEU:HD12	1.89	0.54
4:B:515:HOH:O	1:C:482:ASN:HB2	2.08	0.54
1:D:70:ILE:O	1:D:74:VAL:HG13	2.07	0.54
1:B:182:THR:CG2	1:B:211:ILE:HA	2.38	0.53
1:A:249:GLU:HB2	1:D:252:LYS:HD3	1.90	0.53
1:B:450:HIS:ND1	1:C:137:PHE:HB2	2.24	0.53
1:C:425:ARG:HE	1:D:487:LEU:HD13	1.73	0.53
1:B:356:THR:OG1	1:B:375:ASP:OD2	2.25	0.53
1:D:10:PHE:N	1:D:40:LEU:O	2.42	0.53
1:D:166:VAL:HG13	1:D:198:MET:HG3	1.89	0.53
1:A:88:LYS:HA	1:A:110:VAL:HG11	1.91	0.53
1:B:168:LEU:HD13	1:B:238:THR:HG21	1.91	0.52
1:B:182:THR:HG22	1:B:211:ILE:HA	1.90	0.52
1:D:358:LEU:HD11	1:D:374:ASP:HB2	1.92	0.52
1:D:241:THR:HA	1:D:260:LEU:HD22	1.90	0.52
1:A:266:ALA:HB2	1:A:296:ILE:HD12	1.93	0.51
1:B:449:ASP:O	1:B:466:ILE:HG12	2.11	0.51
1:D:182:THR:CG2	1:D:211:ILE:HA	2.41	0.51
1:D:246:ARG:O	1:D:250:LYS:HG3	2.12	0.50
1:B:133:LYS:HE3	1:B:142:SER:O	2.11	0.50
1:A:156:VAL:HG21	1:A:176:ALA:HB1	1.94	0.50
1:A:408:ASN:CG	1:A:434:GLU:HG3	2.37	0.50
1:D:418:ILE:HD13	1:D:429:ILE:HG22	1.93	0.50
1:B:150:ARG:HD3	1:B:474:THR:OG1	2.12	0.50
1:D:278:ALA:HB3	4:D:531:HOH:O	2.11	0.49
1:D:259:GLU:CG	1:D:464:GLN:NE2	2.73	0.49
1:A:70:ILE:O	1:A:74:VAL:HG13	2.13	0.49
1:A:164:TYR:HB2	1:A:168:LEU:HD12	1.94	0.49
1:B:361:ASN:CG	1:B:361:ASN:O	2.55	0.49
1:A:119:HIS:HD2	1:A:466:ILE:HG21	1.77	0.49
1:B:43:ILE:HD12	1:B:216:THR:HB	1.95	0.49
1:B:398:LYS:HB2	1:B:402:GLU:OE1	2.13	0.49
1:C:426:ALA:HB1	1:C:440:ILE:HD12	1.95	0.49
1:D:280:GLN:OE1	1:D:442:ALA:HA	2.13	0.49
1:B:163:ASN:HB2	1:B:288:TYR:HE2	1.78	0.49
1:B:294:THR:HG22	1:B:414:LEU:CD2	2.36	0.49
1:B:259:GLU:HG3	1:B:259:GLU:O	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:ARG:NH1	3:B:507:SO4:O4	2.38	0.48
1:C:259:GLU:CD	1:C:453:PHE:HE2	2.21	0.48
1:C:241:THR:O	1:C:245:GLU:HG3	2.14	0.48
1:D:93:GLU:HA	1:D:96:MET:HG2	1.94	0.48
1:D:259:GLU:OE2	1:D:464:GLN:OE1	2.31	0.48
1:B:417:SER:O	1:B:418:ILE:HD13	2.13	0.48
1:C:447:GLY:CA	1:C:448:PRO:C	2.87	0.48
1:C:332:THR:HB	1:C:333:PRO:CD	2.43	0.48
1:D:99:VAL:HG11	1:D:162:PHE:O	2.14	0.48
1:A:95:ILE:HD11	1:A:166:VAL:CG2	2.44	0.47
1:C:116:ILE:HD11	1:C:449:ASP:OD1	2.14	0.47
1:D:190:GLN:HG2	1:D:335:ILE:HD11	1.96	0.47
1:A:157:LEU:HB2	1:A:232:ILE:HD12	1.95	0.47
1:B:135:ASP:HB2	1:D:71:HIS:HB3	1.95	0.47
1:A:227:VAL:HG21	1:A:247:ILE:HG12	1.97	0.47
1:D:17:ASN:OD1	1:D:57:GLY:HA3	2.14	0.47
1:A:137:PHE:HB2	1:D:450:HIS:ND1	2.30	0.47
1:A:430:GLY:O	1:D:477:ARG:NH2	2.48	0.47
1:A:217:GLY:O	1:A:218:ARG:C	2.57	0.47
1:D:85:GLU:O	1:D:88:LYS:HG2	2.15	0.47
1:D:246:ARG:HH11	1:D:250:LYS:HZ3	1.62	0.46
1:A:172:LYS:HG2	1:A:236:THR:HG21	1.97	0.46
1:D:104:LYS:HG3	4:D:550:HOH:O	2.15	0.46
1:B:102:PRO:HB2	4:B:550:HOH:O	2.15	0.46
1:A:172:LYS:HD3	1:A:236:THR:HB	1.98	0.46
1:A:135:ASP:C	1:A:137:PHE:H	2.24	0.46
1:B:338:LYS:HG3	4:B:524:HOH:O	2.16	0.46
1:C:14:ILE:HD13	1:C:214:VAL:O	2.15	0.46
1:D:28:ARG:HH22	1:D:44:PRO:HD3	1.80	0.46
1:D:246:ARG:HH11	1:D:250:LYS:NZ	2.13	0.46
1:B:448:PRO:HD2	1:B:451:PHE:CD2	2.51	0.45
1:C:74:VAL:HG13	1:C:121:ALA:HB1	1.97	0.45
1:A:278:ALA:O	1:A:282:VAL:HG23	2.16	0.45
1:D:197:LYS:HD3	1:D:197:LYS:HA	1.84	0.45
1:A:139:GLY:O	1:A:140:GLY:C	2.60	0.45
1:B:237:PHE:HB3	1:B:258:LEU:HD23	1.99	0.45
1:A:176:ALA:O	1:A:181:ASN:HB2	2.17	0.45
1:A:322:VAL:HB	1:A:331:ILE:HB	1.97	0.45
1:C:429:ILE:HG12	1:D:487:LEU:HD22	1.99	0.45
1:A:150:ARG:HG2	1:A:474:THR:HB	1.99	0.45
1:A:177:LEU:HD22	1:A:211:ILE:HG22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:GLU:OE2	1:B:118:ARG:NH2	2.47	0.45
1:C:86:GLU:HB2	4:C:533:HOH:O	2.15	0.45
1:B:310:VAL:HG11	1:B:358:LEU:CD1	2.47	0.45
1:A:229:HIS:HA	1:A:230:PRO:HD3	1.88	0.45
1:B:91:ILE:O	1:B:95:ILE:HG13	2.17	0.45
1:C:336:ASP:O	1:C:339:SER:HB3	2.17	0.45
1:D:14:ILE:HG21	1:D:216:THR:OG1	2.17	0.45
1:D:333:PRO:HD3	1:D:367:LEU:HD13	1.98	0.44
1:B:67:ILE:HD13	1:B:67:ILE:HA	1.83	0.44
1:B:398:LYS:HB3	1:B:398:LYS:HE3	1.42	0.44
1:B:418:ILE:HD11	1:B:433:LEU:HD11	1.99	0.44
1:C:313:ILE:O	1:C:317:VAL:HG23	2.18	0.44
1:B:156:VAL:HG13	1:B:234:MET:HE1	2.00	0.44
1:B:218:ARG:HA	4:B:509:HOH:O	2.16	0.44
1:B:233:ASP:O	1:B:255:PRO:HD2	2.17	0.44
1:D:157:LEU:HD23	1:D:235:ILE:HD13	1.98	0.44
1:D:164:TYR:HB2	1:D:168:LEU:HD12	1.99	0.44
1:B:310:VAL:CG1	1:B:358:LEU:CD1	2.95	0.44
1:A:215:VAL:HG21	1:A:226:LEU:HD21	1.99	0.44
1:A:322:VAL:HG23	1:A:367:LEU:HD21	1.99	0.44
1:A:420:THR:HG22	1:A:426:ALA:HB2	1.99	0.44
1:A:234:MET:HG3	1:A:255:PRO:HB2	2.00	0.44
1:B:234:MET:HE1	1:B:236:THR:OG1	2.18	0.44
1:B:232:ILE:O	1:B:254:ILE:HG21	2.18	0.43
1:D:163:ASN:HD22	1:D:163:ASN:HA	1.60	0.43
1:A:69:PRO:HG2	1:A:72:GLU:HG3	2.00	0.43
1:A:182:THR:CG2	1:A:211:ILE:HA	2.47	0.43
1:A:275:LYS:HA	1:A:275:LYS:HD3	1.88	0.43
1:D:116:ILE:HD11	1:D:449:ASP:CG	2.43	0.43
1:B:297:LYS:O	1:B:393:PRO:HD2	2.19	0.43
1:C:275:LYS:O	1:C:279:SER:HB2	2.19	0.43
1:B:74:VAL:HG21	1:B:125:LEU:HD13	1.99	0.43
1:B:294:THR:HG21	1:B:453:PHE:CE1	2.54	0.43
1:A:476:GLU:OE1	1:D:463:VAL:HG21	2.18	0.43
1:B:234:MET:HE3	1:B:235:ILE:C	2.44	0.43
1:A:297:LYS:NZ	1:A:386:GLU:OE2	2.48	0.43
1:B:218:ARG:HG3	1:B:221:VAL:CG2	2.48	0.43
1:B:377:THR:C	1:B:379:ALA:H	2.26	0.43
1:A:113:THR:HG23	1:A:170:ALA:HB3	2.01	0.43
1:B:74:VAL:HG13	1:B:121:ALA:HB1	2.00	0.43
1:B:420:THR:HG22	1:B:426:ALA:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:MET:CB	1:D:101:LYS:O	2.67	0.42
1:B:287:SER:HB2	1:B:292:ARG:NH1	2.33	0.42
1:A:238:THR:HG23	1:A:259:GLU:HG3	2.01	0.42
1:A:262:GLY:O	1:A:297:LYS:HD2	2.19	0.42
1:B:9:GLN:HB2	1:B:40:LEU:O	2.20	0.42
1:B:10:PHE:CD1	1:B:193:LEU:HD22	2.55	0.42
1:D:28:ARG:HD3	1:D:42:SER:OG	2.20	0.42
1:A:123:GLU:HG2	1:B:126:ARG:HE	1.85	0.42
1:C:96:MET:HE3	1:C:97:HIS:CE1	2.55	0.42
1:C:12:ALA:O	1:C:44:PRO:HD3	2.20	0.42
1:C:68:ARG:O	1:C:73:ARG:NH1	2.53	0.42
1:D:259:GLU:CD	1:D:464:GLN:OE1	2.62	0.42
1:A:348:ASP:HB2	4:A:523:HOH:O	2.19	0.42
1:D:397:VAL:HG12	1:D:398:LYS:N	2.34	0.42
1:B:55:ILE:HD13	1:B:229:HIS:CG	2.55	0.41
1:B:344:GLN:OE1	1:B:363:ARG:HD2	2.20	0.41
1:C:40:LEU:HD11	1:C:97:HIS:HB3	2.02	0.41
1:C:397:VAL:HG22	1:C:402:GLU:HB3	2.01	0.41
1:A:85:GLU:O	1:A:88:LYS:HG2	2.19	0.41
1:B:426:ALA:HB1	1:B:440:ILE:HD12	2.02	0.41
1:C:88:LYS:HA	1:C:110:VAL:HG11	2.02	0.41
1:D:127:LEU:HD23	1:D:127:LEU:HA	1.91	0.41
1:A:24:ARG:NH2	1:A:53:ASP:OD2	2.48	0.41
1:A:243:THR:HB	2:A:506:NAP:O2A	2.20	0.41
1:B:297:LYS:HD3	1:B:387:PRO:HD2	2.02	0.41
1:D:418:ILE:HD13	1:D:429:ILE:CG2	2.49	0.41
1:A:177:LEU:HD22	1:A:211:ILE:CG2	2.51	0.41
1:B:94:LEU:HD23	1:B:197:LYS:HG3	2.02	0.41
1:C:259:GLU:CG	1:C:259:GLU:O	2.68	0.41
1:C:446:ARG:HG3	1:C:453:PHE:CG	2.55	0.41
1:D:238:THR:HA	1:D:259:GLU:O	2.19	0.41
1:D:418:ILE:HD11	1:D:429:ILE:HG22	2.01	0.41
1:A:324:SER:HB3	1:A:327:ASP:OD2	2.20	0.41
1:D:20:TRP:CZ2	1:D:200:GLU:HA	2.55	0.41
1:D:259:GLU:HG2	1:D:464:GLN:NE2	2.36	0.41
1:B:392:LEU:HD12	1:B:393:PRO:HD2	2.02	0.41
1:C:124:ALA:C	1:C:126:ARG:H	2.29	0.41
1:C:159:ILE:HA	1:C:186:LYS:O	2.20	0.41
1:A:248:SER:OG	1:D:252:LYS:HB2	2.21	0.41
1:B:169:ALA:HB1	1:B:198:MET:HE1	2.02	0.41
1:B:265:PRO:HD2	1:B:415:GLN:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:309:LEU:O	1:B:313:ILE:HG13	2.21	0.41
1:D:236:THR:OG1	1:D:257:VAL:HG11	2.20	0.41
1:D:445:GLU:HB2	3:D:507:SO4:O3	2.21	0.41
1:A:96:MET:O	1:A:100:ALA:HA	2.21	0.41
1:B:218:ARG:H	1:B:218:ARG:HG2	1.34	0.41
1:C:273:ASP:OD1	1:C:273:ASP:C	2.63	0.41
1:D:15:LEU:HD21	1:D:212:ILE:HG23	2.02	0.41
1:D:199:VAL:O	1:D:200:GLU:C	2.64	0.41
1:A:417:SER:O	1:A:418:ILE:HD13	2.20	0.40
1:B:14:ILE:HG13	1:B:44:PRO:HG2	2.02	0.40
4:B:515:HOH:O	1:C:482:ASN:CB	2.68	0.40
1:A:25:THR:CG2	1:A:27:GLU:HG2	2.52	0.40
1:A:126:ARG:HD2	1:B:126:ARG:HD3	2.03	0.40
1:C:259:GLU:OE1	1:C:453:PHE:HE2	2.05	0.40
1:D:364:GLN:N	1:D:367:LEU:O	2.50	0.40
1:A:144:LYS:HB3	1:A:482:ASN:OD1	2.20	0.40
1:B:467:LYS:N	1:B:468:PRO:CD	2.85	0.40
1:C:94:LEU:HD23	1:C:194:SER:HA	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	473/505 (94%)	445 (94%)	25 (5%)	3 (1%)	21	27
1	B	474/505 (94%)	453 (96%)	20 (4%)	1 (0%)	43	55
1	C	476/505 (94%)	453 (95%)	22 (5%)	1 (0%)	43	55
1	D	479/505 (95%)	456 (95%)	22 (5%)	1 (0%)	43	55
All	All	1902/2020 (94%)	1807 (95%)	89 (5%)	6 (0%)	36	46

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	GLY
1	A	218	ARG
1	D	360	GLY
1	B	360	GLY
1	C	161	PRO
1	A	413	GLY

5.3.2 Protein sidechains ⓘ

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	386/414 (93%)	362 (94%)	24 (6%)	16	24
1	B	386/414 (93%)	363 (94%)	23 (6%)	17	25
1	C	389/414 (94%)	368 (95%)	21 (5%)	20	29
1	D	391/414 (94%)	354 (90%)	37 (10%)	8	10
All	All	1552/1656 (94%)	1447 (93%)	105 (7%)	14	20

All (105) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	THR
1	A	59	LYS
1	A	70	ILE
1	A	74	VAL
1	A	142	SER
1	A	150	ARG
1	A	209	GLU
1	A	234	MET
1	A	236	THR
1	A	252	LYS
1	A	256	VAL
1	A	259	GLU
1	A	270	ASP
1	A	279	SER
1	A	304	SER
1	A	310	VAL

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Mol	Chain	Res	Type
1	A	322	VAL
1	A	328	ASP
1	A	337	GLU
1	A	368	LEU
1	A	388	PHE
1	A	397	VAL
1	A	435	VAL
1	A	461	LEU
1	B	67	ILE
1	B	89	GLU
1	B	94	LEU
1	B	120	THR
1	B	133	LYS
1	B	145	ILE
1	B	150	ARG
1	B	167	ASN
1	B	178	VAL
1	B	213	GLN
1	B	218	ARG
1	B	256	VAL
1	B	259	GLU
1	B	269	LEU
1	B	294	THR
1	B	319	GLN
1	B	361	ASN
1	B	397	VAL
1	B	398	LYS
1	B	405	SER
1	B	431	LYS
1	B	463	VAL
1	B	471	LEU
1	C	8	GLU
1	C	25	THR
1	C	36	SER
1	C	49	GLU
1	C	89	GLU
1	C	105	SER
1	C	116	ILE
1	C	192	SER
1	C	249	GLU
1	C	259	GLU
1	C	283	SER

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Mol	Chain	Res	Type
1	C	287	SER
1	C	294	THR
1	C	310	VAL
1	C	315	GLU
1	C	328	ASP
1	C	388	PHE
1	C	391	VAL
1	C	398	LYS
1	C	463	VAL
1	C	471	LEU
1	D	25	THR
1	D	36	SER
1	D	49	GLU
1	D	62	GLN
1	D	74	VAL
1	D	89	GLU
1	D	93	GLU
1	D	94	LEU
1	D	133	LYS
1	D	142	SER
1	D	143	LYS
1	D	147	LEU
1	D	150	ARG
1	D	163	ASN
1	D	194	SER
1	D	218	ARG
1	D	245	GLU
1	D	257	VAL
1	D	259	GLU
1	D	269	LEU
1	D	296	ILE
1	D	315	GLU
1	D	328	ASP
1	D	332	THR
1	D	337	GLU
1	D	338	LYS
1	D	364	GLN
1	D	420	THR
1	D	423	THR
1	D	429	ILE
1	D	431	LYS
1	D	463	VAL

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Mol	Chain	Res	Type
1	D	467	LYS
1	D	471	LEU
1	D	473	MET
1	D	481	LEU
1	D	487	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	308	GLN
1	A	401	ASN
1	A	464	GLN
1	B	13	ASN
1	B	119	HIS
1	B	302	GLN
1	B	344	GLN
1	B	361	ASN
1	B	401	ASN
1	C	11	ASN
1	C	48	GLN
1	C	97	HIS
1	C	163	ASN
1	C	312	ASN
1	D	11	ASN
1	D	48	GLN
1	D	62	GLN
1	D	163	ASN
1	D	229	HIS
1	D	302	GLN
1	D	308	GLN
1	D	344	GLN
1	D	366	ASN
1	D	401	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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$
3	SO4	B	507	-	4,4,4	0.24	0	6,6,6	0.13	0
2	NAP	C	506	-	32,33,52	1.36	4 (12%)	50,52,80	1.78	9 (18%)
3	SO4	D	507	-	4,4,4	0.27	0	6,6,6	0.27	0
2	NAP	D	506	-	32,33,52	1.35	4 (12%)	50,52,80	1.74	9 (18%)
3	SO4	C	507	-	4,4,4	0.24	0	6,6,6	0.17	0
2	NAP	B	506	-	32,33,52	1.40	5 (15%)	50,52,80	1.75	8 (16%)
2	NAP	A	506	-	32,33,52	1.33	3 (9%)	50,52,80	1.73	9 (18%)
3	SO4	A	507	-	4,4,4	0.21	0	6,6,6	0.13	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	NAP	B	506	-	-	4/21/37/67	0/3/3/5
2	NAP	D	506	-	-	10/21/37/67	0/3/3/5
2	NAP	C	506	-	-	4/21/37/67	0/3/3/5
2	NAP	A	506	-	-	7/21/37/67	0/3/3/5

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	506	NAP	PN-O1N	3.82	1.62	1.50
2	B	506	NAP	PN-O1N	3.79	1.62	1.50
2	A	506	NAP	PN-O1N	3.77	1.62	1.50
2	C	506	NAP	PN-O1N	3.65	1.61	1.50
2	B	506	NAP	PA-O3	2.82	1.62	1.59
2	C	506	NAP	C2A-N1A	2.68	1.38	1.33
2	A	506	NAP	C2A-N3A	2.68	1.38	1.33
2	A	506	NAP	C2A-N1A	2.67	1.38	1.33
2	B	506	NAP	C2A-N1A	2.65	1.38	1.33
2	C	506	NAP	C2A-N3A	2.63	1.38	1.33
2	B	506	NAP	C2A-N3A	2.63	1.38	1.33
2	D	506	NAP	C2A-N1A	2.53	1.38	1.33
2	D	506	NAP	C2A-N3A	2.45	1.38	1.33
2	D	506	NAP	C8A-N7A	2.29	1.36	1.31
2	C	506	NAP	C8A-N7A	2.18	1.35	1.31
2	B	506	NAP	C8A-N7A	2.10	1.35	1.31

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	506	NAP	N3A-C2A-N1A	-5.76	119.86	128.58
2	D	506	NAP	N3A-C2A-N1A	-5.62	120.08	128.58
2	A	506	NAP	N3A-C2A-N1A	-5.43	120.36	128.58
2	B	506	NAP	N3A-C2A-N1A	-5.43	120.36	128.58
2	B	506	NAP	C5A-C4A-N3A	-4.41	120.65	126.72
2	C	506	NAP	C5A-C4A-N3A	-4.36	120.72	126.72
2	D	506	NAP	C5A-C4A-N3A	-4.12	121.04	126.72
2	A	506	NAP	C5A-C4A-N3A	-4.07	121.11	126.72
2	D	506	NAP	N9A-C8A-N7A	-4.03	108.22	113.94
2	C	506	NAP	N9A-C8A-N7A	-3.97	108.31	113.94
2	A	506	NAP	N9A-C8A-N7A	-3.83	108.50	113.94
2	B	506	NAP	N9A-C8A-N7A	-3.79	108.56	113.94
2	A	506	NAP	O5D-PN-O2N	3.72	121.75	107.80
2	C	506	NAP	C5A-N7A-C8A	3.57	109.06	103.45
2	C	506	NAP	C2A-N3A-C4A	3.47	120.31	111.83
2	D	506	NAP	C5A-N7A-C8A	3.42	108.82	103.45
2	B	506	NAP	C5A-N7A-C8A	3.35	108.72	103.45
2	B	506	NAP	C2A-N3A-C4A	3.30	119.88	111.83
2	D	506	NAP	C2A-N3A-C4A	3.26	119.80	111.83
2	A	506	NAP	C5A-N7A-C8A	3.17	108.44	103.45
2	A	506	NAP	C2A-N3A-C4A	3.13	119.48	111.83
2	C	506	NAP	O5D-PN-O2N	3.09	119.38	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	506	NAP	O5D-PN-O2N	2.99	119.01	107.80
2	B	506	NAP	N3A-C4A-N9A	2.93	132.16	127.17
2	A	506	NAP	N3A-C4A-N9A	2.89	132.07	127.17
2	B	506	NAP	O5D-PN-O2N	2.88	118.62	107.80
2	C	506	NAP	N3A-C4A-N9A	2.65	131.67	127.17
2	D	506	NAP	C4A-C5A-N7A	-2.63	107.58	110.58
2	C	506	NAP	C4A-C5A-N7A	-2.47	107.75	110.58
2	D	506	NAP	N3A-C4A-N9A	2.45	131.33	127.17
2	B	506	NAP	C4A-C5A-N7A	-2.30	107.96	110.58
2	C	506	NAP	O4B-C1B-C2B	-2.27	102.68	106.59
2	A	506	NAP	C4A-N9A-C8A	2.19	108.04	105.74
2	D	506	NAP	O3B-C3B-C4B	-2.11	105.02	111.08
2	A	506	NAP	C4A-C5A-N7A	-2.08	108.20	110.58

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	506	NAP	C5B-O5B-PA-O1A
2	D	506	NAP	C5B-O5B-PA-O2A
2	D	506	NAP	C5B-O5B-PA-O3
2	B	506	NAP	C2B-C1B-N9A-C8A
2	C	506	NAP	C2B-C1B-N9A-C8A
2	D	506	NAP	PN-O3-PA-O1A
2	A	506	NAP	PN-O3-PA-O5B
2	B	506	NAP	PN-O3-PA-O5B
2	D	506	NAP	PN-O3-PA-O5B
2	A	506	NAP	C2B-C1B-N9A-C8A
2	C	506	NAP	C2B-C1B-N9A-C4A
2	D	506	NAP	C2B-C1B-N9A-C8A
2	A	506	NAP	C5B-O5B-PA-O1A
2	A	506	NAP	C2B-C1B-N9A-C4A
2	B	506	NAP	C2B-C1B-N9A-C4A
2	D	506	NAP	C2B-O2B-P2B-O3X
2	D	506	NAP	C2B-C1B-N9A-C4A
2	C	506	NAP	C2B-O2B-P2B-O1X
2	A	506	NAP	O4B-C1B-N9A-C8A
2	D	506	NAP	PA-O3-PN-O1N
2	B	506	NAP	O4B-C1B-N9A-C8A
2	C	506	NAP	O4B-C1B-N9A-C8A
2	D	506	NAP	O4B-C1B-N9A-C8A
2	A	506	NAP	PN-O3-PA-O1A

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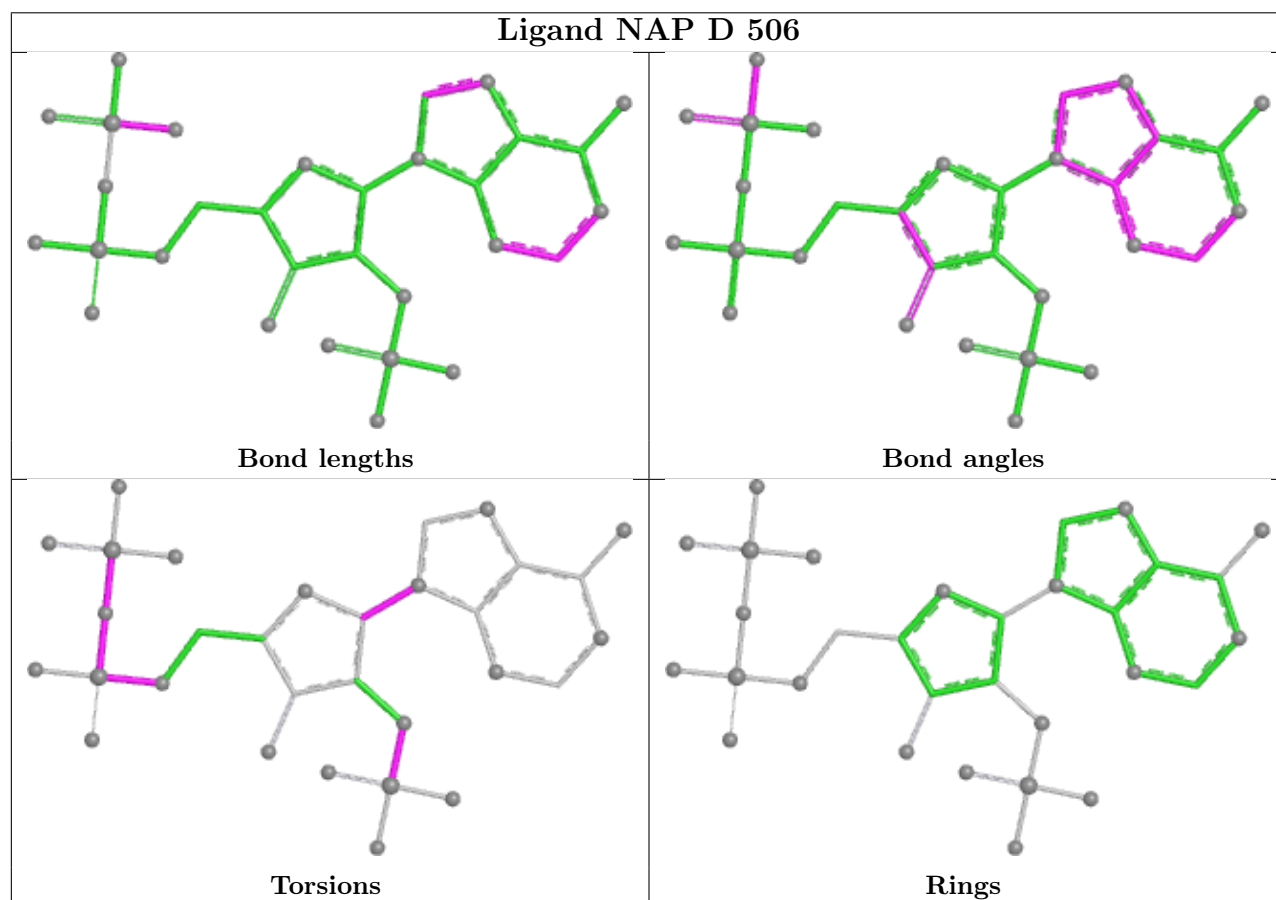
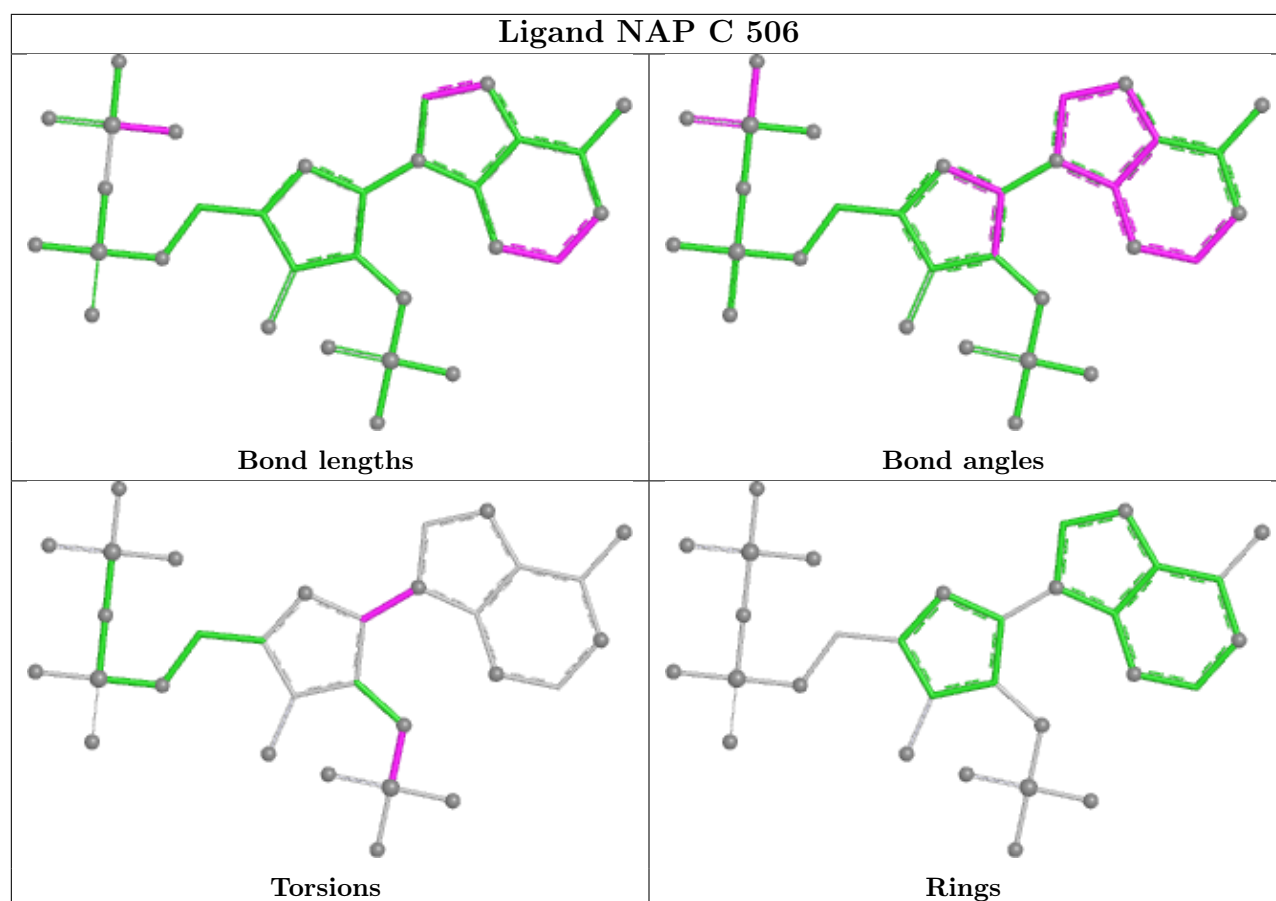
Mol	Chain	Res	Type	Atoms
2	A	506	NAP	PN-O3-PA-O2A

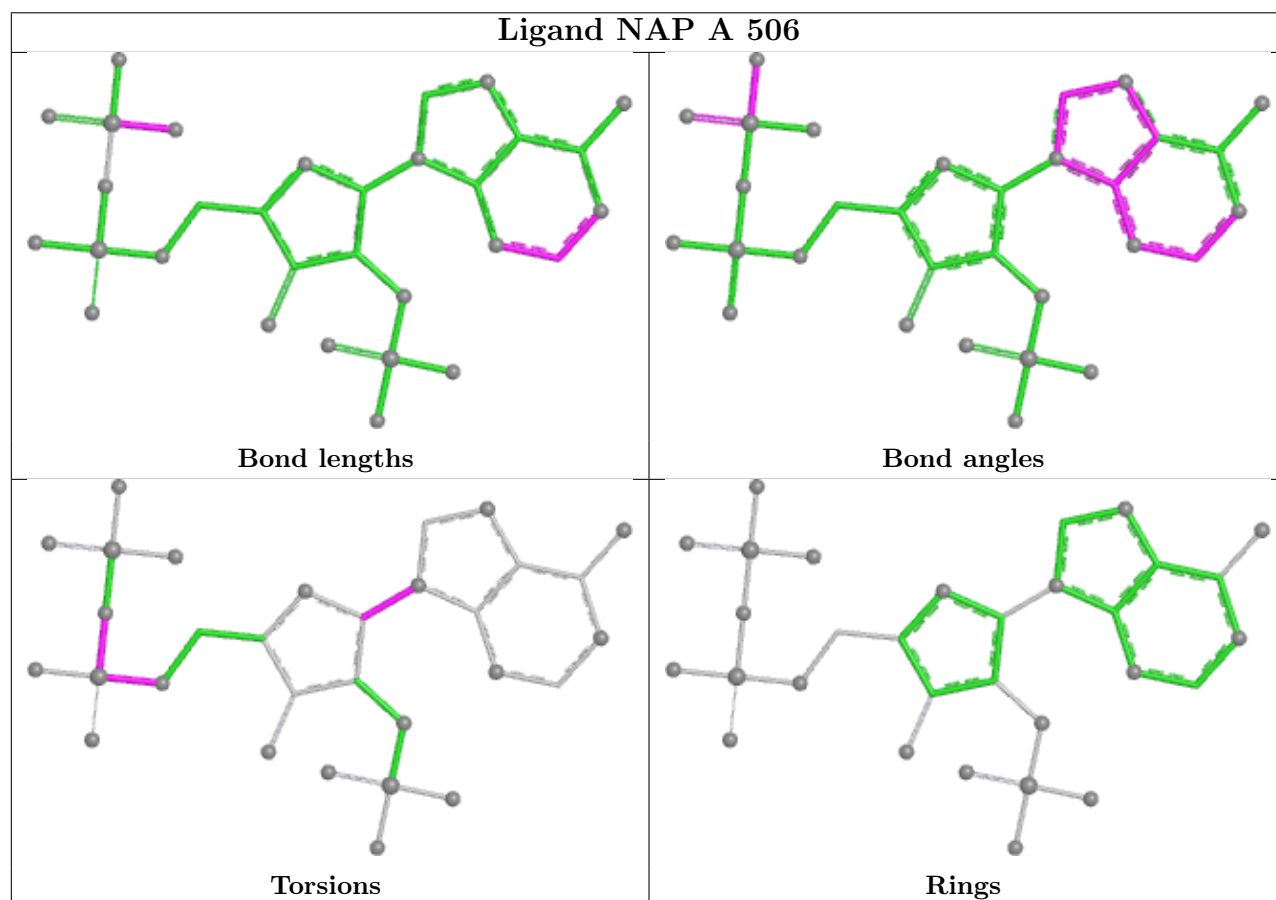
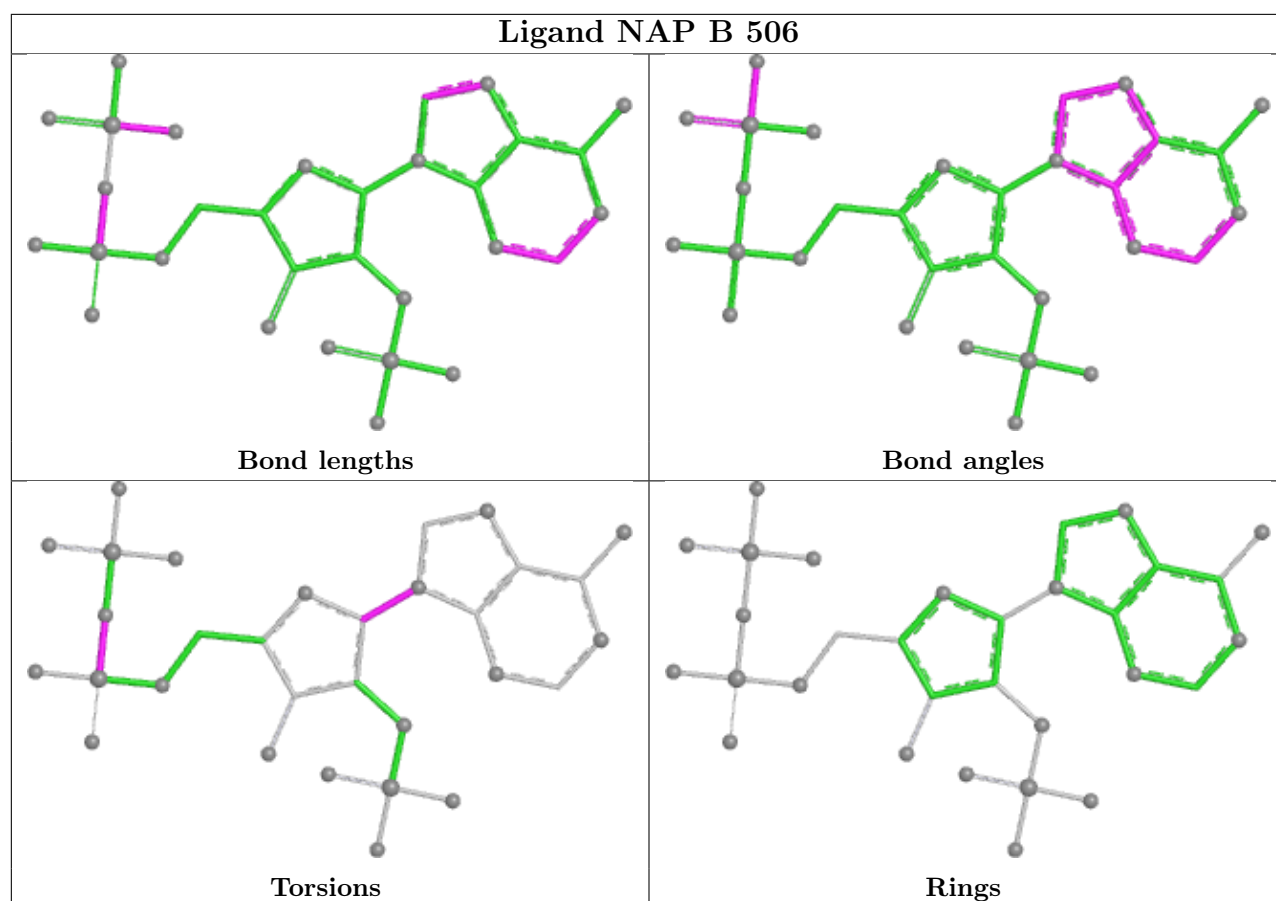
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	507	SO4	1	0
3	D	507	SO4	1	0
2	A	506	NAP	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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	475/505 (94%)	0.04	1 (0%) 91 91	33, 47, 64, 78	0
1	B	476/505 (94%)	-0.05	0 100 100	22, 45, 60, 68	0
1	C	476/505 (94%)	0.13	0 100 100	24, 52, 74, 89	2 (0%)
1	D	480/505 (95%)	0.21	9 (1%) 66 68	26, 56, 74, 90	1 (0%)
All	All	1907/2020 (94%)	0.08	10 (0%) 87 87	22, 49, 70, 90	3 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	483	LEU	3.3
1	D	453	PHE	2.8
1	D	220	SER	2.6
1	D	487	LEU	2.5
1	D	334	VAL	2.4
1	D	61	ALA	2.4
1	D	333	PRO	2.3
1	D	488	TYR	2.2
1	D	239	GLY	2.1
1	D	25	THR	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 oligosaccharides in this entry.

6.4 Ligands

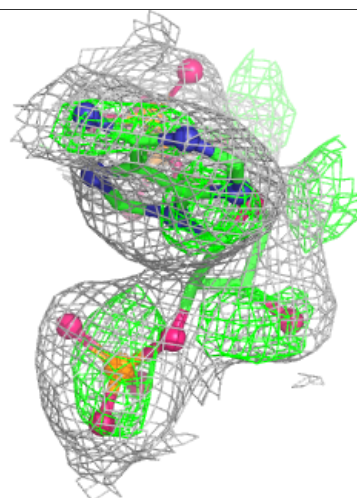
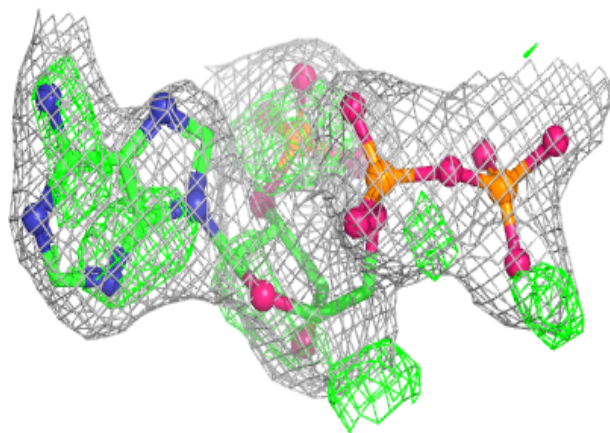
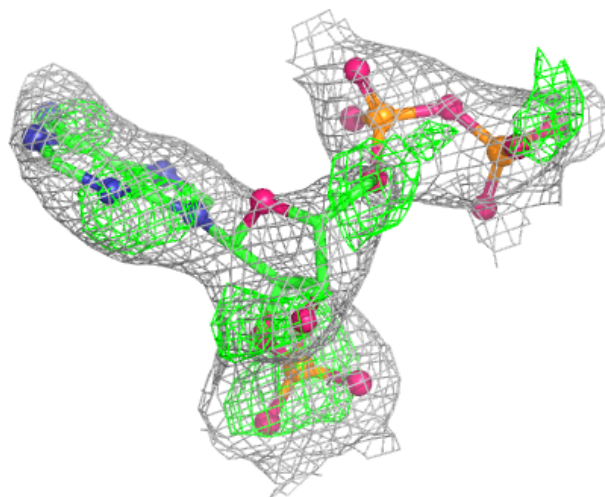
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	D	507	5/5	0.75	0.35	27,27,28,41	5
2	NAP	A	506	31/48	0.86	0.16	41,42,44,48	31
2	NAP	C	506	31/48	0.87	0.15	42,43,47,48	31
2	NAP	D	506	31/48	0.88	0.19	36,37,39,40	31
2	NAP	B	506	31/48	0.89	0.14	41,41,42,44	31
3	SO4	B	507	5/5	0.90	0.10	42,42,42,43	5
3	SO4	C	507	5/5	0.92	0.11	30,36,37,38	5
3	SO4	A	507	5/5	0.94	0.15	37,37,37,44	5

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.

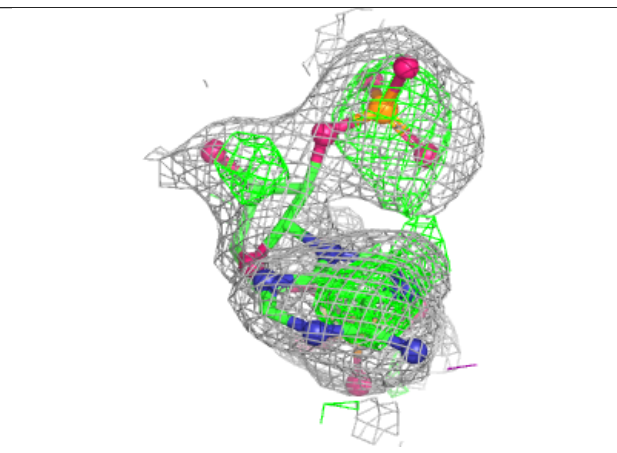
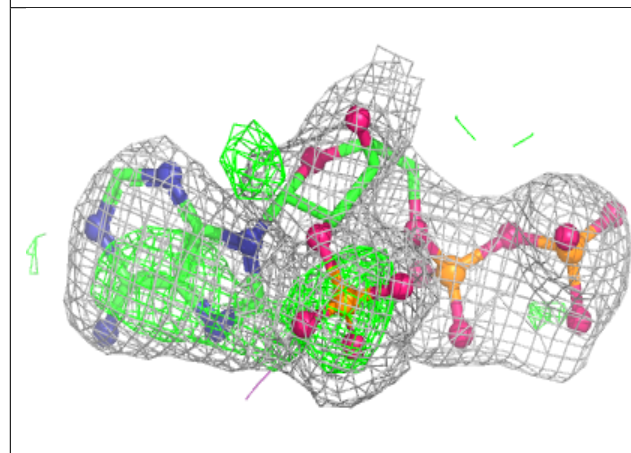
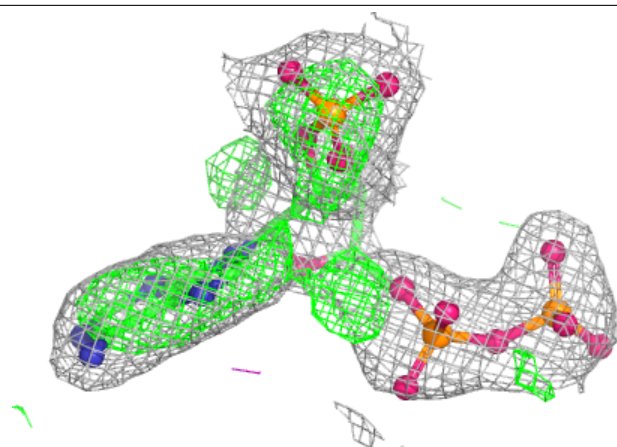
Electron density around NAP A 506:

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



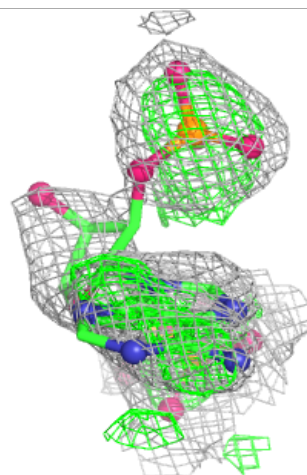
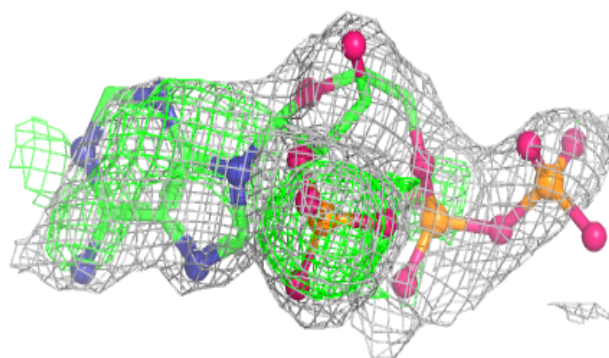
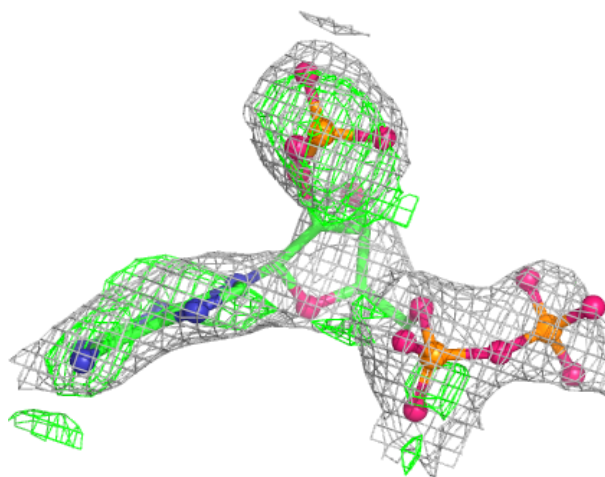
Electron density around NAP C 506:

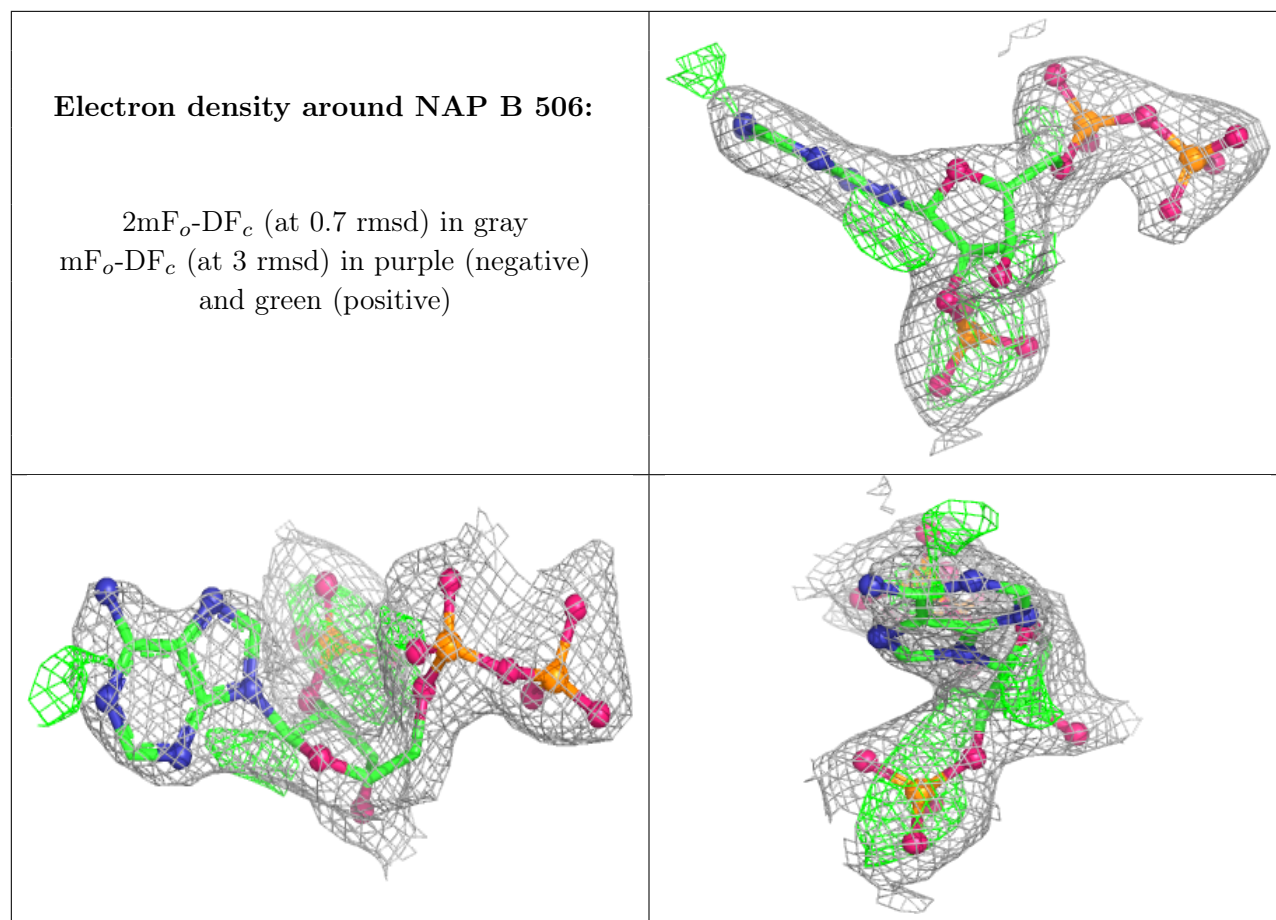
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around NAP D 506:

$2mF_o-DF_c$ (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.