



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:54 PM UTC

PDB ID : 2Y53 / pdb_00002y53
Title : Crystal structure of E257Q mutant of the box pathway encoded ALDH from Burkholderia xenovorans LB400
Authors : Bains, J.; Leon, R.; Temke, K.G.; Boulanger, M.J.
Deposited on : 2011-01-11
Resolution : 1.40 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)
Xtrriage (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

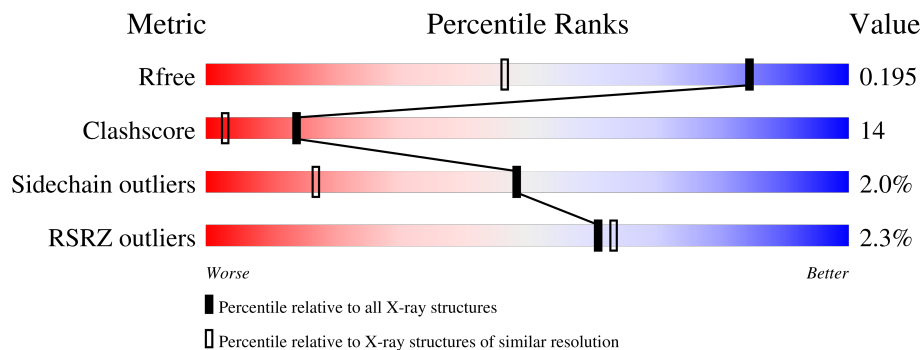
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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	2563 (1.40-1.40)
Clashscore	190562	2660 (1.40-1.40)
Sidechain outliers	187428	2610 (1.40-1.40)
RSRZ outliers	180081	2561 (1.40-1.40)

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	534	
1	B	534	

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
3	NAP	A	1534	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAP	B	1527	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9444 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 ALDEHYDE DEHYDROGENASE (BOX PATHWAY).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	529	3905	2451	706	738	10	0	1	0
1	B	524	3884	2434	703	737	10	0	3	0

There are 8 discrepancies between the modelled and reference sequences:

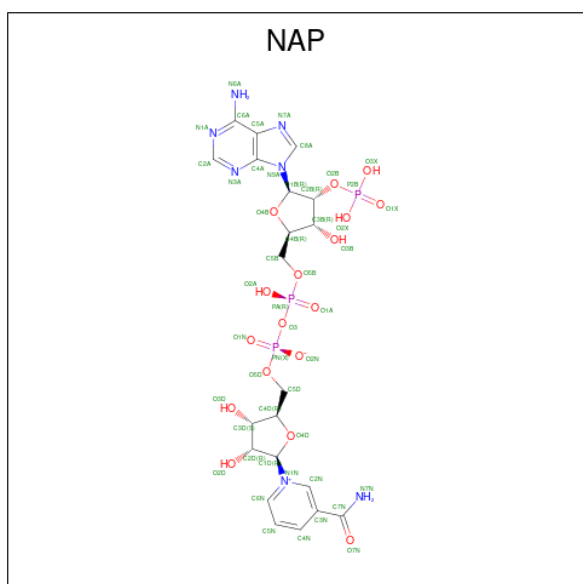
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q13WK4
A	-1	SER	-	expression tag	UNP Q13WK4
A	0	HIS	-	expression tag	UNP Q13WK4
A	257	GLN	GLU	engineered mutation	UNP Q13WK4
B	-2	GLY	-	expression tag	UNP Q13WK4
B	-1	SER	-	expression tag	UNP Q13WK4
B	0	HIS	-	expression tag	UNP Q13WK4
B	257	GLN	GLU	engineered mutation	UNP Q13WK4

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

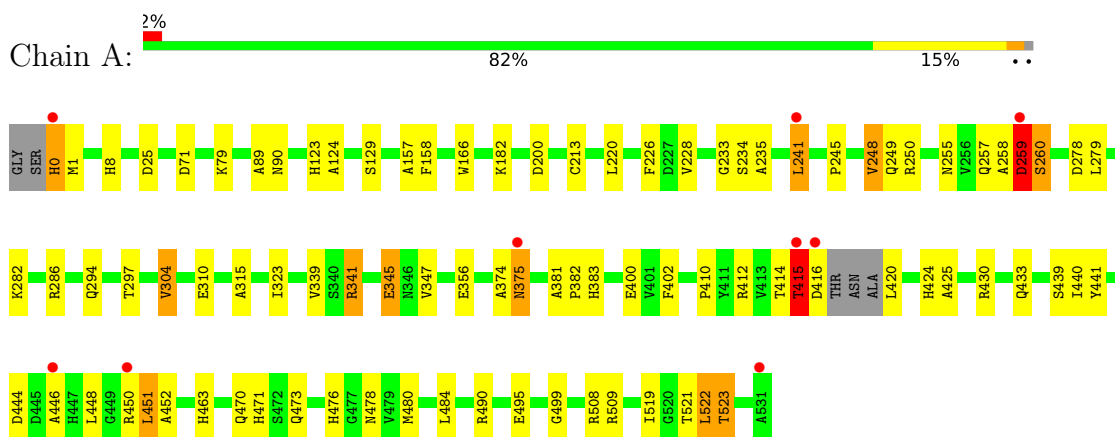
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	797	Total	O	0	0
			797	797		
4	B	738	Total	O	0	0
			738	738		

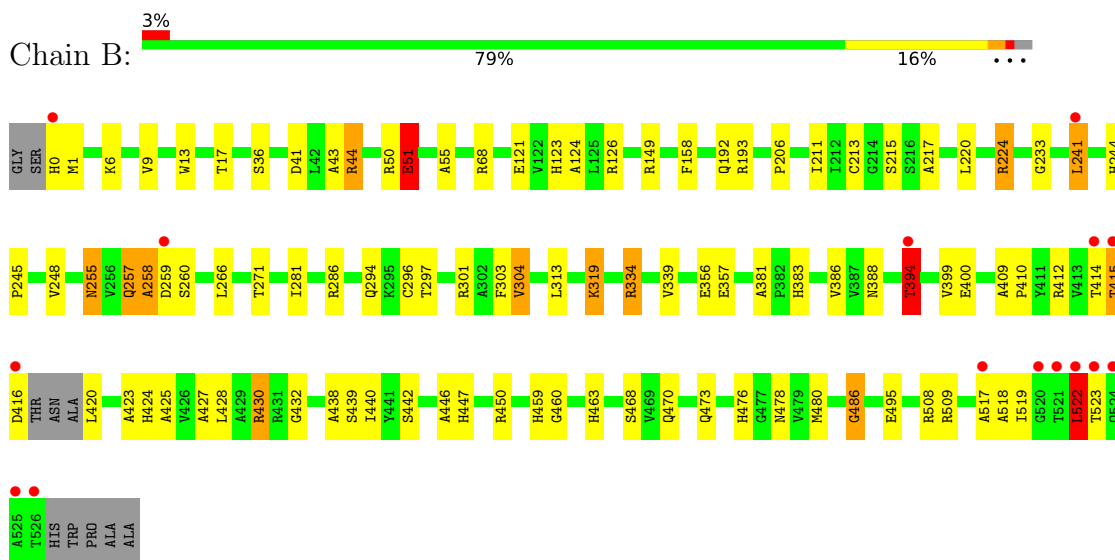
3 Residue-property plots

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: ALDEHYDE DEHYDROGENASE (BOX PATHWAY)



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.24Å 67.81Å 77.51Å 111.22° 90.45° 113.51°	Depositor
Resolution (Å)	23.37 – 1.40 23.37 – 1.40	Depositor EDS
% Data completeness (in resolution range)	95.6 (23.37-1.40) 95.6 (23.37-1.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 1.40Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.156 , 0.185 (Not available) , 0.195	Depositor DCC
R_{free} test set	9501 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	8.9	Xtrriage
Anisotropy	0.011	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.014 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9444	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.93% 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: NAP, GOL

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	1.74	44/3982 (1.1%)	1.48	29/5425 (0.5%)
1	B	1.75	39/3957 (1.0%)	1.43	14/5387 (0.3%)
All	All	1.75	83/7939 (1.0%)	1.45	43/10812 (0.4%)

All (83) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	341	ARG	CZ-NH2	11.52	1.48	1.33
1	A	129	SER	CB-OG	-11.03	1.20	1.42
1	A	412	ARG	CZ-NH2	-10.86	1.19	1.33
1	A	446	ALA	CA-CB	-9.92	1.37	1.53
1	B	394	THR	C-O	8.37	1.33	1.24
1	B	304	VAL	CA-C	8.17	1.61	1.52
1	A	375	ASN	CB-CG	7.99	1.72	1.52
1	A	425	ALA	CA-CB	-7.98	1.40	1.53
1	A	341	ARG	CD-NE	7.82	1.57	1.46
1	A	341	ARG	CZ-NH1	7.73	1.43	1.32
1	B	304	VAL	C-O	-7.67	1.18	1.24
1	A	451	LEU	CA-C	7.65	1.62	1.52
1	A	446	ALA	C-O	7.57	1.32	1.24
1	B	121	GLU	N-CA	-7.40	1.37	1.46
1	A	341	ARG	CB-CG	7.27	1.74	1.52
1	A	452	ALA	CA-CB	-6.96	1.42	1.53
1	A	508	ARG	CD-NE	-6.93	1.36	1.46
1	B	36	SER	C-O	-6.88	1.14	1.23
1	A	508	ARG	CZ-NH2	-6.73	1.24	1.33
1	B	424	HIS	CA-C	6.64	1.61	1.52
1	B	518	ALA	CA-CB	-6.60	1.43	1.53
1	B	486	GLY	N-CA	6.51	1.52	1.45
1	B	425	ALA	CA-CB	-6.49	1.43	1.53
1	B	259	ASP	C-O	-6.49	1.15	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	271	THR	CA-CB	6.46	1.61	1.53
1	B	13	TRP	C-O	6.42	1.32	1.24
1	A	424	HIS	CA-C	6.38	1.60	1.52
1	B	399	VAL	C-O	-6.33	1.17	1.24
1	B	412	ARG	N-CA	6.30	1.54	1.46
1	B	257	GLN	CA-C	-6.24	1.45	1.52
1	B	425	ALA	N-CA	6.19	1.53	1.46
1	B	440	ILE	N-CA	6.18	1.53	1.46
1	A	441	TYR	N-CA	6.06	1.53	1.46
1	A	382	PRO	C-O	6.00	1.30	1.23
1	B	206	PRO	N-CA	-5.96	1.40	1.47
1	A	304	VAL	CA-C	5.94	1.58	1.52
1	B	68	ARG	CD-NE	5.92	1.54	1.46
1	A	415	THR	N-CA	5.91	1.53	1.46
1	B	9	VAL	C-O	5.90	1.29	1.24
1	A	166	TRP	CA-C	-5.88	1.44	1.52
1	B	427	ALA	N-CA	-5.87	1.39	1.46
1	A	450	ARG	CA-C	5.86	1.60	1.52
1	B	432	GLY	C-O	-5.85	1.17	1.23
1	A	124	ALA	CA-C	-5.79	1.45	1.52
1	A	226	PHE	CD1-CE1	5.76	1.55	1.38
1	A	259	ASP	C-O	-5.74	1.16	1.23
1	A	444	ASP	N-CA	5.70	1.54	1.46
1	A	451	LEU	N-CA	-5.68	1.39	1.46
1	A	250	ARG	CZ-NH2	5.68	1.40	1.33
1	B	51	GLU	CB-CG	-5.67	1.35	1.52
1	B	193	ARG	CZ-NH2	5.65	1.40	1.33
1	A	260	SER	N-CA	-5.60	1.39	1.46
1	A	374	ALA	CA-C	-5.55	1.44	1.52
1	B	124	ALA	C-O	5.50	1.30	1.23
1	A	248	VAL	CB-CG2	-5.50	1.34	1.52
1	A	345	GLU	CD-OE1	-5.47	1.15	1.25
1	A	228	VAL	C-O	5.45	1.29	1.24
1	A	523	THR	N-CA	-5.43	1.39	1.46
1	B	427	ALA	CA-CB	5.39	1.61	1.53
1	B	258	ALA	CA-CB	-5.35	1.44	1.53
1	B	17	THR	CA-CB	5.34	1.62	1.53
1	B	43	ALA	CA-CB	5.34	1.61	1.53
1	B	430	ARG	CB-CG	-5.34	1.36	1.52
1	B	447	HIS	C-O	-5.32	1.17	1.24
1	A	213	CYS	C-O	5.31	1.30	1.24
1	A	323	ILE	N-CA	5.27	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	301	ARG	CA-C	-5.27	1.46	1.52
1	B	255	ASN	CA-C	-5.23	1.46	1.52
1	B	266	LEU	CA-C	5.22	1.59	1.52
1	A	260	SER	CB-OG	-5.16	1.31	1.42
1	A	499	GLY	N-CA	5.13	1.52	1.45
1	A	129	SER	C-O	5.12	1.31	1.23
1	A	90	ASN	CG-ND2	-5.12	1.22	1.33
1	A	424	HIS	C-O	-5.12	1.18	1.24
1	A	347	VAL	CA-CB	5.12	1.60	1.54
1	A	315	ALA	N-CA	-5.06	1.40	1.46
1	B	517	ALA	N-CA	5.05	1.52	1.46
1	B	213	CYS	C-O	5.04	1.30	1.24
1	B	215	SER	C-O	-5.04	1.17	1.23
1	B	50	ARG	CZ-NH1	5.04	1.39	1.32
1	B	55	ALA	CA-C	5.02	1.59	1.52
1	A	412	ARG	CA-CB	-5.01	1.46	1.53
1	A	8	HIS	C-O	5.01	1.29	1.24

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	341	ARG	NE-CZ-NH1	-16.92	104.58	121.50
1	A	412	ARG	NE-CZ-NH2	-16.05	104.76	119.20
1	A	412	ARG	NE-CZ-NH1	13.80	135.30	121.50
1	A	259	ASP	CA-CB-CG	-9.99	102.61	112.60
1	B	304	VAL	O-C-N	8.75	127.20	121.69
1	A	341	ARG	CB-CG-CD	8.27	130.32	111.30
1	A	259	ASP	N-CA-CB	8.07	121.99	110.44
1	A	341	ARG	NE-CZ-NH2	8.02	126.42	119.20
1	A	414	THR	CA-C-N	-7.51	112.41	123.00
1	A	414	THR	C-N-CA	-7.51	112.41	123.00
1	B	0	HIS	N-CA-C	-7.50	90.00	111.00
1	A	260	SER	N-CA-CB	-7.48	97.93	110.57
1	A	341	ARG	NH1-CZ-NH2	7.43	128.96	119.30
1	A	412	ARG	CD-NE-CZ	7.16	134.43	124.40
1	A	450	ARG	N-CA-C	7.13	118.84	111.14
1	A	375	ASN	CB-CA-C	7.06	121.89	110.09
1	B	0	HIS	CA-C-N	7.01	134.31	121.70
1	B	0	HIS	C-N-CA	7.01	134.31	121.70
1	B	442	SER	CA-C-O	6.85	128.78	121.45
1	A	345	GLU	CG-CD-OE1	6.84	134.13	118.40
1	A	375	ASN	N-CA-CB	-6.72	100.54	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	428	LEU	O-C-N	6.61	129.12	122.12
1	A	241	LEU	CA-CB-CG	6.43	138.81	116.30
1	B	468	SER	N-CA-C	6.18	119.92	112.38
1	B	303	PHE	CA-C-O	6.15	127.08	120.38
1	A	450	ARG	CA-C-O	6.14	127.02	120.70
1	B	260	SER	N-CA-CB	-5.87	100.96	110.71
1	B	266	LEU	O-C-N	5.76	129.75	123.13
1	A	345	GLU	CG-CD-OE2	-5.67	105.35	118.40
1	B	423	ALA	N-CA-C	5.47	117.67	111.11
1	A	129	SER	CA-CB-OG	-5.46	100.18	111.10
1	B	522	LEU	CB-CA-C	-5.39	102.18	110.81
1	A	304	VAL	O-C-N	5.37	125.07	121.69
1	A	79	LYS	CA-CB-CG	-5.35	103.40	114.10
1	A	415	THR	N-CA-CB	5.27	118.88	110.65
1	A	425	ALA	CA-C-O	-5.23	115.00	120.55
1	B	409	ALA	CA-C-O	5.21	125.66	120.03
1	A	279	LEU	N-CA-C	-5.16	105.74	111.36
1	A	259	ASP	N-CA-C	5.15	117.18	110.53
1	B	388	ASN	N-CA-C	5.09	117.73	111.82
1	A	424	HIS	O-C-N	5.07	127.29	122.07
1	A	375	ASN	CA-C-O	5.05	125.26	119.35
1	A	508	ARG	NE-CZ-NH2	-5.03	114.67	119.20

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	3905	0	3908	115	0
1	B	3884	0	3892	97	0
2	A	12	0	16	0	0
2	B	12	0	16	0	0
3	A	48	0	24	23	0
3	B	48	0	24	22	0
4	A	797	0	0	54	3

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	738	0	0	41	4
All	All	9444	0	7880	225	4

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 14.

All (225) 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:508[A]:ARG:HG2	4:B:2709:HOH:O	1.26	1.32
1:A:402:PHE:CZ	3:A:1534:NAP:O2D	1.84	1.30
1:B:286:ARG:HD3	4:B:2462:HOH:O	1.29	1.30
1:A:416:ASP:HB3	4:A:2686:HOH:O	1.16	1.27
1:A:249:GLN:HG2	4:A:2480:HOH:O	1.28	1.26
1:A:310:GLU:HG3	4:A:2525:HOH:O	1.30	1.22
3:B:1527:NAP:H6N	4:B:2724:HOH:O	1.36	1.22
1:A:278:ASP:HB3	4:A:2503:HOH:O	1.37	1.22
1:A:241:LEU:HD12	4:A:2470:HOH:O	1.35	1.20
1:A:0:HIS:CE1	4:A:2004:HOH:O	1.99	1.15
1:A:259:ASP:HB2	3:A:1534:NAP:C3N	1.78	1.13
1:B:523:THR:HG21	4:B:2324:HOH:O	1.49	1.10
1:B:286:ARG:NH2	1:B:478[A]:ASN:HD21	1.51	1.09
3:A:1534:NAP:C5D	4:A:2795:HOH:O	2.02	1.08
1:A:241:LEU:HB2	4:A:2470:HOH:O	1.52	1.08
1:A:286:ARG:NH2	1:A:478[A]:ASN:HD21	1.52	1.05
1:B:357[B]:GLU:OE2	1:B:394:THR:OG1	1.76	1.03
1:A:430:ARG:NH1	4:A:2695:HOH:O	1.93	1.01
1:A:248:VAL:HB	4:A:2476:HOH:O	0.81	0.99
1:B:296:CYS:SG	3:B:1527:NAP:O7N	2.22	0.98
1:A:420:LEU:N	4:A:2690:HOH:O	1.95	0.98
1:B:509:ARG:NH2	4:B:2716:HOH:O	1.96	0.98
1:A:478[B]:ASN:ND2	1:A:480:MET:CE	2.27	0.98
3:B:1527:NAP:O1N	3:B:1527:NAP:C5B	2.12	0.96
1:A:478[B]:ASN:ND2	1:A:480:MET:HE3	1.79	0.96
3:B:1527:NAP:O3B	4:B:2730:HOH:O	1.85	0.95
1:A:484:LEU:HD21	1:B:508[A]:ARG:HH22	1.34	0.93
4:A:2467:HOH:O	1:B:248:VAL:HB	0.74	0.92
1:A:519:ILE:O	1:A:523:THR:HG23	1.68	0.92
1:A:356:GLU:HG3	4:A:2588:HOH:O	1.71	0.90
1:A:0:HIS:ND1	1:A:0:HIS:O	2.03	0.90
1:B:519:ILE:O	1:B:523:THR:HG23	1.70	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:ASN:HB3	4:A:2635:HOH:O	1.72	0.89
3:A:1534:NAP:H52N	4:A:2795:HOH:O	1.66	0.89
1:A:235:ALA:H	3:A:1534:NAP:H71N	1.21	0.89
1:B:233:GLY:O	3:B:1527:NAP:N7N	2.08	0.86
1:A:484:LEU:HD21	1:B:508[A]:ARG:NH2	1.90	0.85
3:B:1527:NAP:O1N	3:B:1527:NAP:H51A	1.77	0.84
1:B:286:ARG:NH2	1:B:478[A]:ASN:ND2	2.27	0.83
1:B:420:LEU:N	4:B:2626:HOH:O	2.10	0.83
1:B:478[B]:ASN:ND2	1:B:480:MET:CE	2.41	0.83
1:A:375:ASN:CB	4:A:2635:HOH:O	2.28	0.82
1:A:495:GLU:OE2	1:B:508[A]:ARG:NH1	2.12	0.81
1:B:248:VAL:HG21	4:B:2427:HOH:O	1.80	0.81
1:A:286:ARG:NH2	1:A:478[A]:ASN:ND2	2.29	0.80
1:B:149:ARG:NH1	4:B:2333:HOH:O	1.76	0.80
1:A:158:PHE:CD2	3:A:1534:NAP:H51N	2.17	0.79
1:B:450:ARG:NE	4:B:2652:HOH:O	2.10	0.79
1:A:255:ASN:HD21	1:A:495:GLU:H	1.31	0.78
1:B:6:LYS:NZ	4:B:2024:HOH:O	2.15	0.78
1:B:478[B]:ASN:ND2	1:B:480:MET:HE1	1.99	0.78
1:A:157:ALA:HB1	3:A:1534:NAP:H3D	1.66	0.78
1:A:478[B]:ASN:HD21	1:A:480:MET:HE1	1.49	0.78
1:A:356:GLU:CG	4:A:2588:HOH:O	2.29	0.77
1:A:286:ARG:HH22	1:A:478[A]:ASN:HD21	1.33	0.76
1:A:0:HIS:HE1	4:A:2004:HOH:O	1.51	0.76
1:B:383:HIS:HE1	4:B:2555:HOH:O	1.69	0.74
1:B:508[A]:ARG:CG	4:B:2709:HOH:O	1.98	0.74
1:B:255:ASN:HD21	1:B:495:GLU:H	1.32	0.74
1:B:416:ASP:C	4:B:2621:HOH:O	2.29	0.74
1:A:341:ARG:HG2	4:A:2342:HOH:O	1.85	0.74
1:A:478[B]:ASN:ND2	1:A:480:MET:HE1	2.02	0.74
1:B:258:ALA:HA	3:B:1527:NAP:H72N	1.52	0.74
3:B:1527:NAP:O1N	3:B:1527:NAP:H52A	1.88	0.73
3:B:1527:NAP:O2A	3:B:1527:NAP:H52N	1.89	0.73
1:B:523:THR:CG2	4:B:2324:HOH:O	2.21	0.73
3:B:1527:NAP:C6N	4:B:2724:HOH:O	2.11	0.72
1:A:259:ASP:CB	3:A:1534:NAP:C3N	2.59	0.72
1:A:478[A]:ASN:ND2	4:A:2740:HOH:O	2.23	0.72
1:B:478[B]:ASN:ND2	1:B:480:MET:HE3	2.05	0.72
1:A:235:ALA:N	3:A:1534:NAP:H71N	1.87	0.71
1:A:356:GLU:OE2	4:A:2588:HOH:O	2.07	0.71
1:A:234:SER:HA	3:A:1534:NAP:H71N	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:ASN:CG	4:A:2635:HOH:O	2.32	0.70
1:A:282:LYS:HE3	4:A:2503:HOH:O	1.91	0.70
1:A:375:ASN:CB	4:A:2634:HOH:O	2.39	0.70
1:B:244:HIS:O	1:B:248:VAL:HG22	1.91	0.70
1:A:375:ASN:HB3	4:A:2634:HOH:O	1.92	0.70
1:B:478[A]:ASN:ND2	4:B:2681:HOH:O	2.24	0.70
1:A:259:ASP:HB2	3:A:1534:NAP:C2N	2.21	0.69
1:A:383:HIS:HE1	4:A:2609:HOH:O	1.75	0.69
1:B:508[A]:ARG:NH1	4:B:2711:HOH:O	2.24	0.69
1:A:286:ARG:HH22	1:A:478[A]:ASN:ND2	1.88	0.68
1:A:490:ARG:HD3	4:A:2749:HOH:O	1.92	0.68
1:B:245:PRO:HA	1:B:248:VAL:CG2	2.23	0.68
1:A:523:THR:HG21	4:A:2359:HOH:O	1.94	0.67
1:B:450:ARG:NH2	4:B:2653:HOH:O	2.26	0.67
1:A:158:PHE:CE2	3:A:1534:NAP:H4D	2.31	0.66
1:A:439:SER:OG	1:A:476:HIS:HD2	1.80	0.65
1:B:51:GLU:OE2	4:B:2135:HOH:O	2.13	0.65
1:A:402:PHE:CE1	3:A:1534:NAP:O2D	2.45	0.64
1:A:259:ASP:OD2	4:A:2491:HOH:O	2.14	0.64
1:A:341:ARG:CG	4:A:2571:HOH:O	2.45	0.64
1:B:258:ALA:HA	3:B:1527:NAP:N7N	2.14	0.63
1:A:286:ARG:HH21	1:A:478[A]:ASN:HD21	1.43	0.63
1:A:375:ASN:ND2	4:A:2638:HOH:O	2.21	0.63
3:B:1527:NAP:O2A	3:B:1527:NAP:C5D	2.47	0.62
1:A:402:PHE:CE2	3:A:1534:NAP:O2D	2.46	0.62
1:B:439:SER:OG	1:B:476:HIS:HD2	1.83	0.61
1:B:245:PRO:HA	1:B:248:VAL:HG23	1.82	0.61
1:A:415:THR:C	1:A:416:ASP:OD1	2.44	0.61
3:B:1527:NAP:H52N	4:B:2727:HOH:O	2.00	0.60
1:A:341:ARG:HG2	4:A:2571:HOH:O	2.02	0.59
1:A:0:HIS:O	1:A:0:HIS:CG	2.55	0.59
1:B:123:HIS:HE1	4:B:2273:HOH:O	1.85	0.59
1:A:123:HIS:HE1	4:A:2311:HOH:O	1.85	0.59
1:A:294:GLN:HE22	1:A:339:VAL:H	1.48	0.58
1:B:224:ARG:NH1	4:B:2401:HOH:O	2.36	0.58
1:A:440:ILE:HG13	1:A:448:LEU:HD22	1.84	0.58
1:B:220:LEU:HD11	1:B:241:LEU:HG	1.84	0.58
1:A:381:ALA:O	1:A:383:HIS:HD2	1.87	0.58
1:A:248:VAL:CG2	4:A:2476:HOH:O	2.13	0.57
1:A:356:GLU:CD	4:A:2588:HOH:O	2.47	0.57
1:A:470:GLN:HG2	4:A:2734:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:381:ALA:O	1:B:383:HIS:HD2	1.88	0.56
1:A:234:SER:CA	3:A:1534:NAP:H71N	2.18	0.56
1:A:0:HIS:CG	4:A:2003:HOH:O	2.58	0.56
1:A:71:ASP:OD1	4:A:2205:HOH:O	2.18	0.56
1:A:241:LEU:CB	4:A:2470:HOH:O	2.25	0.56
1:A:258:ALA:HA	3:A:1534:NAP:C5N	2.36	0.56
1:A:258:ALA:HA	3:A:1534:NAP:H5N	1.89	0.55
1:B:281:ILE:HG22	1:B:319:LYS:HE3	1.88	0.55
1:B:478[B]:ASN:HD21	1:B:480:MET:HE1	1.69	0.55
1:A:255:ASN:HD21	1:A:495:GLU:N	2.03	0.54
1:A:463:HIS:HE1	1:A:473:GLN:OE1	1.90	0.54
1:A:522:LEU:HD12	1:A:522:LEU:C	2.33	0.53
1:B:415:THR:HG23	4:B:2620:HOH:O	2.07	0.53
1:B:294:GLN:HE22	1:B:339:VAL:H	1.56	0.53
1:A:345:GLU:OE2	4:A:2577:HOH:O	0.60	0.53
1:B:313:LEU:HD21	1:B:386:VAL:HG22	1.90	0.53
1:A:200:ASP:OD1	4:A:2419:HOH:O	2.19	0.53
1:A:258:ALA:CA	3:A:1534:NAP:H5N	2.34	0.52
1:A:304:VAL:O	1:A:410:PRO:HA	2.09	0.52
1:B:463:HIS:HE1	1:B:473:GLN:OE1	1.94	0.51
1:A:0:HIS:CD2	4:A:2003:HOH:O	2.62	0.51
1:A:241:LEU:CD1	4:A:2470:HOH:O	2.05	0.51
1:B:400:GLU:HA	4:B:2603:HOH:O	2.11	0.51
1:A:158:PHE:CD1	1:A:339:VAL:HG21	2.47	0.50
1:B:415:THR:O	1:B:416:ASP:CB	2.58	0.50
1:B:508[A]:ARG:HG3	1:B:509:ARG:N	2.25	0.50
1:A:249:GLN:NE2	4:A:2478:HOH:O	2.44	0.50
1:A:375:ASN:HB3	4:A:2637:HOH:O	2.10	0.50
1:B:415:THR:O	1:B:416:ASP:HB2	2.12	0.50
3:A:1534:NAP:H51N	4:A:2795:HOH:O	1.90	0.50
1:A:182:LYS:NZ	3:A:1534:NAP:O1X	2.45	0.49
1:A:471:HIS:CD2	4:A:2733:HOH:O	2.66	0.49
3:A:1534:NAP:O5B	3:A:1534:NAP:O5D	2.30	0.49
1:B:217:ALA:HB2	4:B:2383:HOH:O	2.12	0.49
1:B:470:GLN:HG2	4:B:2665:HOH:O	2.13	0.49
1:B:224:ARG:CG	1:B:224:ARG:HH11	2.25	0.49
1:B:334:ARG:NH1	1:B:334:ARG:HG3	2.27	0.49
1:B:334:ARG:NH2	4:B:2506:HOH:O	2.46	0.49
1:A:415:THR:O	1:A:416:ASP:OD1	2.30	0.49
1:B:319:LYS:HB2	1:B:319:LYS:HE2	1.66	0.48
1:B:356:GLU:HA	1:B:356:GLU:OE1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:THR:HG21	1:B:446:ALA:HB1	1.94	0.48
1:A:341:ARG:HG3	4:A:2571:HOH:O	2.09	0.47
1:B:439:SER:OG	1:B:476:HIS:CD2	2.66	0.47
1:A:420:LEU:HD11	1:A:451:LEU:HA	1.96	0.47
1:A:310:GLU:CG	4:A:2525:HOH:O	2.13	0.47
1:B:286:ARG:CD	4:B:2462:HOH:O	2.14	0.47
1:B:233:GLY:HA2	3:B:1527:NAP:H2N	1.97	0.46
1:A:259:ASP:CG	1:A:260:SER:H	2.23	0.46
1:A:478[B]:ASN:HD22	1:A:480:MET:HE3	1.73	0.46
1:B:478[A]:ASN:CG	4:B:2681:HOH:O	2.56	0.46
1:B:509:ARG:CZ	4:B:2716:HOH:O	2.54	0.46
1:A:521:THR:HG21	1:B:446:ALA:CB	2.45	0.46
1:A:220:LEU:HD11	1:A:241:LEU:HG	1.97	0.46
1:B:297:THR:O	1:B:476:HIS:CE1	2.69	0.45
1:B:286:ARG:HH21	1:B:478[A]:ASN:HD21	1.55	0.45
1:B:297:THR:O	1:B:476:HIS:HE1	1.99	0.45
1:B:334:ARG:NH1	1:B:334:ARG:CG	2.79	0.45
3:B:1527:NAP:O2A	3:B:1527:NAP:O5D	2.35	0.45
1:A:233:GLY:O	1:A:258:ALA:HA	2.17	0.45
1:B:255:ASN:HD21	1:B:495:GLU:N	2.09	0.45
1:B:258:ALA:CA	3:B:1527:NAP:H72N	2.23	0.45
1:A:297:THR:O	1:A:476:HIS:CE1	2.70	0.45
1:A:433:GLN:NE2	4:A:2701:HOH:O	2.50	0.45
1:B:439:SER:CB	1:B:476:HIS:HD2	2.29	0.45
3:B:1527:NAP:O1X	4:B:2728:HOH:O	2.21	0.45
1:A:297:THR:O	1:A:476:HIS:HE1	1.99	0.45
1:A:439:SER:CB	1:A:476:HIS:HD2	2.30	0.44
1:A:158:PHE:CE2	3:A:1534:NAP:H51N	2.52	0.44
3:B:1527:NAP:O1N	3:B:1527:NAP:H3D	2.18	0.44
1:A:375:ASN:CG	4:A:2634:HOH:O	2.59	0.44
1:B:470:GLN:CD	4:B:2669:HOH:O	2.60	0.44
1:A:0:HIS:HA	1:A:1:MET:HA	1.60	0.44
1:A:286:ARG:NH1	4:A:2511:HOH:O	2.22	0.44
1:B:334:ARG:CG	1:B:334:ARG:HH11	2.30	0.44
1:B:192:GLN:HA	1:B:211:ILE:HD13	2.00	0.44
1:A:381:ALA:O	1:A:383:HIS:CD2	2.70	0.43
3:B:1527:NAP:H3D	4:B:2734:HOH:O	2.16	0.43
1:A:245:PRO:HA	1:A:248:VAL:HG22	2.00	0.43
1:A:478[A]:ASN:CG	4:A:2740:HOH:O	2.57	0.43
1:B:430:ARG:HG3	4:B:2662:HOH:O	2.17	0.43
1:B:123:HIS:CE1	4:B:2273:HOH:O	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:ARG:NH1	1:B:495:GLU:OE1	2.52	0.43
1:B:158:PHE:CD1	1:B:339:VAL:HG21	2.55	0.42
1:B:304:VAL:O	1:B:410:PRO:HA	2.19	0.42
1:B:1:MET:HE3	4:B:2038:HOH:O	2.19	0.42
1:A:439:SER:OG	1:A:476:HIS:CD2	2.66	0.42
1:A:234:SER:HA	3:A:1534:NAP:N7N	2.30	0.42
1:A:25:ASP:HA	1:A:89:ALA:O	2.20	0.42
1:B:126:ARG:HD3	1:B:523:THR:HG22	2.02	0.42
1:B:522:LEU:HD23	1:B:522:LEU:HA	1.82	0.41
1:B:258:ALA:C	3:B:1527:NAP:N7N	2.78	0.41
1:A:415:THR:CG2	4:A:2675:HOH:O	2.69	0.41
1:A:341:ARG:HH11	1:A:341:ARG:HD2	1.44	0.41
1:A:439:SER:HA	1:A:463:HIS:O	2.21	0.41
1:B:286:ARG:NH1	4:B:2460:HOH:O	2.49	0.41
1:B:313:LEU:HD21	1:B:386:VAL:CG2	2.49	0.41
1:B:438:ALA:HB2	1:B:459:HIS:CD2	2.56	0.41
1:B:414:THR:HG22	1:B:416:ASP:H	1.85	0.41
3:B:1527:NAP:C5B	3:B:1527:NAP:PN	3.09	0.40
3:B:1527:NAP:C5D	4:B:2727:HOH:O	2.64	0.40
1:B:245:PRO:HA	1:B:248:VAL:HG22	2.00	0.40
1:B:460:GLY:HA3	1:B:486:GLY:O	2.21	0.40
1:B:41:ASP:CG	1:B:44:ARG:HB2	2.46	0.40
1:B:233:GLY:O	1:B:258:ALA:HA	2.21	0.40
1:B:394:THR:OG1	4:B:2594:HOH:O	2.21	0.40
1:B:400:GLU:CD	4:B:2599:HOH:O	2.65	0.40
1:A:420:LEU:HD13	1:A:451:LEU:HD13	2.04	0.40
1:B:281:ILE:CG2	1:B:319:LYS:HE3	2.49	0.40
1:B:381:ALA:O	1:B:383:HIS:CD2	2.72	0.40

All (4) 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)
4:A:2238:HOH:O	4:B:2192:HOH:O[1_444]	2.05	0.15
4:A:2109:HOH:O	4:B:2003:HOH:O[1_444]	2.11	0.09
4:B:2431:HOH:O	4:B:2665:HOH:O[1_455]	2.17	0.03
4:A:2239:HOH:O	4:B:2189:HOH:O[1_444]	2.18	0.02

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	397/399 (100%)	391 (98%)	6 (2%)	57	26
1	B	396/399 (99%)	386 (98%)	10 (2%)	42	11
All	All	793/798 (99%)	777 (98%)	16 (2%)	48	17

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	257	GLN
1	A	259	ASP
1	A	400	GLU
1	A	415	THR
1	A	522	LEU
1	B	44	ARG
1	B	51	GLU
1	B	224	ARG
1	B	241	LEU
1	B	257	GLN
1	B	319	LYS
1	B	334	ARG
1	B	394	THR
1	B	415	THR
1	B	522	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	123	HIS
1	A	142	GLN
1	A	249	GLN
1	A	255	ASN
1	A	294	GLN
1	A	346	ASN
1	A	383	HIS
1	A	433	GLN
1	A	463	HIS
1	A	476	HIS
1	A	527	HIS
1	B	12	GLN
1	B	123	HIS
1	B	255	ASN
1	B	343	GLN
1	B	346	ASN
1	B	383	HIS
1	B	463	HIS
1	B	476	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	B	1528	-	5,5,5	0.81	0	5,5,5	1.03	1 (20%)
3	NAP	B	1527	-	50,52,52	3.27	19 (38%)	71,80,80	3.33	29 (40%)
3	NAP	A	1534	-	50,52,52	1.58	11 (22%)	71,80,80	1.78	14 (19%)
2	GOL	B	1529	-	5,5,5	0.92	0	5,5,5	0.65	0
2	GOL	A	1532	-	5,5,5	0.96	0	5,5,5	0.56	0
2	GOL	A	1533	-	5,5,5	0.79	0	5,5,5	0.87	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	GOL	B	1528	-	-	0/4/4/4	-
3	NAP	B	1527	-	-	14/35/67/67	0/5/5/5
3	NAP	A	1534	-	-	22/35/67/67	0/5/5/5
2	GOL	B	1529	-	-	0/4/4/4	-
2	GOL	A	1532	-	-	0/4/4/4	-
2	GOL	A	1533	-	-	0/4/4/4	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1527	NAP	PA-O3	-9.25	1.49	1.59
3	B	1527	NAP	O3B-C3B	-8.45	1.22	1.43
3	B	1527	NAP	O3D-C3D	6.98	1.60	1.43
3	B	1527	NAP	O4D-C4D	-5.64	1.32	1.45
3	B	1527	NAP	C5A-N7A	-5.38	1.29	1.39
3	B	1527	NAP	O4B-C1B	-5.34	1.29	1.42
3	B	1527	NAP	C2N-N1N	5.19	1.40	1.35
3	B	1527	NAP	O4D-C1D	5.11	1.47	1.40
3	A	1534	NAP	C2N-N1N	4.90	1.40	1.35
3	B	1527	NAP	C8A-N7A	4.45	1.40	1.31
3	B	1527	NAP	C4A-N3A	-4.32	1.26	1.34
3	B	1527	NAP	C5A-C4A	-4.29	1.31	1.39
3	B	1527	NAP	PN-O3	-4.07	1.55	1.59
3	B	1527	NAP	P2B-O2B	3.99	1.66	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1534	NAP	P2B-O1X	3.77	1.62	1.50
3	B	1527	NAP	C3D-C4D	3.61	1.62	1.53
3	A	1534	NAP	PA-O1A	3.34	1.62	1.50
3	A	1534	NAP	PN-O1N	3.32	1.62	1.50
3	B	1527	NAP	P2B-O1X	3.25	1.60	1.50
3	B	1527	NAP	PA-O1A	3.16	1.61	1.50
3	A	1534	NAP	C5A-N7A	-3.16	1.33	1.39
3	B	1527	NAP	C8A-N9A	-3.15	1.32	1.37
3	B	1527	NAP	O4B-C4B	-3.08	1.38	1.45
3	A	1534	NAP	PA-O3	2.66	1.62	1.59
3	B	1527	NAP	C1B-N9A	-2.46	1.39	1.46
3	A	1534	NAP	O4D-C1D	2.32	1.43	1.40
3	A	1534	NAP	C6N-N1N	2.21	1.40	1.35
3	A	1534	NAP	C8A-N9A	-2.19	1.33	1.37
3	A	1534	NAP	C4A-N9A	-2.09	1.33	1.37
3	A	1534	NAP	P2B-O2X	2.00	1.62	1.54

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1527	NAP	C4D-O4D-C1D	-15.86	95.40	109.92
3	B	1527	NAP	C5B-C4B-C3B	-7.98	86.50	115.21
3	B	1527	NAP	N9A-C8A-N7A	-7.89	102.74	113.94
3	B	1527	NAP	O4D-C4D-C5D	-6.71	87.85	109.33
3	A	1534	NAP	C4D-O4D-C1D	-6.40	104.06	109.92
3	B	1527	NAP	O4D-C4D-C3D	-5.56	94.12	105.15
3	A	1534	NAP	C5A-C4A-N3A	-5.32	119.39	126.72
3	B	1527	NAP	C4A-N9A-C8A	5.27	111.27	105.74
3	B	1527	NAP	N3A-C2A-N1A	-5.01	121.00	128.58
3	B	1527	NAP	O3D-C3D-C2D	4.87	127.42	111.82
3	B	1527	NAP	O2D-C2D-C3D	4.85	127.37	111.82
3	A	1534	NAP	N3A-C2A-N1A	-4.53	121.73	128.58
3	B	1527	NAP	C5D-C4D-C3D	4.47	131.31	115.21
3	B	1527	NAP	C5A-N7A-C8A	4.43	110.41	103.45
3	B	1527	NAP	O2N-PN-O5D	3.91	125.30	107.57
3	B	1527	NAP	C2A-N3A-C4A	3.76	121.01	111.83
3	A	1534	NAP	N3A-C4A-N9A	3.63	133.34	127.17
3	A	1534	NAP	C2A-N3A-C4A	3.52	120.42	111.83
3	B	1527	NAP	C5A-C4A-N3A	-3.44	121.97	126.72
3	B	1527	NAP	O4B-C1B-C2B	-3.41	100.72	106.59
3	B	1527	NAP	C6N-C5N-C4N	3.35	124.28	119.45
3	A	1534	NAP	N9A-C8A-N7A	-3.27	109.30	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1527	NAP	N3A-C4A-N9A	3.06	132.37	127.17
3	B	1527	NAP	O3D-C3D-C4D	2.99	119.67	111.08
3	B	1527	NAP	C6N-N1N-C2N	-2.98	119.34	121.88
3	B	1527	NAP	O2A-PA-O3	2.88	115.07	107.27
3	A	1534	NAP	C5A-N7A-C8A	2.84	107.91	103.45
3	B	1527	NAP	C5N-C6N-N1N	-2.82	116.54	120.38
3	A	1534	NAP	C4A-C5A-N7A	-2.70	107.50	110.58
3	B	1527	NAP	PA-O5B-C5B	2.59	136.19	121.35
3	B	1527	NAP	O3-PA-O1A	-2.59	102.92	110.70
3	A	1534	NAP	C2B-C1B-N9A	-2.59	109.50	113.75
3	A	1534	NAP	C4B-O4B-C1B	-2.44	104.08	109.47
3	A	1534	NAP	O2N-PN-O3	2.42	113.81	107.27
3	B	1527	NAP	C4A-N9A-C1B	-2.41	121.00	126.63
3	A	1534	NAP	C6N-N1N-C2N	-2.36	119.87	121.88
3	A	1534	NAP	C3N-C7N-N7N	-2.24	114.97	117.74
3	A	1534	NAP	C4A-N9A-C8A	2.19	108.04	105.74
3	B	1527	NAP	PN-O5D-C5D	2.15	133.69	121.35
3	B	1527	NAP	O2B-P2B-O1X	-2.13	101.73	109.33
3	B	1527	NAP	O2B-C2B-C3B	2.08	119.12	111.68
3	B	1527	NAP	C5A-C6A-N6A	-2.07	118.16	123.29
2	B	1528	GOL	O2-C2-C3	-2.02	100.82	109.18
3	B	1527	NAP	C2D-C3D-C4D	-2.01	98.73	102.61

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1534	NAP	C5D-O5D-PN-O3
3	A	1534	NAP	O4D-C4D-C5D-O5D
3	A	1534	NAP	O4D-C1D-N1N-C2N
3	A	1534	NAP	O4D-C1D-N1N-C6N
3	A	1534	NAP	C2D-C1D-N1N-C2N
3	A	1534	NAP	C2D-C1D-N1N-C6N
3	A	1534	NAP	C2N-C3N-C7N-O7N
3	A	1534	NAP	C2N-C3N-C7N-N7N
3	A	1534	NAP	C4N-C3N-C7N-N7N
3	B	1527	NAP	C5D-O5D-PN-O3
3	B	1527	NAP	C5D-O5D-PN-O1N
3	B	1527	NAP	C5D-O5D-PN-O2N
3	B	1527	NAP	C4D-C5D-O5D-PN
3	B	1527	NAP	O4D-C1D-N1N-C2N
3	B	1527	NAP	O4D-C1D-N1N-C6N

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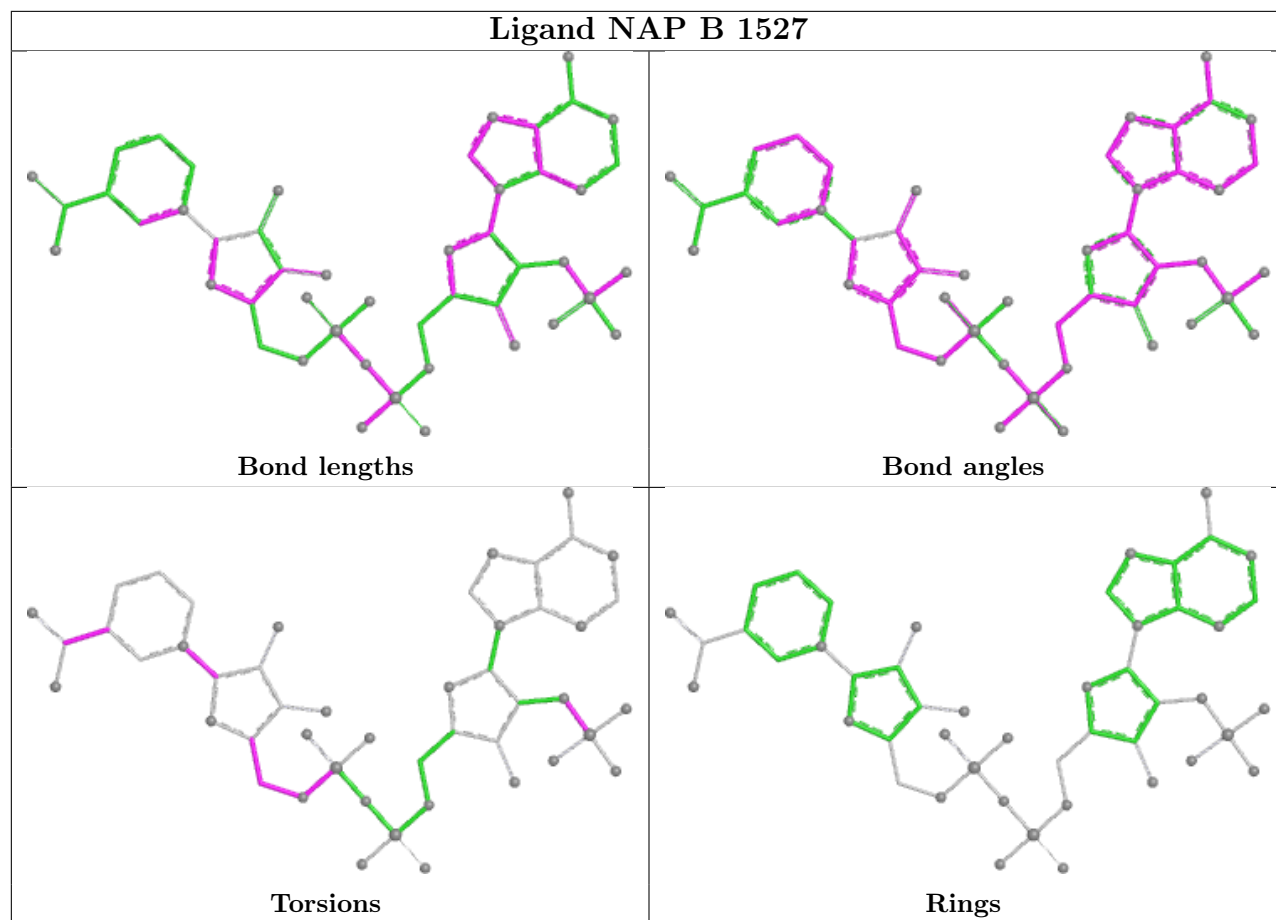
Mol	Chain	Res	Type	Atoms
3	B	1527	NAP	C2D-C1D-N1N-C2N
3	B	1527	NAP	C2D-C1D-N1N-C6N
3	B	1527	NAP	C2N-C3N-C7N-O7N
3	B	1527	NAP	C2N-C3N-C7N-N7N
3	A	1534	NAP	C4N-C3N-C7N-O7N
3	B	1527	NAP	C4N-C3N-C7N-O7N
3	B	1527	NAP	C4N-C3N-C7N-N7N
3	A	1534	NAP	C3D-C4D-C5D-O5D
3	A	1534	NAP	C4D-C5D-O5D-PN
3	A	1534	NAP	C4B-C5B-O5B-PA
3	A	1534	NAP	C3B-C4B-C5B-O5B
3	A	1534	NAP	PN-O3-PA-O5B
3	A	1534	NAP	PA-O3-PN-O5D
3	A	1534	NAP	O4B-C4B-C5B-O5B
3	A	1534	NAP	C5B-O5B-PA-O1A
3	A	1534	NAP	C5D-O5D-PN-O1N
3	B	1527	NAP	O4D-C4D-C5D-O5D
3	A	1534	NAP	PA-O3-PN-O2N
3	A	1534	NAP	C2B-O2B-P2B-O3X
3	A	1534	NAP	C2B-O2B-P2B-O1X
3	B	1527	NAP	C2B-O2B-P2B-O1X

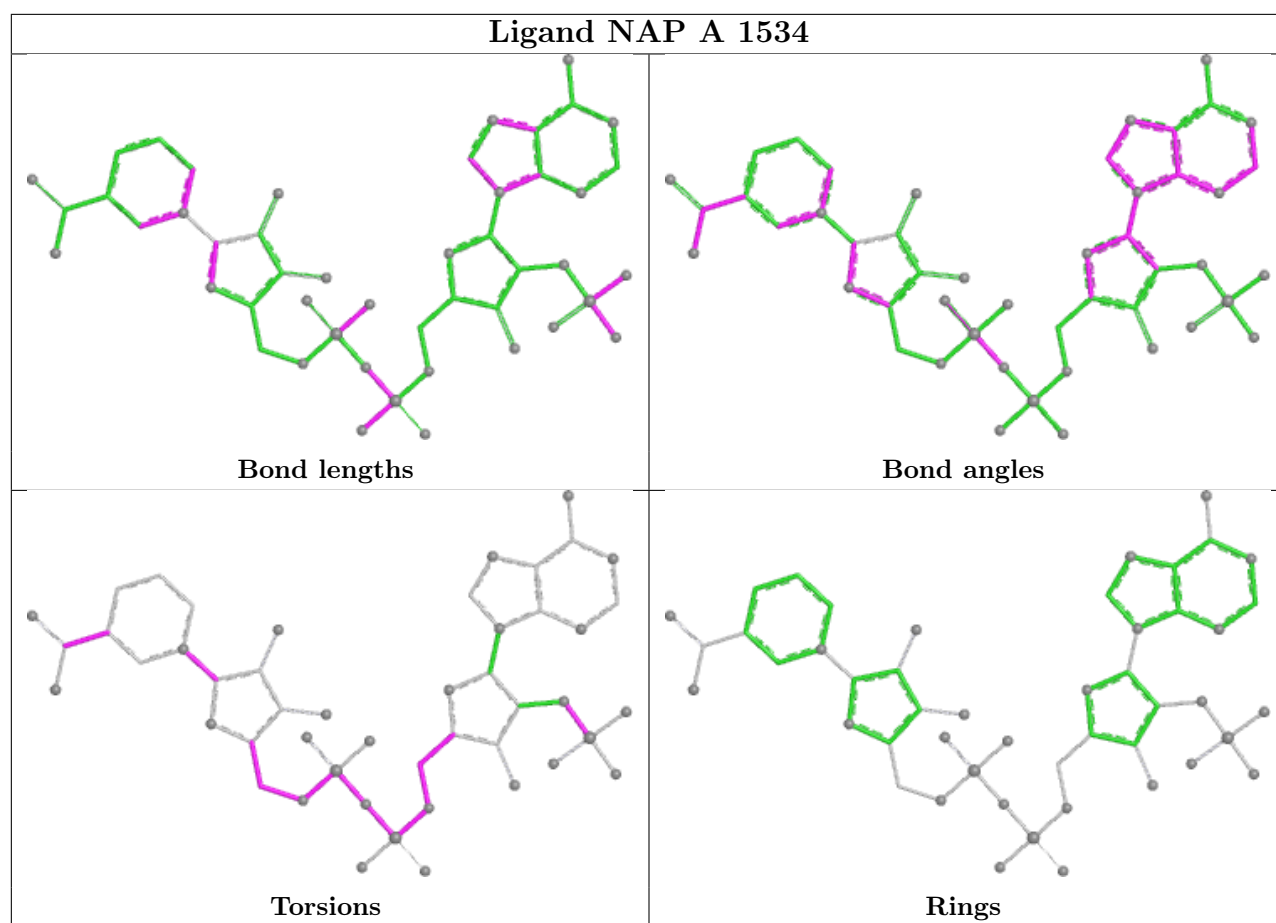
There are no ring outliers.

2 monomers are involved in 45 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1527	NAP	22	0
3	A	1534	NAP	23	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	529/534 (99%)	-0.58	9 (1%) 69 72	4, 8, 18, 43	1 (0%)
1	B	524/534 (98%)	-0.37	15 (2%) 53 56	4, 9, 22, 50	3 (0%)
All	All	1053/1068 (98%)	-0.48	24 (2%) 61 63	4, 9, 21, 50	4 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	HIS	4.7
1	B	523	THR	4.5
1	A	375	ASN	4.0
1	B	522	LEU	3.9
1	B	0	HIS	3.7
1	B	521	THR	3.6
1	A	446	ALA	3.4
1	B	394	THR	3.3
1	B	415	THR	3.3
1	B	526	THR	3.2
1	A	416	ASP	2.8
1	B	520	GLY	2.6
1	B	525	ALA	2.5
1	A	241	LEU	2.4
1	B	524	GLN	2.4
1	A	450	ARG	2.4
1	A	259	ASP	2.3
1	A	415	THR	2.3
1	B	517	ALA	2.2
1	B	259	ASP	2.2
1	B	241	LEU	2.2
1	B	414	THR	2.1
1	B	416	ASP	2.1
1	A	531	ALA	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](#)

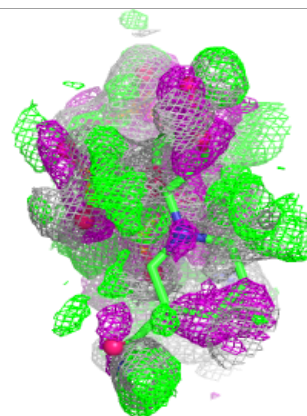
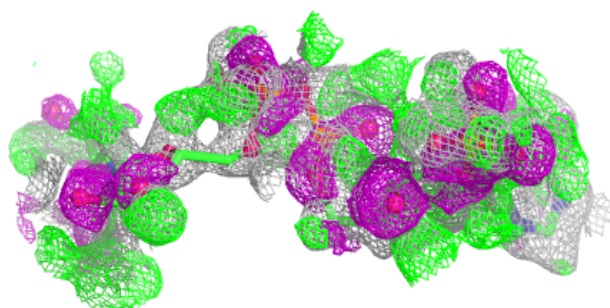
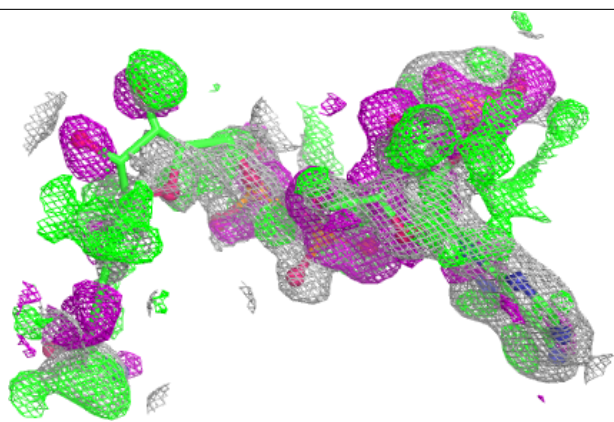
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAP	A	1534	48/48	0.71	0.22	12,19,28,33	18
3	NAP	B	1527	48/48	0.87	0.16	5,22,33,38	17
2	GOL	A	1533	6/6	0.96	0.05	8,11,11,12	0
2	GOL	B	1528	6/6	0.97	0.06	10,11,13,13	0
2	GOL	A	1532	6/6	0.98	0.05	5,9,12,14	0
2	GOL	B	1529	6/6	0.98	0.06	6,10,12,15	0

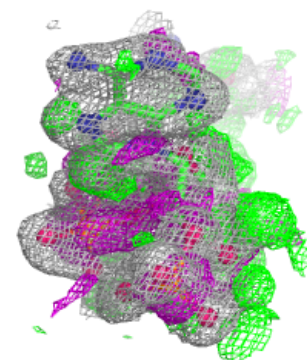
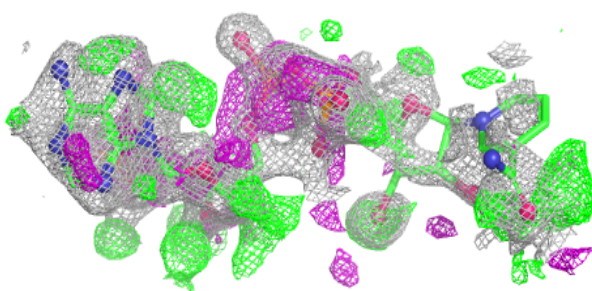
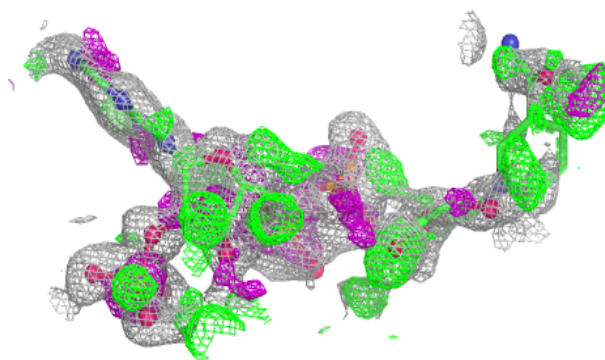
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAP A 1534:

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

**Electron density around NAP B 1527:**

$2mF_o-DF_c$ (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.