



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:28 AM UTC

PDB ID : 3SC6 / pdb_00003sc6
Title : 2.65 Angstrom resolution crystal structure of dTDP-4-dehydrorhamnose reductase (rfbD) from Bacillus anthracis str. Ames in complex with NADP
Authors : Halavaty, A.S.; Kuhn, M.; Shuvalova, L.; Minasov, G.; Peterson, S.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2011-06-07
Resolution : 2.65 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:

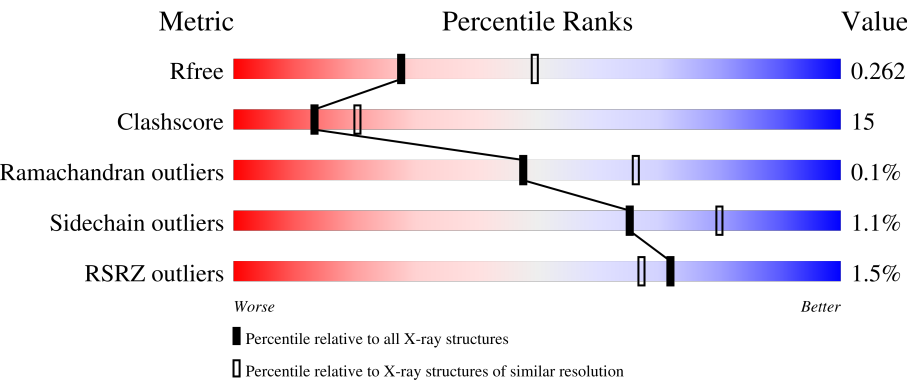
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	287	2% 69% 29% .
1	B	287	% 67% 30% ..
1	C	287	% 74% 23% ..
1	D	287	2% 67% 31% ..

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	287	2% 66% 30% ..
1	F	287	% 67% 29% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14334 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 dTDP-4-dehydrorhamnose reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	3	0
			2288	1470	386	425	7			
1	B	280	Total	C	N	O	S	0	1	0
			2260	1451	383	418	8			
1	C	281	Total	C	N	O	S	0	3	0
			2296	1474	391	423	8			
1	D	283	Total	C	N	O	S	0	2	0
			2293	1472	387	426	8			
1	E	278	Total	C	N	O	S	0	1	0
			2261	1455	381	418	7			
1	F	281	Total	C	N	O	S	0	1	0
			2274	1462	383	421	8			

There are 18 discrepancies between the modelled and reference sequences:

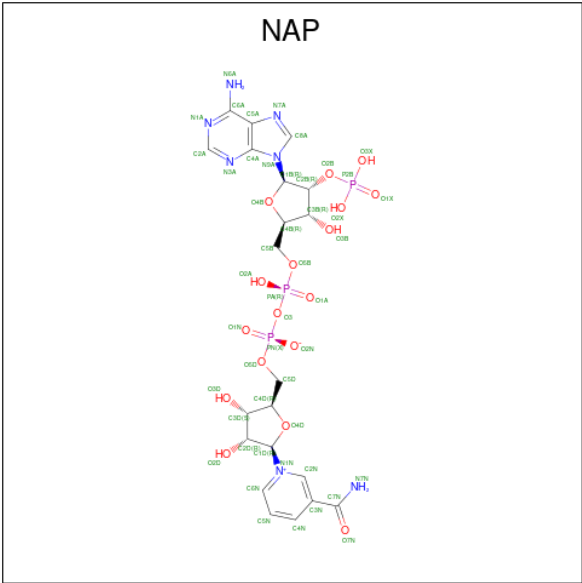
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q81TN9
A	-1	ASN	-	expression tag	UNP Q81TN9
A	0	ALA	-	expression tag	UNP Q81TN9
B	-2	SER	-	expression tag	UNP Q81TN9
B	-1	ASN	-	expression tag	UNP Q81TN9
B	0	ALA	-	expression tag	UNP Q81TN9
C	-2	SER	-	expression tag	UNP Q81TN9
C	-1	ASN	-	expression tag	UNP Q81TN9
C	0	ALA	-	expression tag	UNP Q81TN9
D	-2	SER	-	expression tag	UNP Q81TN9
D	-1	ASN	-	expression tag	UNP Q81TN9
D	0	ALA	-	expression tag	UNP Q81TN9
E	-2	SER	-	expression tag	UNP Q81TN9
E	-1	ASN	-	expression tag	UNP Q81TN9
E	0	ALA	-	expression tag	UNP Q81TN9
F	-2	SER	-	expression tag	UNP Q81TN9
F	-1	ASN	-	expression tag	UNP Q81TN9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	ALA	-	expression tag	UNP Q81TN9

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	1
			96	42	14	34	6		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	E	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	F	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		

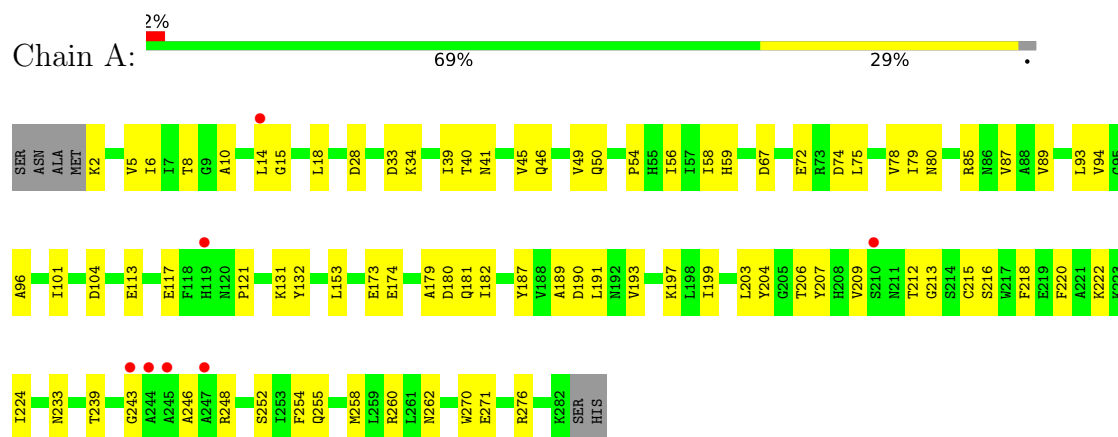
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total	O	0	0
			33	33		
4	B	50	Total	O	0	1
			50	50		
4	C	46	Total	O	0	0
			46	46		
4	D	32	Total	O	0	0
			32	32		
4	E	32	Total	O	0	0
			32	32		
4	F	47	Total	O	0	1
			48	48		

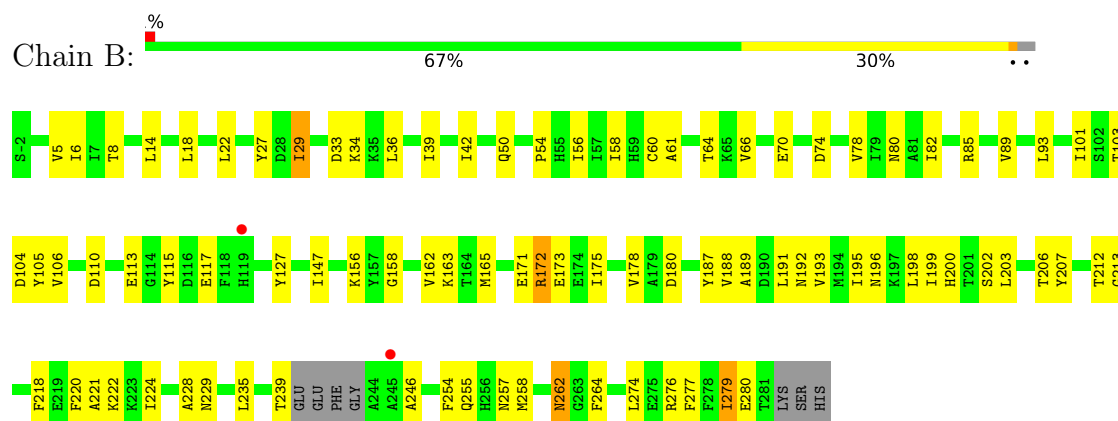
3 Residue-property plots

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.

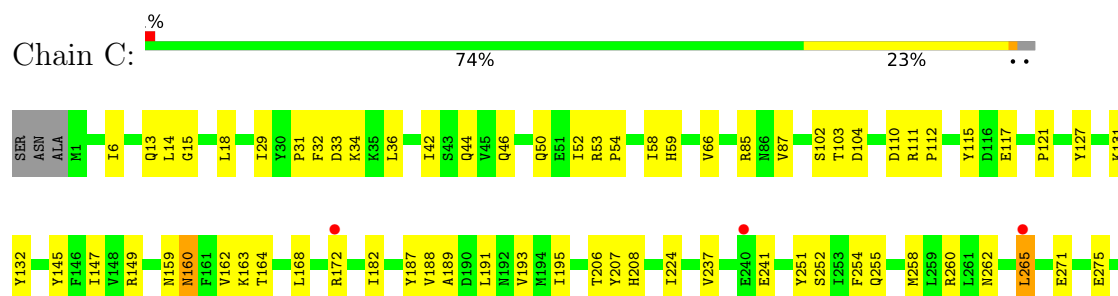
• Molecule 1: dTDP-4-dehydrorhamnose reductase



• Molecule 1: dTDP-4-dehydrorhamnose reductase

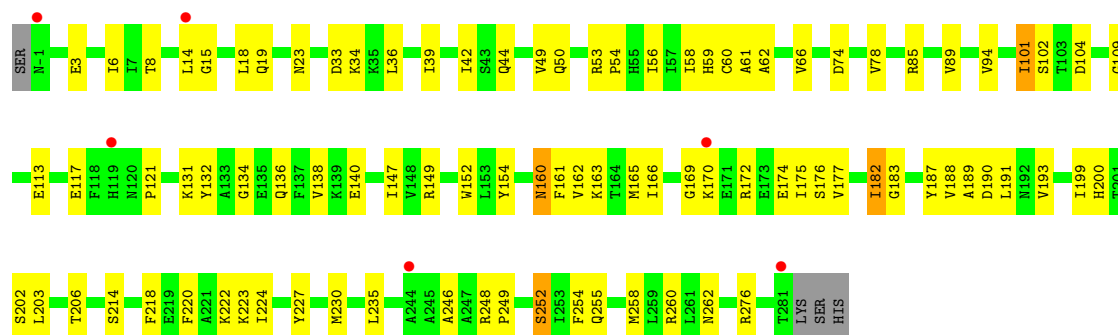


• Molecule 1: dTDP-4-dehydrorhamnose reductase

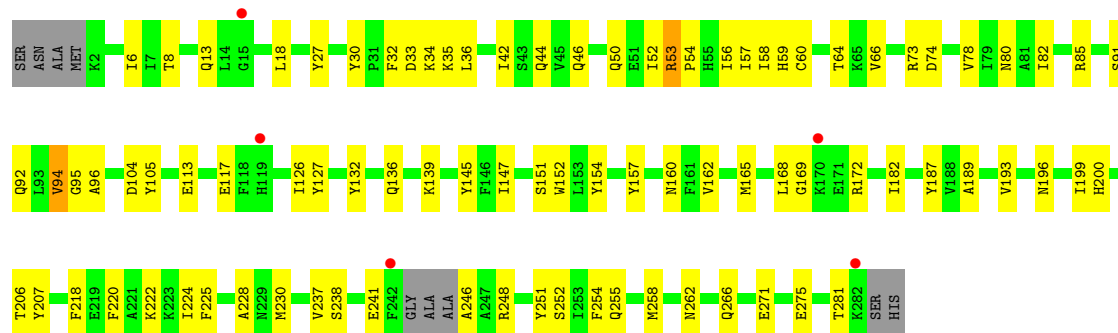




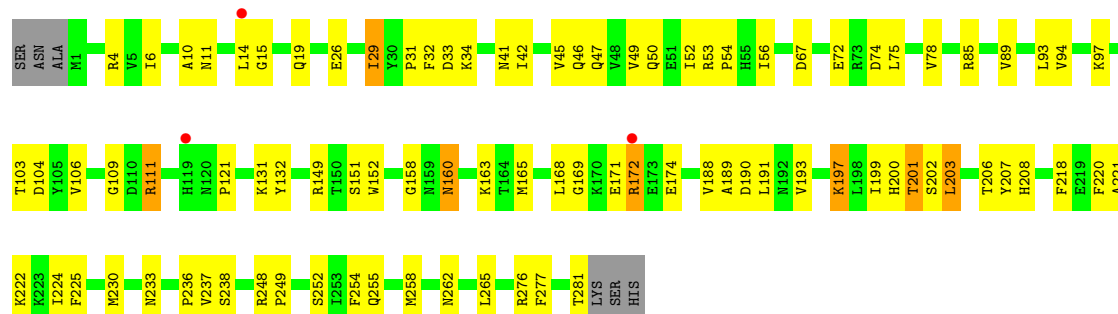
• Molecule 1: dTDP-4-dehydrorhamnose reductase



• Molecule 1: dTDP-4-dehydrorhamnose reductase



• Molecule 1: dTDP-4-dehydrorhamnose reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.45Å 113.34Å 144.95Å 90.00° 91.18° 90.00°	Depositor
Resolution (Å)	29.90 – 2.65 29.90 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.90-2.65) 99.9 (29.90-2.65)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 2.64Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.215 , 0.258 0.219 , 0.262	Depositor DCC
R_{free} test set	3745 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	56.4	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.044 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14334	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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: SO4, NAP

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	0.42	0/2342	0.90	3/3171 (0.1%)
1	B	0.46	0/2312	0.96	8/3130 (0.3%)
1	C	0.43	0/2350	0.92	3/3180 (0.1%)
1	D	0.44	0/2347	0.93	4/3179 (0.1%)
1	E	0.45	0/2314	0.92	4/3131 (0.1%)
1	F	0.45	0/2328	0.96	5/3152 (0.2%)
All	All	0.44	0/13993	0.93	27/18943 (0.1%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	255	GLN	N-CA-C	-8.18	96.42	109.59
1	A	173	GLU	N-CA-C	-7.42	104.35	113.41
1	A	255	GLN	N-CA-C	-6.77	98.21	109.24
1	F	172	ARG	N-CA-C	6.47	119.00	109.24
1	A	80	ASN	N-CA-C	6.12	118.03	111.36
1	D	66	VAL	N-CA-C	5.97	116.63	110.36
1	F	201	THR	N-CA-C	5.96	123.50	110.80
1	E	255	GLN	N-CA-C	-5.88	99.61	109.07
1	B	172	ARG	N-CA-C	5.75	125.17	113.20
1	C	66	VAL	N-CA-C	5.75	116.40	110.36
1	B	255	GLN	N-CA-C	-5.69	99.97	109.24
1	B	66	VAL	N-CA-C	5.67	116.12	110.23
1	D	255	GLN	N-CA-C	-5.58	99.31	108.41
1	E	145	TYR	N-CA-C	5.54	117.39	109.14
1	D	53	ARG	CA-C-N	5.48	125.46	120.03
1	D	53	ARG	C-N-CA	5.48	125.46	120.03
1	C	255	GLN	N-CA-C	-5.41	100.36	109.07
1	F	203	LEU	N-CA-C	-5.41	101.51	109.62

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	80	ASN	N-CA-C	5.38	118.06	111.82
1	B	257	ASN	N-CA-C	5.34	116.91	111.14
1	E	80	ASN	N-CA-C	5.29	117.95	111.82
1	E	169	GLY	N-CA-C	-5.16	108.26	114.66
1	F	171	GLU	N-CA-C	5.06	118.23	111.75
1	B	279	ILE	CB-CA-C	-5.05	105.32	112.14
1	B	106	VAL	N-CA-C	-5.05	106.21	113.07
1	C	145	TYR	N-CA-C	5.03	116.63	109.14
1	B	29	ILE	N-CA-C	5.00	115.95	108.54

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	2288	0	2253	64	0
1	B	2260	0	2232	81	0
1	C	2296	0	2265	59	0
1	D	2293	0	2257	79	0
1	E	2261	0	2228	70	0
1	F	2274	0	2241	73	0
2	A	48	0	25	3	0
2	B	96	0	50	5	0
2	C	48	0	25	2	0
2	D	48	0	25	0	0
2	E	48	0	25	0	0
2	F	48	0	25	1	0
3	A	10	0	0	0	0
3	B	20	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	1	0
3	E	20	0	0	0	0
3	F	15	0	0	0	0
4	A	33	0	0	0	0
4	B	50	0	0	2	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	46	0	0	1	0
4	D	32	0	0	2	0
4	E	32	0	0	4	0
4	F	48	0	0	2	0
All	All	14334	0	13651	411	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 15.

All (411) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:ASP:OD1	1:A:252:SER:HB2	1.63	0.97
1:F:104:ASP:OD1	1:F:252:SER:HB2	1.80	0.81
1:B:165:MET:HE2	1:B:221:ALA:HB1	1.64	0.78
1:C:160:ASN:C	1:C:160:ASN:HD22	1.91	0.78
1:E:281:THR:HG22	1:E:281:THR:O	1.84	0.75
1:F:160:ASN:ND2	1:F:163:LYS:H	1.85	0.75
1:C:160:ASN:ND2	1:C:163:LYS:H	1.83	0.75
1:A:39:ILE:HG22	2:A:285:NAP:N6A	2.02	0.74
1:E:220:PHE:O	1:E:224:ILE:HG13	1.87	0.74
1:D:147:ILE:HB	1:D:206:THR:HG22	1.70	0.73
1:E:73:ARG:NH2	1:E:126:ILE:HD11	2.04	0.72
1:D:14:LEU:HD23	1:D:188:VAL:HG13	1.72	0.72
1:E:53:ARG:HH11	1:E:53:ARG:CG	2.02	0.72
1:A:182:ILE:HG23	1:A:215:CYS:O	1.90	0.71
1:B:117:GLU:OE1	1:B:117:GLU:N	2.19	0.71
1:C:187:TYR:CE2	1:C:189:ALA:HB3	2.27	0.70
1:F:111:ARG:HH11	1:F:111:ARG:CG	2.04	0.70
1:F:169:GLY:HA3	1:F:230:MET:HE1	1.74	0.70
1:D:190:ASP:OD1	1:D:276:ARG:NH1	2.25	0.69
1:F:160:ASN:C	1:F:160:ASN:HD22	2.00	0.69
1:A:207:TYR:OH	1:A:262:ASN:ND2	2.25	0.69
1:A:2:LYS:HE3	1:A:28:ASP:OD1	1.92	0.69
1:E:165:MET:HE2	1:E:165:MET:HA	1.72	0.69
1:B:178:VAL:HG23	4:B:326:HOH:O	1.91	0.68
1:A:260:ARG:NH2	1:B:74:ASP:OD2	2.26	0.68
1:C:260:ARG:NH2	4:C:331:HOH:O	2.27	0.68
1:F:189:ALA:O	1:F:193:VAL:HG23	1.94	0.68
1:C:160:ASN:HD21	1:C:163:LYS:H	1.40	0.67
1:C:14:LEU:HD23	1:C:188:VAL:HG13	1.77	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:104:ASP:HB3	1:D:254:PHE:CZ	2.29	0.67
1:F:14:LEU:HD23	1:F:188:VAL:HG13	1.76	0.66
1:A:101:ILE:O	1:A:131:LYS:NZ	2.23	0.66
1:E:237:VAL:CG1	1:E:241:GLU:HB2	2.25	0.66
1:B:56:ILE:HD13	1:B:199:ILE:HB	1.78	0.66
1:B:189:ALA:O	1:B:193:VAL:HG23	1.96	0.66
1:C:104:ASP:HB3	1:C:254:PHE:CZ	2.31	0.66
1:D:56:ILE:HD13	1:D:199:ILE:HB	1.78	0.65
1:A:67:ASP:OD2	1:A:248:ARG:NH1	2.29	0.65
1:D:117:GLU:OE1	1:D:117:GLU:N	2.27	0.65
1:F:26:GLU:OE2	1:F:200:HIS:NE2	2.22	0.65
1:D:74:ASP:O	1:D:78:VAL:HG23	1.96	0.65
1:F:89:VAL:O	1:F:93:LEU:HD13	1.97	0.65
1:A:39:ILE:HG22	2:A:285:NAP:C6A	2.27	0.65
1:A:218:PHE:CE2	1:A:222:LYS:HE2	2.32	0.64
1:B:218:PHE:CZ	1:B:222:LYS:HE3	2.32	0.64
1:B:189:ALA:HB1	1:B:276:ARG:HH12	1.61	0.64
2:B:285[A]:NAP:N7N	2:B:285[A]:NAP:O1N	2.30	0.64
1:E:225:PHE:HD1	1:E:230:MET:HE3	1.62	0.64
1:D:160:ASN:C	1:D:160:ASN:HD22	2.06	0.64
1:B:191:LEU:HD21	1:B:195:ILE:HD11	1.79	0.64
1:B:191:LEU:CD2	1:B:195:ILE:HD11	2.28	0.64
1:B:82:ILE:HG23	1:B:85:ARG:HH12	1.62	0.63
1:E:18:LEU:HD21	1:E:58:ILE:HG21	1.81	0.63
1:C:191:LEU:HD21	1:C:195:ILE:HD11	1.80	0.63
1:B:196:ASN:O	1:B:200:HIS:ND1	2.32	0.62
1:F:193:VAL:HG12	1:F:197:LYS:HE2	1.81	0.62
1:B:50:GLN:NE2	1:B:93:LEU:CD2	2.63	0.62
1:B:74:ASP:O	1:B:78:VAL:HG23	1.98	0.62
1:F:160:ASN:HD21	1:F:163:LYS:H	1.44	0.62
1:D:160:ASN:ND2	1:D:163:LYS:H	1.97	0.62
1:D:248:ARG:NH2	3:D:287:SO4:O1	2.32	0.62
1:F:207:TYR:OH	1:F:262:ASN:ND2	2.34	0.61
1:C:191:LEU:HD23	1:C:195:ILE:HG13	1.83	0.61
1:D:203:LEU:HD23	1:E:42:ILE:HD12	1.82	0.61
1:F:131:LYS:NZ	2:F:285:NAP:O3D	2.29	0.61
1:D:147:ILE:HD12	1:D:206:THR:HG22	1.82	0.61
1:D:189:ALA:O	1:D:193:VAL:HG23	2.00	0.61
1:A:206:THR:O	1:A:258:MET:HG3	2.02	0.60
1:F:74:ASP:O	1:F:78:VAL:HG23	2.01	0.60
1:F:104:ASP:HB3	1:F:254:PHE:CZ	2.37	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:147:ILE:HB	1:C:206:THR:HG22	1.84	0.60
1:A:74:ASP:O	1:A:78:VAL:HG23	2.02	0.60
1:E:53:ARG:HH11	1:E:53:ARG:HG3	1.65	0.60
1:C:237:VAL:HG13	1:C:241:GLU:OE1	2.02	0.60
1:D:161:PHE:CZ	1:D:165:MET:HE3	2.36	0.60
1:A:180:ASP:OD2	1:A:239:THR:HG23	2.02	0.59
1:E:206:THR:O	1:E:258:MET:HG3	2.03	0.59
1:E:207:TYR:OH	1:E:262:ASN:ND2	2.34	0.59
1:F:103:THR:O	1:F:106:VAL:HG22	2.02	0.59
1:B:147:ILE:HD12	1:B:206:THR:HG22	1.84	0.59
1:E:266:GLN:NE2	4:E:321:HOH:O	2.35	0.59
1:A:203:LEU:HD22	1:B:85:ARG:HH21	1.67	0.59
1:C:191:LEU:CD2	1:C:195:ILE:HD11	2.33	0.59
1:F:29:ILE:HD12	1:F:29:ILE:N	2.18	0.59
1:A:189:ALA:O	1:A:193:VAL:HG23	2.03	0.59
1:D:182:ILE:N	1:D:182:ILE:HD12	2.18	0.59
1:A:5:VAL:HG12	1:A:56:ILE:HB	1.85	0.58
1:F:14:LEU:HD21	1:F:191:LEU:HD23	1.84	0.58
1:C:191:LEU:HD23	1:C:191:LEU:C	2.29	0.58
1:F:165:MET:HE2	1:F:221:ALA:HB1	1.85	0.58
1:A:104:ASP:HB3	1:A:254:PHE:CZ	2.38	0.58
1:A:6:ILE:HD12	1:A:54:PRO:HG3	1.85	0.58
1:B:191:LEU:C	1:B:191:LEU:HD23	2.28	0.58
1:F:206:THR:O	1:F:258:MET:HG3	2.03	0.58
1:E:196:ASN:O	1:E:200:HIS:CD2	2.57	0.58
1:B:27:TYR:OH	1:B:200:HIS:NE2	2.36	0.57
1:C:110:ASP:OD2	1:C:115:TYR:OH	2.23	0.57
1:B:110:ASP:OD2	1:B:115:TYR:OH	2.21	0.57
1:B:147:ILE:HB	1:B:206:THR:HG22	1.87	0.57
1:B:198:LEU:HD23	1:B:264:PHE:CZ	2.38	0.57
1:D:101:ILE:HD12	1:D:101:ILE:N	2.20	0.57
1:A:49:VAL:HG12	1:A:94:VAL:HG11	1.87	0.57
1:D:246:ALA:HB3	1:D:248:ARG:NH1	2.19	0.57
1:B:189:ALA:CB	1:B:276:ARG:HH12	2.18	0.57
1:B:18:LEU:HD21	1:B:58:ILE:HD13	1.86	0.56
1:B:64:THR:HG22	2:B:285[B]:NAP:H2D	1.86	0.56
1:B:14:LEU:HD12	1:B:188:VAL:HG13	1.87	0.56
1:C:46:GLN:O	1:C:50:GLN:HG2	2.05	0.56
1:D:220:PHE:O	1:D:224:ILE:HG13	2.04	0.56
1:E:56:ILE:HD13	1:E:199:ILE:HB	1.85	0.56
1:E:187:TYR:CE2	1:E:189:ALA:HB3	2.39	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:237:VAL:HG11	1:E:241:GLU:HB2	1.87	0.56
1:B:180:ASP:OD1	1:B:239:THR:HG22	2.06	0.56
1:D:117:GLU:HG2	1:D:258:MET:HG3	1.87	0.56
1:C:189:ALA:O	1:C:193:VAL:HG23	2.06	0.56
1:D:39:ILE:HD13	1:D:61:ALA:HB3	1.87	0.56
1:E:182:ILE:HD11	1:E:251:TYR:CD2	2.41	0.55
1:A:56:ILE:HD13	1:A:199:ILE:HB	1.87	0.55
1:B:70:GLU:HG3	1:B:246:ALA:HB2	1.88	0.55
1:C:265:LEU:N	1:C:265:LEU:HD12	2.20	0.55
1:D:191:LEU:O	1:D:191:LEU:HD23	2.06	0.55
1:D:202:SER:OG	1:E:42:ILE:HD11	2.04	0.55
1:D:36:LEU:O	1:D:44:GLN:NE2	2.40	0.55
1:F:149:ARG:HD2	1:F:208:HIS:CE1	2.41	0.55
1:A:117:GLU:OE1	1:A:117:GLU:N	2.35	0.55
1:B:156:LYS:HE3	1:B:280:GLU:HB3	1.88	0.55
1:C:18:LEU:HD21	1:C:58:ILE:HG21	1.88	0.55
1:C:162:VAL:HA	1:C:224:ILE:HD13	1.89	0.55
1:D:218:PHE:CZ	1:D:222:LYS:HE3	2.42	0.55
1:E:104:ASP:OD1	1:E:252:SER:HB3	2.07	0.55
1:E:225:PHE:CD1	1:E:230:MET:HE3	2.42	0.55
1:D:191:LEU:HD23	1:D:191:LEU:C	2.32	0.55
1:C:33:ASP:OD1	1:C:34:LYS:N	2.39	0.54
1:D:42:ILE:HG21	1:F:203:LEU:HD21	1.90	0.54
1:E:33:ASP:OD1	1:E:34:LYS:N	2.40	0.54
1:D:117:GLU:HG2	1:D:258:MET:CG	2.37	0.54
1:B:36:LEU:HD12	4:B:295:HOH:O	2.06	0.54
1:B:165:MET:HE1	1:B:175:ILE:HG21	1.89	0.54
1:D:23:ASN:N	4:D:317:HOH:O	2.39	0.54
1:E:6:ILE:HD11	1:E:52:ILE:HG21	1.89	0.54
1:A:18:LEU:HD21	1:A:58:ILE:HD13	1.90	0.54
1:C:172[A]:ARG:HD3	1:C:172[A]:ARG:N	2.23	0.54
1:D:169:GLY:C	1:D:230:MET:HE2	2.32	0.54
1:D:203:LEU:HD22	1:E:85:ARG:NH2	2.23	0.54
1:F:11:ASN:O	1:F:11:ASN:ND2	2.41	0.54
1:E:182:ILE:HD11	1:E:251:TYR:HD2	1.73	0.53
1:E:281:THR:O	1:E:281:THR:CG2	2.55	0.53
1:D:182:ILE:CG2	1:D:214:SER:OG	2.56	0.53
1:A:220:PHE:O	1:A:224:ILE:HG13	2.09	0.53
1:B:191:LEU:CD2	1:B:195:ILE:CD1	2.87	0.53
1:A:40:THR:CG2	1:A:79:ILE:HD13	2.38	0.53
1:E:18:LEU:CD2	1:E:58:ILE:HG21	2.39	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:149:ARG:HH11	1:F:208:HIS:HE1	1.55	0.53
1:A:33:ASP:OD1	1:A:34:LYS:N	2.35	0.53
1:A:46:GLN:O	1:A:50:GLN:HG2	2.08	0.53
1:A:41:ASN:O	1:A:45:VAL:HG23	2.09	0.52
1:F:67:ASP:OD1	1:F:248:ARG:NH1	2.27	0.52
1:B:50:GLN:NE2	1:B:93:LEU:HD21	2.25	0.52
1:D:33:ASP:OD1	1:D:34:LYS:N	2.37	0.52
1:D:262:ASN:HD21	1:E:82:ILE:HD12	1.74	0.52
1:F:104:ASP:C	1:F:106:VAL:H	2.18	0.52
1:F:165:MET:HE3	1:F:225:PHE:CZ	2.44	0.52
1:C:6:ILE:HD12	1:C:54:PRO:HG3	1.92	0.52
1:F:49:VAL:HG12	1:F:94:VAL:HG21	1.92	0.52
1:A:220:PHE:CE2	1:A:224:ILE:HD11	2.45	0.52
1:C:18:LEU:CD2	1:C:58:ILE:HG21	2.40	0.52
1:A:215:CYS:SG	1:A:220:PHE:HB2	2.50	0.52
1:B:5:VAL:HG22	1:B:56:ILE:HB	1.92	0.52
1:D:152:TRP:CZ2	1:D:183:GLY:HA3	2.44	0.52
1:E:94:VAL:O	1:E:94:VAL:CG1	2.58	0.52
1:A:14:LEU:HD21	1:A:153:LEU:HD21	1.91	0.51
1:B:64:THR:CG2	2:B:285[B]:NAP:H2D	2.40	0.51
1:F:111:ARG:HH11	1:F:111:ARG:HG2	1.74	0.51
1:C:271:GLU:O	1:C:275:GLU:HG3	2.09	0.51
1:E:246:ALA:HB3	1:E:248:ARG:NH1	2.26	0.51
1:A:181:GLN:HE21	1:A:239:THR:CG2	2.24	0.51
1:C:191:LEU:CD2	1:C:195:ILE:CD1	2.88	0.51
1:D:85:ARG:O	1:D:89:VAL:HG23	2.10	0.51
1:C:117:GLU:OE1	1:C:117:GLU:N	2.38	0.51
1:C:160:ASN:O	1:C:164:THR:HG23	2.10	0.51
1:D:8:THR:O	1:D:60:CYS:HB2	2.10	0.51
1:B:189:ALA:HB1	1:B:276:ARG:NH1	2.24	0.51
1:E:117:GLU:OE1	1:E:117:GLU:N	2.40	0.51
1:B:70:GLU:CG	1:B:246:ALA:HB2	2.40	0.50
1:C:237:VAL:CG1	1:C:241:GLU:HB2	2.41	0.50
1:F:111:ARG:HH11	1:F:111:ARG:HG3	1.76	0.50
1:A:59:HIS:CE1	1:A:87:VAL:HG21	2.46	0.50
1:C:191:LEU:HD23	1:C:191:LEU:O	2.11	0.50
1:B:220:PHE:O	1:B:224:ILE:HG13	2.12	0.50
1:F:218:PHE:CZ	1:F:222:LYS:HE3	2.45	0.50
1:D:3:GLU:N	1:D:3:GLU:OE1	2.44	0.50
1:D:161:PHE:CE1	1:D:165:MET:HE2	2.47	0.50
1:B:104:ASP:HB3	1:B:254:PHE:CZ	2.47	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:THR:OG1	1:D:59:HIS:HA	2.11	0.50
1:B:191:LEU:HD23	1:B:195:ILE:HG13	1.93	0.50
1:E:271:GLU:O	1:E:275:GLU:HG3	2.12	0.50
1:F:169:GLY:CA	1:F:230:MET:HE1	2.40	0.50
1:B:113:GLU:OE2	1:B:113:GLU:N	2.41	0.49
1:D:223:LYS:HE3	1:D:227:TYR:CE2	2.47	0.49
1:B:189:ALA:CB	1:B:276:ARG:NH1	2.75	0.49
1:B:202:SER:OG	1:C:42:ILE:HD11	2.12	0.49
1:E:228:ALA:CB	1:E:230:MET:HE2	2.43	0.49
1:A:203:LEU:HD22	1:B:85:ARG:NH2	2.27	0.49
1:F:46:GLN:O	1:F:50:GLN:HG2	2.12	0.49
1:A:180:ASP:CG	1:A:239:THR:HG23	2.37	0.49
1:A:246:ALA:HB3	1:A:248:ARG:NH1	2.28	0.49
1:D:174:GLU:OE1	1:D:235:LEU:HD11	2.12	0.49
1:D:187:TYR:CE2	1:D:189:ALA:HB3	2.47	0.49
1:E:105:TYR:HD2	1:E:127:TYR:CD2	2.30	0.49
1:E:168:LEU:HD22	1:E:172:ARG:HH22	1.77	0.49
1:B:235:LEU:HD23	1:F:236:PRO:HG3	1.95	0.49
1:D:49:VAL:HG12	1:D:94:VAL:HG11	1.92	0.49
1:B:33:ASP:OD1	1:B:34:LYS:N	2.44	0.49
1:B:206:THR:O	1:B:258:MET:HG3	2.12	0.49
1:C:237:VAL:HG13	1:C:241:GLU:HB2	1.95	0.49
1:D:154:TYR:CD2	1:D:162:VAL:HG21	2.47	0.49
1:E:13:GLN:HG2	1:E:157:TYR:O	2.12	0.48
1:F:29:ILE:HG22	1:F:31:PRO:HD3	1.94	0.48
1:B:262:ASN:N	1:B:262:ASN:HD22	2.11	0.48
1:E:6:ILE:HD13	1:E:30:TYR:HB2	1.95	0.48
1:F:47:GLN:HG2	4:F:296:HOH:O	2.13	0.48
1:D:160:ASN:HD21	1:D:163:LYS:H	1.60	0.48
1:F:10:ALA:HA	1:F:15:GLY:HA3	1.95	0.48
1:F:14:LEU:CD2	1:F:188:VAL:HG13	2.42	0.48
1:E:95:GLY:HA2	4:E:304:HOH:O	2.13	0.48
1:C:14:LEU:CD2	1:C:188:VAL:HG13	2.43	0.48
1:D:161:PHE:CE1	1:D:165:MET:CE	2.96	0.48
1:C:18:LEU:HD21	1:C:58:ILE:HD13	1.96	0.48
1:E:32:PHE:HZ	1:E:52:ILE:HD11	1.79	0.48
1:E:104:ASP:HB3	1:E:254:PHE:CZ	2.48	0.48
1:F:109:GLY:HA3	1:F:249:PRO:HB3	1.96	0.48
1:C:36:LEU:O	1:C:44:GLN:NE2	2.47	0.47
1:C:102:SER:HB3	1:C:149:ARG:HG2	1.95	0.47
1:A:203:LEU:HD21	1:B:42:ILE:HG21	1.95	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:ILE:HD13	1:B:61:ALA:CB	2.43	0.47
1:A:204:TYR:O	1:B:85:ARG:HD2	2.14	0.47
1:D:121:PRO:HG2	1:D:132:TYR:CE1	2.50	0.47
1:B:101:ILE:HG22	2:B:285[A]:NAP:H1D	1.97	0.47
1:C:160:ASN:C	1:C:160:ASN:ND2	2.64	0.47
1:C:168:LEU:HD22	1:C:172[A]:ARG:NH2	2.29	0.47
1:D:252:SER:OG	1:D:252:SER:O	2.31	0.47
1:F:103:THR:OG1	1:F:104:ASP:N	2.47	0.47
1:B:8:THR:O	1:B:60:CYS:HB2	2.14	0.47
1:B:173:GLU:N	1:B:173:GLU:OE1	2.48	0.47
1:C:149:ARG:HD2	1:C:208:HIS:CD2	2.50	0.47
1:D:109:GLY:HA3	1:D:249:PRO:CB	2.45	0.47
1:E:8:THR:OG1	1:E:59:HIS:HA	2.15	0.47
1:E:154:TYR:CD2	1:E:162:VAL:HG21	2.50	0.47
1:F:190:ASP:OD1	1:F:276:ARG:NH2	2.47	0.47
1:A:39:ILE:HG22	2:A:285:NAP:H62A	1.80	0.47
1:D:166:ILE:O	1:D:230:MET:HE1	2.14	0.47
1:D:113:GLU:OE2	1:D:113:GLU:N	2.36	0.46
1:A:174:GLU:HG2	1:A:233:ASN:HB2	1.97	0.46
1:B:274:LEU:O	1:B:277:PHE:HB3	2.15	0.46
1:A:181:GLN:HE21	1:A:239:THR:HG22	1.79	0.46
1:B:203:LEU:HD22	1:C:85:ARG:NH2	2.31	0.46
1:D:147:ILE:HD12	1:D:206:THR:CG2	2.45	0.46
1:D:260:ARG:NH2	1:E:74:ASP:OD2	2.46	0.46
1:B:228:ALA:O	1:B:229:ASN:C	2.58	0.46
1:F:41:ASN:O	1:F:45:VAL:HG23	2.15	0.46
1:A:215:CYS:SG	1:A:216:SER:N	2.87	0.46
1:D:14:LEU:O	1:D:15:GLY:C	2.59	0.46
1:B:22:LEU:HB2	1:B:29:ILE:HD11	1.98	0.46
1:D:152:TRP:CE2	1:D:183:GLY:HA3	2.51	0.46
1:A:191:LEU:HD12	1:A:209:VAL:HG11	1.98	0.45
1:D:42:ILE:HD12	1:F:203:LEU:HD23	1.97	0.45
1:E:237:VAL:HG12	1:E:238:SER:N	2.31	0.45
1:F:94:VAL:HG12	1:F:94:VAL:O	2.15	0.45
1:A:89:VAL:O	1:A:93:LEU:HD13	2.16	0.45
1:C:265:LEU:CD1	1:C:265:LEU:H	2.29	0.45
1:D:131:LYS:HE2	1:D:131:LYS:HA	1.98	0.45
1:E:53:ARG:HH11	1:E:53:ARG:HG2	1.80	0.45
1:F:168:LEU:HD13	1:F:172:ARG:HH21	1.81	0.45
1:F:237:VAL:HG12	1:F:238:SER:N	2.31	0.45
1:E:33:ASP:OD2	1:E:35:LYS:NZ	2.49	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:GLU:OE2	1:A:113:GLU:N	2.36	0.45
1:E:27:TYR:OH	1:E:200:HIS:CE1	2.70	0.45
1:E:271:GLU:OE1	1:E:271:GLU:N	2.46	0.45
1:D:165:MET:HE1	1:D:177:VAL:HG22	1.99	0.45
1:E:139:LYS:HG3	1:E:147:ILE:HD11	1.99	0.45
1:E:189:ALA:O	1:E:193:VAL:HG23	2.17	0.45
1:A:190:ASP:OD1	1:A:276:ARG:NH1	2.44	0.45
1:E:222:LYS:NZ	1:E:222:LYS:HB3	2.31	0.45
1:E:218:PHE:CZ	1:E:222:LYS:HD2	2.52	0.45
1:F:72:GLU:OE2	1:F:75:LEU:HD23	2.16	0.45
1:F:160:ASN:ND2	1:F:160:ASN:C	2.69	0.44
1:C:265:LEU:HD12	1:C:265:LEU:H	1.81	0.44
1:F:19:GLN:NE2	1:F:31:PRO:HG3	2.32	0.44
1:A:39:ILE:CD1	1:A:59:HIS:CE1	3.00	0.44
1:A:212:THR:HG22	1:A:213:GLY:N	2.31	0.44
1:E:57:ILE:HD12	1:E:91:SER:OG	2.18	0.44
1:D:154:TYR:O	1:D:188:VAL:HG23	2.17	0.44
1:E:73:ARG:NE	4:E:300:HOH:O	2.49	0.44
1:E:151:SER:O	1:E:152:TRP:C	2.61	0.44
1:F:169:GLY:C	1:F:230:MET:HE1	2.41	0.44
1:C:207:TYR:OH	1:C:262:ASN:ND2	2.51	0.44
1:F:97:LYS:NZ	1:F:202:SER:HB3	2.32	0.44
1:E:64:THR:HG22	1:E:64:THR:O	2.17	0.44
1:B:203:LEU:HD23	1:C:42:ILE:HD12	1.99	0.44
1:C:13:GLN:NE2	1:C:159:ASN:O	2.50	0.44
1:F:151:SER:O	1:F:152:TRP:C	2.60	0.44
1:F:201:THR:HB	1:F:203:LEU:HG	1.99	0.44
1:C:32:PHE:HZ	1:C:52:ILE:HD11	1.83	0.44
1:F:42:ILE:HG22	4:F:318:HOH:O	2.17	0.44
1:B:103:THR:OG1	1:B:105:TYR:HD2	2.01	0.43
1:B:191:LEU:HD23	1:B:191:LEU:O	2.18	0.43
1:B:85:ARG:O	1:B:89:VAL:HG23	2.18	0.43
1:B:191:LEU:CD2	1:B:191:LEU:C	2.91	0.43
1:D:18:LEU:HD21	1:D:58:ILE:HG21	1.99	0.43
1:C:131:LYS:NZ	2:C:285:NAP:O3D	2.48	0.43
1:C:265:LEU:N	1:C:265:LEU:CD1	2.82	0.43
1:C:59:HIS:CE1	1:C:87:VAL:HG21	2.54	0.43
1:E:168:LEU:HD22	1:E:172:ARG:NH2	2.33	0.43
1:F:201:THR:O	1:F:202:SER:OG	2.23	0.43
1:F:220:PHE:O	1:F:224:ILE:HG13	2.18	0.43
1:B:14:LEU:CD1	1:B:188:VAL:HG13	2.48	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:LEU:O	1:C:15:GLY:C	2.61	0.43
1:D:39:ILE:HD13	1:D:61:ALA:CB	2.48	0.43
1:A:33:ASP:CG	1:A:34:LYS:H	2.24	0.43
1:C:121:PRO:HG2	1:C:132:TYR:CE1	2.54	0.43
1:C:206:THR:O	1:C:258:MET:HG3	2.19	0.43
1:D:175:ILE:HG22	1:D:176:SER:N	2.33	0.43
1:E:18:LEU:HD21	1:E:58:ILE:HD13	2.00	0.43
1:F:111:ARG:CG	1:F:111:ARG:NH1	2.73	0.43
1:B:218:PHE:CE2	1:B:222:LYS:HE3	2.53	0.43
1:F:52:ILE:O	1:F:53:ARG:C	2.61	0.43
1:B:187:TYR:CE2	1:B:189:ALA:HB3	2.53	0.43
1:F:29:ILE:HD12	1:F:29:ILE:H	1.83	0.43
1:F:174:GLU:HG2	1:F:233:ASN:HB3	1.99	0.43
1:D:109:GLY:HA3	1:D:249:PRO:HB3	2.01	0.43
1:E:66:VAL:HG12	1:E:246:ALA:CB	2.49	0.43
1:D:160:ASN:C	1:D:160:ASN:ND2	2.77	0.42
1:E:8:THR:O	1:E:60:CYS:HB2	2.19	0.42
1:E:92:GLN:C	1:E:92:GLN:OE1	2.62	0.42
1:E:132:TYR:OH	1:E:136:GLN:NE2	2.51	0.42
1:B:39:ILE:HD13	1:B:61:ALA:HB3	2.00	0.42
1:A:218:PHE:CZ	1:A:222:LYS:HE2	2.53	0.42
1:D:134:GLY:O	1:D:138:VAL:HG23	2.19	0.42
1:E:52:ILE:O	1:E:53:ARG:C	2.62	0.42
1:A:203:LEU:HD23	1:B:42:ILE:HD12	2.01	0.42
1:B:207:TYR:CE1	1:B:258:MET:HB3	2.54	0.42
1:D:50:GLN:NE2	1:D:50:GLN:HA	2.35	0.42
1:D:61:ALA:O	1:D:62:ALA:HB2	2.19	0.42
1:D:182:ILE:N	1:D:182:ILE:CD1	2.82	0.42
1:A:18:LEU:CD2	1:A:58:ILE:HG21	2.50	0.42
1:B:18:LEU:HD13	1:B:192:ASN:CG	2.45	0.42
1:B:165:MET:HE2	1:B:221:ALA:CB	2.44	0.42
1:F:56:ILE:HD13	1:F:199:ILE:HB	2.01	0.42
1:A:8:THR:OG1	1:A:59:HIS:HA	2.20	0.42
1:B:50:GLN:NE2	1:B:93:LEU:HD23	2.33	0.42
1:B:180:ASP:CG	1:B:239:THR:HG22	2.45	0.42
1:D:102:SER:HB3	1:D:149:ARG:HG2	2.01	0.42
1:A:193:VAL:O	1:A:197:LYS:HG3	2.20	0.42
1:C:111:ARG:HA	1:C:112:PRO:HD3	1.93	0.42
1:E:54:PRO:O	1:E:96:ALA:HB2	2.19	0.42
1:A:85:ARG:C	1:A:85:ARG:HD2	2.45	0.42
1:B:50:GLN:HE21	1:B:93:LEU:CD2	2.32	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:GLN:O	1:D:140:GLU:HG3	2.20	0.42
1:E:193:VAL:HG23	4:E:320:HOH:O	2.20	0.42
1:A:10:ALA:HA	1:A:15:GLY:HA3	2.02	0.41
1:F:121:PRO:HG2	1:F:132:TYR:CE1	2.55	0.41
1:D:169:GLY:HA3	1:D:230:MET:HE2	2.02	0.41
1:A:54:PRO:O	1:A:96:ALA:HB2	2.20	0.41
1:C:52:ILE:O	1:C:53:ARG:C	2.63	0.41
1:D:101:ILE:N	1:D:101:ILE:CD1	2.83	0.41
1:A:121:PRO:HG2	1:A:132:TYR:CE1	2.56	0.41
1:C:131:LYS:NZ	2:C:285:NAP:O2D	2.49	0.41
1:D:170:LYS:HG3	1:D:170:LYS:O	2.20	0.41
1:D:191:LEU:C	1:D:191:LEU:CD2	2.93	0.41
1:E:13:GLN:OE1	1:E:160:ASN:HB3	2.21	0.41
1:F:33:ASP:OD1	1:F:34:LYS:N	2.48	0.41
1:F:85:ARG:HD2	1:F:85:ARG:C	2.45	0.41
1:F:277:PHE:O	1:F:281:THR:OG1	2.32	0.41
1:C:103:THR:HG21	1:C:127:TYR:CE2	2.56	0.41
1:E:36:LEU:O	1:E:44:GLN:NE2	2.54	0.41
1:E:94:VAL:O	1:E:94:VAL:HG13	2.20	0.41
1:A:187:TYR:CE2	1:A:189:ALA:HB3	2.55	0.41
1:C:104:ASP:OD1	1:C:252:SER:HB3	2.20	0.41
1:F:4:ARG:HB2	1:F:54:PRO:HA	2.02	0.41
1:F:158:GLY:O	1:F:163:LYS:NZ	2.51	0.41
1:C:191:LEU:CD2	1:C:191:LEU:C	2.94	0.41
1:A:179:ALA:HB2	1:A:218:PHE:HB2	2.02	0.41
1:A:6:ILE:CD1	1:A:54:PRO:HG3	2.49	0.41
1:A:72:GLU:OE2	1:A:75:LEU:HD23	2.20	0.41
1:B:6:ILE:HD12	1:B:54:PRO:HG3	2.02	0.41
1:B:171:GLU:O	1:B:172:ARG:HB2	2.21	0.41
1:B:276:ARG:HA	1:B:279:ILE:HD12	2.03	0.41
1:D:50:GLN:HA	1:D:50:GLN:HE21	1.86	0.41
1:E:78:VAL:O	1:E:82:ILE:HB	2.21	0.41
1:F:53:ARG:HG3	1:F:94:VAL:HG13	2.03	0.41
1:B:103:THR:HG21	1:B:127:TYR:CE2	2.56	0.40
2:B:285[B]:NAP:O2A	2:B:285[B]:NAP:O3B	2.35	0.40
1:C:182:ILE:O	1:C:251:TYR:HA	2.21	0.40
1:F:32:PHE:HZ	1:F:52:ILE:HD11	1.87	0.40
1:B:162:VAL:HA	1:B:224:ILE:HD13	2.03	0.40
1:D:15:GLY:O	1:D:19:GLN:HB2	2.21	0.40
1:D:23:ASN:HB3	4:D:317:HOH:O	2.21	0.40
1:E:46:GLN:O	1:E:50:GLN:HG2	2.20	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:THR:HG22	1:B:213:GLY:N	2.36	0.40
1:C:29:ILE:HG22	1:C:31:PRO:HD3	2.03	0.40
1:F:6:ILE:HD12	1:F:54:PRO:HG3	2.03	0.40
1:A:5:VAL:O	1:A:5:VAL:HG23	2.21	0.40
1:A:39:ILE:HD11	1:A:59:HIS:CE1	2.55	0.40
1:A:270:TRP:CE2	1:A:271:GLU:HG3	2.56	0.40
1:B:158:GLY:O	1:B:163:LYS:NZ	2.53	0.40
1:D:6:ILE:HD12	1:D:54:PRO:HG3	2.03	0.40
1:F:165:MET:HE3	1:F:225:PHE:CE2	2.56	0.40
1:B:207:TYR:CD1	1:B:258:MET:HB3	2.56	0.40
1:D:199:ILE:HG13	1:D:200:HIS:CD2	2.57	0.40
1:F:29:ILE:N	1:F:29:ILE:CD1	2.83	0.40
1:F:104:ASP:C	1:F:106:VAL:N	2.78	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	282/287 (98%)	265 (94%)	16 (6%)	1 (0%)	30	45
1	B	277/287 (96%)	266 (96%)	11 (4%)	0	100	100
1	C	282/287 (98%)	274 (97%)	8 (3%)	0	100	100
1	D	283/287 (99%)	266 (94%)	17 (6%)	0	100	100
1	E	275/287 (96%)	264 (96%)	11 (4%)	0	100	100
1	F	280/287 (98%)	266 (95%)	14 (5%)	0	100	100
All	All	1679/1722 (98%)	1601 (95%)	77 (5%)	1 (0%)	48	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	GLY

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	245/247 (99%)	245 (100%)	0	100	100
1	B	242/247 (98%)	241 (100%)	1 (0%)	84	93
1	C	245/247 (99%)	243 (99%)	2 (1%)	73	85
1	D	245/247 (99%)	240 (98%)	5 (2%)	48	69
1	E	243/247 (98%)	240 (99%)	3 (1%)	63	79
1	F	243/247 (98%)	238 (98%)	5 (2%)	47	69
All	All	1463/1482 (99%)	1447 (99%)	16 (1%)	65	80

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

Mol	Chain	Res	Type
1	B	262	ASN
1	C	160	ASN
1	C	265	LEU
1	D	101	ILE
1	D	160	ASN
1	D	172	ARG
1	D	182	ILE
1	D	252	SER
1	E	53	ARG
1	E	94	VAL
1	E	113	GLU
1	F	29	ILE
1	F	111	ARG
1	F	160	ASN
1	F	197	LYS
1	F	265	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	47	GLN
1	A	59	HIS
1	A	68	GLN
1	A	92	GLN
1	A	181	GLN
1	A	211	ASN
1	A	255	GLN
1	A	262	ASN
1	B	13	GLN
1	B	47	GLN
1	B	92	GLN
1	B	181	GLN
1	B	211	ASN
1	B	229	ASN
1	B	266	GLN
1	C	47	GLN
1	C	92	GLN
1	C	108	GLN
1	C	160	ASN
1	C	255	GLN
1	C	266	GLN
1	D	92	GLN
1	D	160	ASN
1	D	181	GLN
1	D	211	ASN
1	D	266	GLN
1	E	47	GLN
1	E	120	ASN
1	E	208	HIS
1	E	211	ASN
1	F	11	ASN
1	F	19	GLN
1	F	68	GLN
1	F	136	GLN
1	F	160	ASN
1	F	192	ASN
1	F	208	HIS
1	F	211	ASN
1	F	255	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 ⓘ

24 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	E	287	-	4,4,4	0.23	0	6,6,6	0.11	0
3	SO4	A	287	-	4,4,4	0.24	0	6,6,6	0.07	0
2	NAP	C	285	-	50,52,52	1.00	6 (12%)	71,80,80	1.46	9 (12%)
3	SO4	B	286	-	4,4,4	0.24	0	6,6,6	0.17	0
3	SO4	D	287	-	4,4,4	0.23	0	6,6,6	0.14	0
3	SO4	E	288	-	4,4,4	0.25	0	6,6,6	0.11	0
2	NAP	B	285[A]	-	50,52,52	0.97	5 (10%)	71,80,80	1.56	10 (14%)
2	NAP	D	285	-	50,52,52	1.04	6 (12%)	71,80,80	1.58	10 (14%)
3	SO4	F	287	-	4,4,4	0.25	0	6,6,6	0.15	0
3	SO4	F	286	-	4,4,4	0.22	0	6,6,6	0.13	0
3	SO4	B	289	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	C	286	-	4,4,4	0.23	0	6,6,6	0.14	0
3	SO4	C	287	-	4,4,4	0.24	0	6,6,6	0.12	0
3	SO4	A	286	-	4,4,4	0.23	0	6,6,6	0.09	0
3	SO4	B	288	-	4,4,4	0.26	0	6,6,6	0.08	0
2	NAP	B	285[B]	-	50,52,52	1.00	6 (12%)	71,80,80	1.52	10 (14%)
3	SO4	D	286	-	4,4,4	0.23	0	6,6,6	0.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	E	286	-	4,4,4	0.25	0	6,6,6	0.13	0
2	NAP	E	285	-	50,52,52	1.01	6 (12%)	71,80,80	1.54	11 (15%)
3	SO4	F	288	-	4,4,4	0.25	0	6,6,6	0.11	0
2	NAP	F	285	-	50,52,52	1.04	6 (12%)	71,80,80	1.49	9 (12%)
2	NAP	A	285	-	50,52,52	1.03	6 (12%)	71,80,80	1.54	9 (12%)
3	SO4	B	287	-	4,4,4	0.25	0	6,6,6	0.09	0
3	SO4	E	289	-	4,4,4	0.22	0	6,6,6	0.12	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	C	285	-	-	13/35/67/67	0/5/5/5
2	NAP	D	285	-	-	6/35/67/67	0/5/5/5
2	NAP	B	285[B]	-	-	11/35/67/67	0/5/5/5
2	NAP	F	285	-	-	10/35/67/67	0/5/5/5
2	NAP	A	285	-	-	12/35/67/67	0/5/5/5
2	NAP	B	285[A]	-	-	3/35/67/67	0/5/5/5
2	NAP	E	285	-	-	5/35/67/67	0/5/5/5

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	285	NAP	C5A-C6A	2.75	1.48	1.41
2	A	285	NAP	PN-O3	2.74	1.62	1.59
2	B	285[B]	NAP	C5A-C6A	2.69	1.48	1.41
2	D	285	NAP	PA-O3	2.68	1.62	1.59
2	C	285	NAP	C5A-C6A	2.64	1.48	1.41
2	F	285	NAP	C5A-N7A	-2.61	1.34	1.39
2	D	285	NAP	C5A-C6A	2.60	1.48	1.41
2	B	285[A]	NAP	C5A-C6A	2.59	1.48	1.41
2	F	285	NAP	C5A-C6A	2.57	1.48	1.41
2	E	285	NAP	C5A-C6A	2.56	1.48	1.41
2	F	285	NAP	PN-O3	2.53	1.62	1.59
2	D	285	NAP	PN-O3	2.53	1.62	1.59
2	B	285[B]	NAP	PN-O3	2.48	1.62	1.59
2	F	285	NAP	PA-O3	2.45	1.62	1.59
2	C	285	NAP	PN-O3	2.43	1.62	1.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	285	NAP	O4D-C1D	2.41	1.44	1.40
2	E	285	NAP	PA-O3	2.37	1.62	1.59
2	A	285	NAP	PA-O3	2.32	1.62	1.59
2	D	285	NAP	C5A-N7A	-2.31	1.34	1.39
2	B	285[A]	NAP	PA-O3	2.30	1.62	1.59
2	B	285[A]	NAP	O4D-C1D	2.29	1.43	1.40
2	D	285	NAP	C8A-N7A	2.28	1.36	1.31
2	B	285[A]	NAP	C8A-N7A	2.27	1.36	1.31
2	A	285	NAP	O4D-C1D	2.24	1.43	1.40
2	B	285[B]	NAP	PA-O3	2.23	1.61	1.59
2	E	285	NAP	C8A-N7A	2.22	1.35	1.31
2	A	285	NAP	C5A-N7A	-2.21	1.35	1.39
2	E	285	NAP	O4D-C1D	2.20	1.43	1.40
2	C	285	NAP	C5A-N7A	-2.20	1.35	1.39
2	B	285[B]	NAP	O4D-C1D	2.19	1.43	1.40
2	B	285[B]	NAP	C5A-N7A	-2.18	1.35	1.39
2	C	285	NAP	O4D-C1D	2.17	1.43	1.40
2	F	285	NAP	C8A-N7A	2.17	1.35	1.31
2	E	285	NAP	PN-O3	2.16	1.61	1.59
2	E	285	NAP	C5A-N7A	-2.12	1.35	1.39
2	C	285	NAP	C8A-N7A	2.12	1.35	1.31
2	A	285	NAP	C8A-N7A	2.10	1.35	1.31
2	B	285[B]	NAP	C8A-N7A	2.09	1.35	1.31
2	C	285	NAP	PA-O3	2.06	1.61	1.59
2	B	285[A]	NAP	C5A-N7A	-2.02	1.35	1.39
2	D	285	NAP	O4D-C1D	2.00	1.43	1.40

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	285[B]	NAP	C5A-C4A-N3A	-5.35	119.35	126.72
2	D	285	NAP	C5A-C4A-N3A	-5.19	119.57	126.72
2	A	285	NAP	C5A-C4A-N3A	-5.15	119.62	126.72
2	F	285	NAP	C5A-C4A-N3A	-5.04	119.77	126.72
2	C	285	NAP	C5A-C4A-N3A	-5.04	119.77	126.72
2	E	285	NAP	C5A-C4A-N3A	-4.89	119.98	126.72
2	B	285[A]	NAP	C5A-C4A-N3A	-4.79	120.13	126.72
2	E	285	NAP	N3A-C2A-N1A	-4.65	121.54	128.58
2	B	285[A]	NAP	N3A-C2A-N1A	-4.59	121.64	128.58
2	D	285	NAP	N3A-C2A-N1A	-4.58	121.64	128.58
2	B	285[B]	NAP	N3A-C2A-N1A	-4.47	121.81	128.58
2	A	285	NAP	N3A-C2A-N1A	-4.42	121.88	128.58

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	285	NAP	N3A-C2A-N1A	-4.35	122.00	128.58
2	F	285	NAP	N3A-C2A-N1A	-4.25	122.15	128.58
2	B	285[B]	NAP	C2A-N3A-C4A	4.05	121.72	111.83
2	D	285	NAP	C2A-N3A-C4A	4.05	121.72	111.83
2	D	285	NAP	O4B-C1B-N9A	4.03	115.82	108.09
2	B	285[A]	NAP	O4B-C1B-N9A	3.99	115.76	108.09
2	A	285	NAP	C2A-N3A-C4A	3.97	121.52	111.83
2	C	285	NAP	C2A-N3A-C4A	3.95	121.47	111.83
2	B	285[A]	NAP	C2A-N3A-C4A	3.94	121.46	111.83
2	E	285	NAP	C2A-N3A-C4A	3.93	121.43	111.83
2	B	285[A]	NAP	N9A-C8A-N7A	-3.89	108.41	113.94
2	E	285	NAP	N9A-C8A-N7A	-3.82	108.52	113.94
2	F	285	NAP	C2A-N3A-C4A	3.79	121.09	111.83
2	A	285	NAP	O4B-C1B-N9A	3.72	115.22	108.09
2	A	285	NAP	N9A-C8A-N7A	-3.69	108.70	113.94
2	B	285[B]	NAP	N3A-C4A-N9A	3.56	133.23	127.17
2	C	285	NAP	N9A-C8A-N7A	-3.53	108.92	113.94
2	B	285[B]	NAP	N9A-C8A-N7A	-3.52	108.94	113.94
2	D	285	NAP	N9A-C8A-N7A	-3.52	108.94	113.94
2	F	285	NAP	N9A-C8A-N7A	-3.43	109.06	113.94
2	F	285	NAP	O4B-C1B-N9A	3.41	114.64	108.09
2	A	285	NAP	N3A-C4A-N9A	3.29	132.76	127.17
2	F	285	NAP	N3A-C4A-N9A	3.25	132.70	127.17
2	D	285	NAP	N3A-C4A-N9A	3.23	132.66	127.17
2	C	285	NAP	O4B-C1B-N9A	3.21	114.26	108.09
2	E	285	NAP	O4B-C1B-N9A	3.17	114.17	108.09
2	E	285	NAP	N3A-C4A-N9A	3.16	132.55	127.17
2	C	285	NAP	N3A-C4A-N9A	3.10	132.45	127.17
2	B	285[A]	NAP	N3A-C4A-N9A	3.01	132.29	127.17
2	B	285[A]	NAP	C4A-N9A-C8A	2.94	108.82	105.74
2	E	285	NAP	C4A-N9A-C8A	2.87	108.75	105.74
2	B	285[B]	NAP	C2B-C1B-N9A	-2.83	109.10	113.75
2	A	285	NAP	C5A-N7A-C8A	2.64	107.60	103.45
2	B	285[B]	NAP	O4B-C1B-N9A	2.54	112.97	108.09
2	B	285[A]	NAP	C5A-N7A-C8A	2.53	107.43	103.45
2	A	285	NAP	C4A-C5A-N7A	-2.51	107.71	110.58
2	A	285	NAP	C4A-N9A-C8A	2.50	108.36	105.74
2	B	285[B]	NAP	C4A-N9A-C8A	2.50	108.36	105.74
2	B	285[A]	NAP	C4A-C5A-N7A	-2.47	107.76	110.58
2	B	285[B]	NAP	C5A-N7A-C8A	2.46	107.32	103.45
2	E	285	NAP	C5A-N7A-C8A	2.44	107.29	103.45
2	D	285	NAP	C5A-N7A-C8A	2.43	107.28	103.45

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	285	NAP	C5A-N7A-C8A	2.40	107.23	103.45
2	B	285[A]	NAP	C2B-C1B-N9A	-2.39	109.82	113.75
2	C	285	NAP	C5A-N7A-C8A	2.38	107.19	103.45
2	C	285	NAP	C4A-N9A-C8A	2.38	108.23	105.74
2	C	285	NAP	C4A-C5A-N7A	-2.37	107.87	110.58
2	D	285	NAP	C4A-C5A-N7A	-2.35	107.89	110.58
2	F	285	NAP	C4A-N9A-C8A	2.31	108.16	105.74
2	B	285[B]	NAP	C4A-C5A-N7A	-2.30	107.95	110.58
2	E	285	NAP	C4A-C5A-N7A	-2.29	107.96	110.58
2	D	285	NAP	C4A-N9A-C8A	2.26	108.11	105.74
2	F	285	NAP	C4A-C5A-N7A	-2.25	108.01	110.58
2	E	285	NAP	C3N-C7N-N7N	2.20	120.44	117.74
2	D	285	NAP	C3N-C7N-N7N	2.16	120.41	117.74
2	E	285	NAP	O2B-P2B-O1X	-2.01	102.16	109.33

There are no chirality outliers.

All (60) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	285	NAP	C5B-O5B-PA-O1A
2	A	285	NAP	C5B-O5B-PA-O3
2	A	285	NAP	PN-O3-PA-O5B
2	A	285	NAP	O4D-C1D-N1N-C2N
2	B	285[B]	NAP	C5B-O5B-PA-O1A
2	B	285[B]	NAP	C5B-O5B-PA-O2A
2	B	285[B]	NAP	C5D-O5D-PN-O3
2	B	285[B]	NAP	C5D-O5D-PN-O1N
2	B	285[B]	NAP	C5D-O5D-PN-O2N
2	C	285	NAP	C5B-O5B-PA-O1A
2	C	285	NAP	C5D-O5D-PN-O3
2	C	285	NAP	O4D-C1D-N1N-C2N
2	D	285	NAP	C5D-O5D-PN-O1N
2	D	285	NAP	O4D-C1D-N1N-C2N
2	E	285	NAP	C5D-O5D-PN-O3
2	E	285	NAP	C5D-O5D-PN-O1N
2	E	285	NAP	C5D-O5D-PN-O2N
2	F	285	NAP	C5D-O5D-PN-O3
2	F	285	NAP	C5D-O5D-PN-O1N
2	F	285	NAP	O4D-C1D-N1N-C2N
2	B	285[B]	NAP	C4B-C5B-O5B-PA
2	A	285	NAP	O4D-C4D-C5D-O5D
2	A	285	NAP	C3D-C4D-C5D-O5D

Continued on next page...

Continued from previous page...

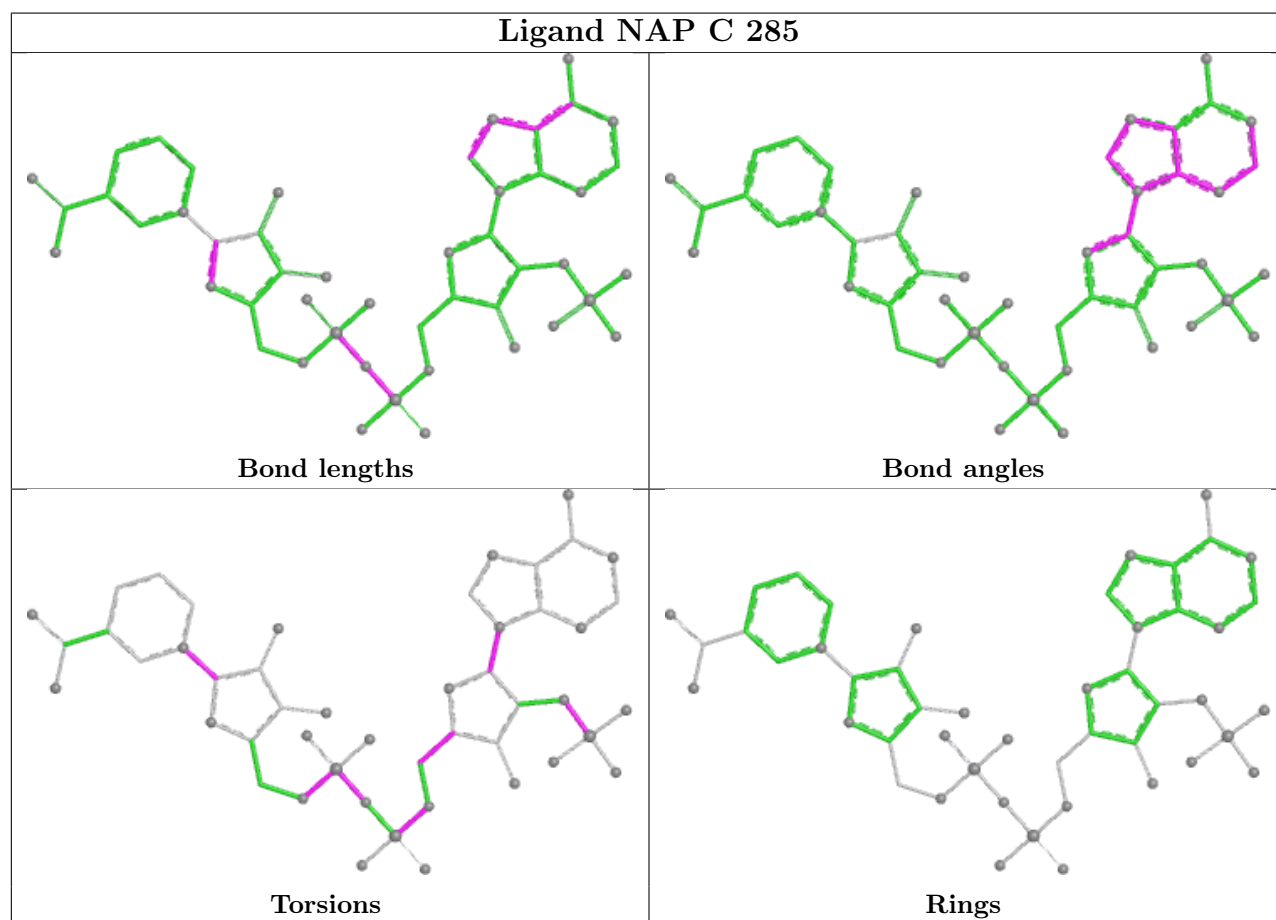
Mol	Chain	Res	Type	Atoms
2	B	285[B]	NAP	O4B-C4B-C5B-O5B
2	B	285[B]	NAP	C3B-C4B-C5B-O5B
2	B	285[A]	NAP	O4B-C4B-C5B-O5B
2	F	285	NAP	O4D-C4D-C5D-O5D
2	A	285	NAP	C2B-C1B-N9A-C8A
2	B	285[B]	NAP	C2B-C1B-N9A-C8A
2	E	285	NAP	C2B-C1B-N9A-C8A
2	F	285	NAP	C2B-C1B-N9A-C8A
2	F	285	NAP	C3D-C4D-C5D-O5D
2	A	285	NAP	C5B-O5B-PA-O2A
2	B	285[B]	NAP	C5B-O5B-PA-O3
2	C	285	NAP	C5B-O5B-PA-O3
2	C	285	NAP	C5D-O5D-PN-O1N
2	D	285	NAP	C5D-O5D-PN-O3
2	C	285	NAP	PA-O3-PN-O2N
2	C	285	NAP	C2B-C1B-N9A-C8A
2	C	285	NAP	C2B-O2B-P2B-O3X
2	A	285	NAP	O4D-C1D-N1N-C6N
2	C	285	NAP	O4D-C1D-N1N-C6N
2	D	285	NAP	O4D-C1D-N1N-C6N
2	F	285	NAP	O4D-C1D-N1N-C6N
2	C	285	NAP	C2B-O2B-P2B-O1X
2	A	285	NAP	PA-O3-PN-O1N
2	A	285	NAP	PA-O3-PN-O2N
2	D	285	NAP	PA-O3-PN-O2N
2	A	285	NAP	O4B-C1B-N9A-C8A
2	B	285[B]	NAP	O4B-C1B-N9A-C8A
2	E	285	NAP	O4B-C1B-N9A-C8A
2	B	285[A]	NAP	C2B-C1B-N9A-C8A
2	C	285	NAP	O4B-C4B-C5B-O5B
2	B	285[A]	NAP	PA-O3-PN-O2N
2	C	285	NAP	PA-O3-PN-O1N
2	F	285	NAP	PA-O3-PN-O2N
2	D	285	NAP	C2B-C1B-N9A-C8A
2	C	285	NAP	O4B-C1B-N9A-C8A
2	F	285	NAP	O4B-C1B-N9A-C8A
2	F	285	NAP	PA-O3-PN-O1N

There are no ring outliers.

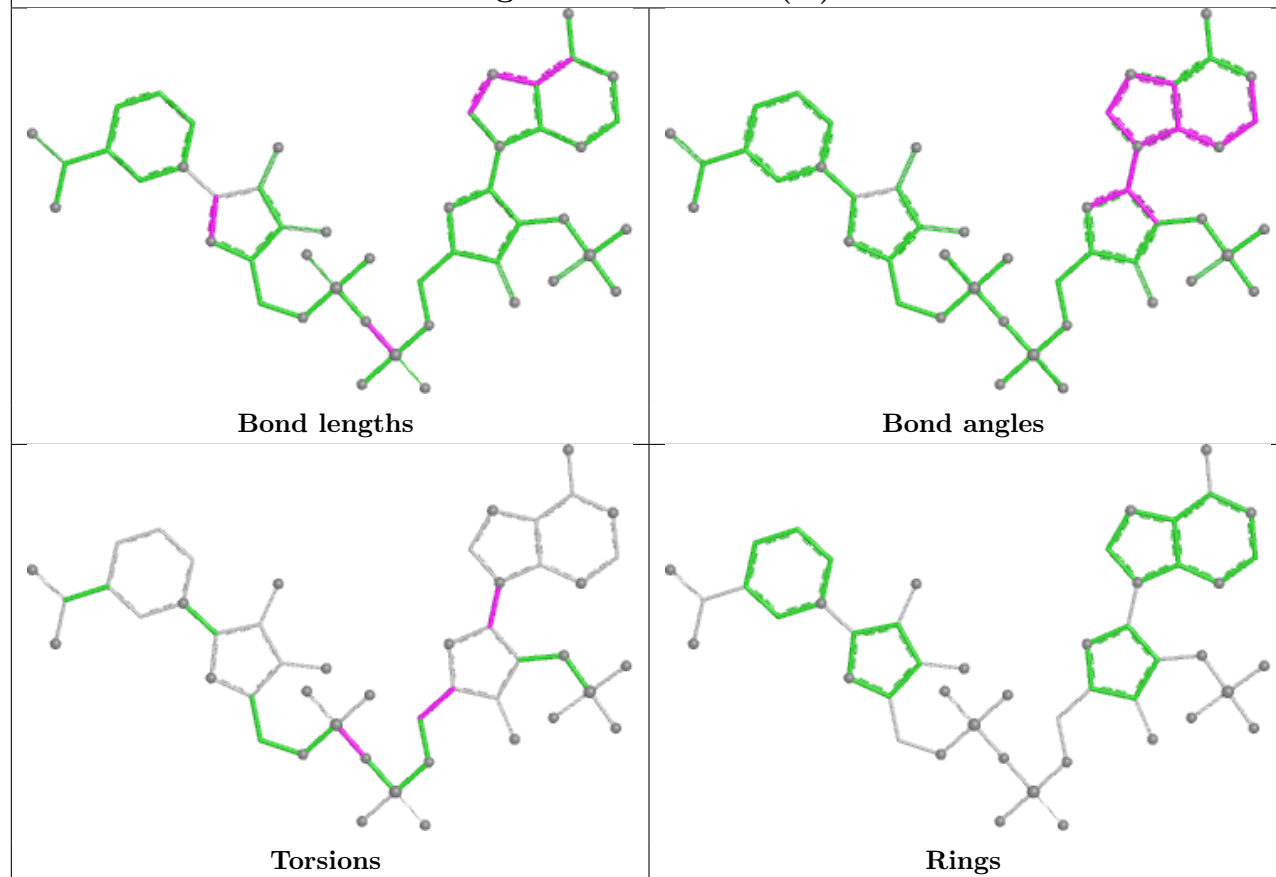
6 monomers are involved in 12 short contacts:

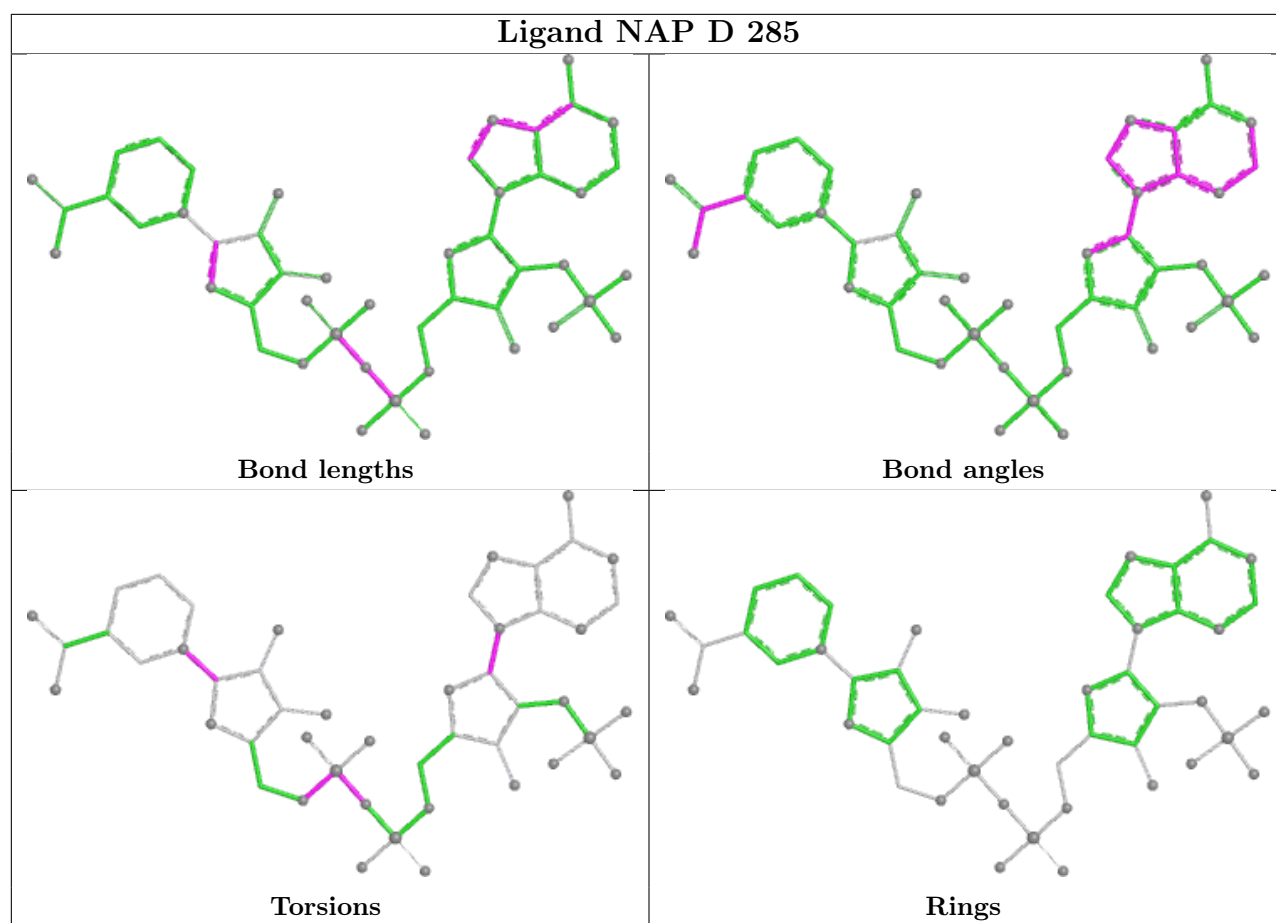
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	285	NAP	2	0
3	D	287	SO4	1	0
2	B	285[A]	NAP	2	0
2	B	285[B]	NAP	3	0
2	F	285	NAP	1	0
2	A	285	NAP	3	0

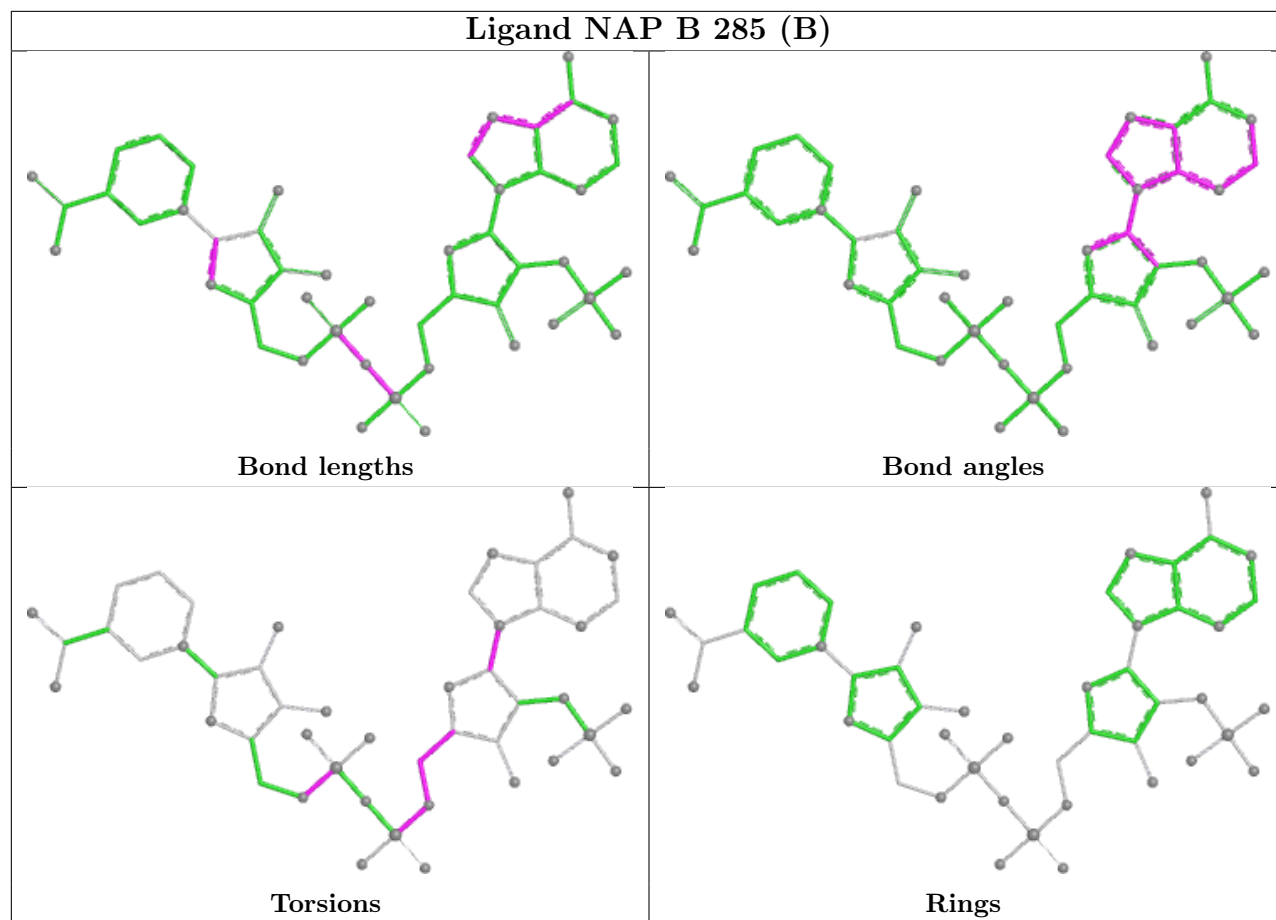
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

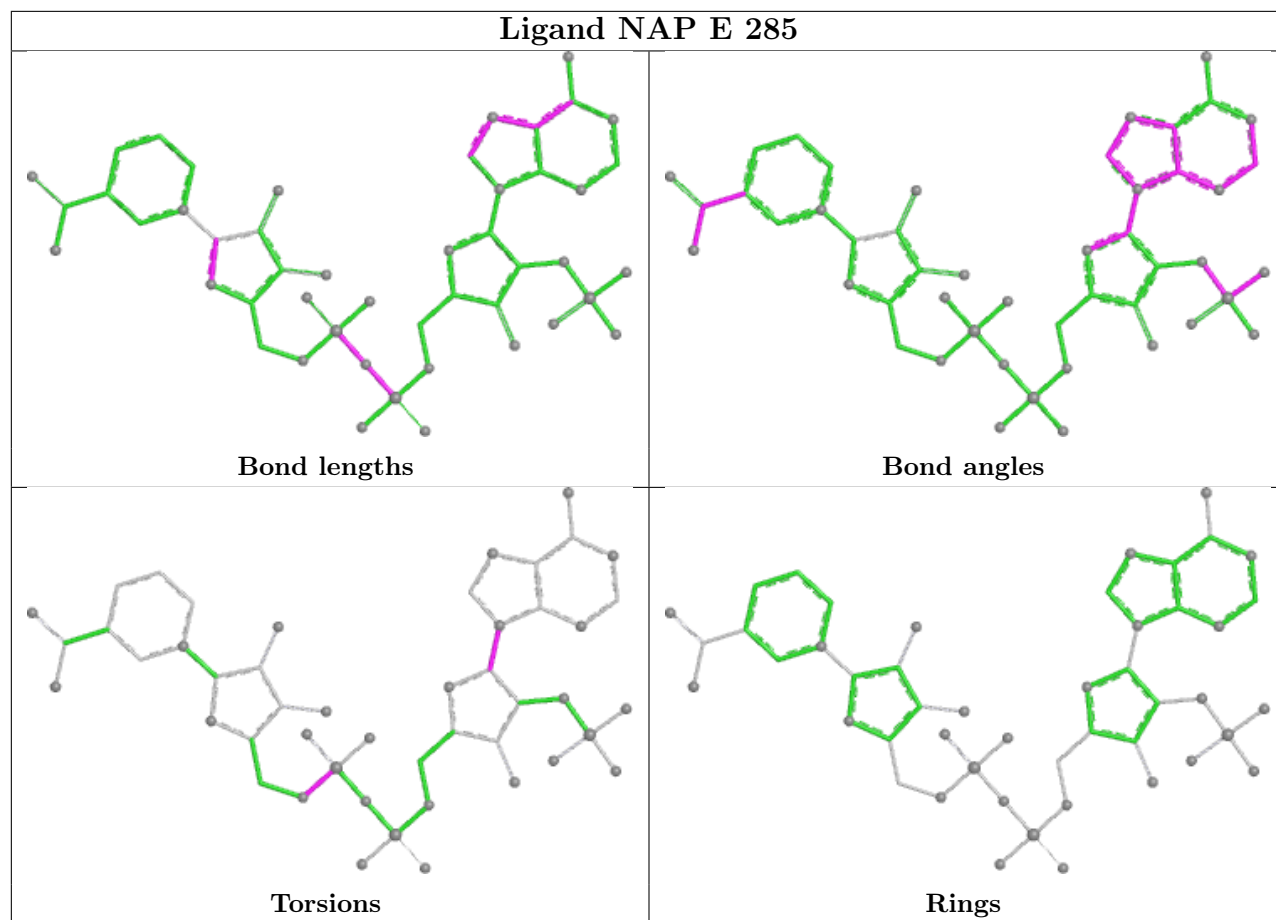


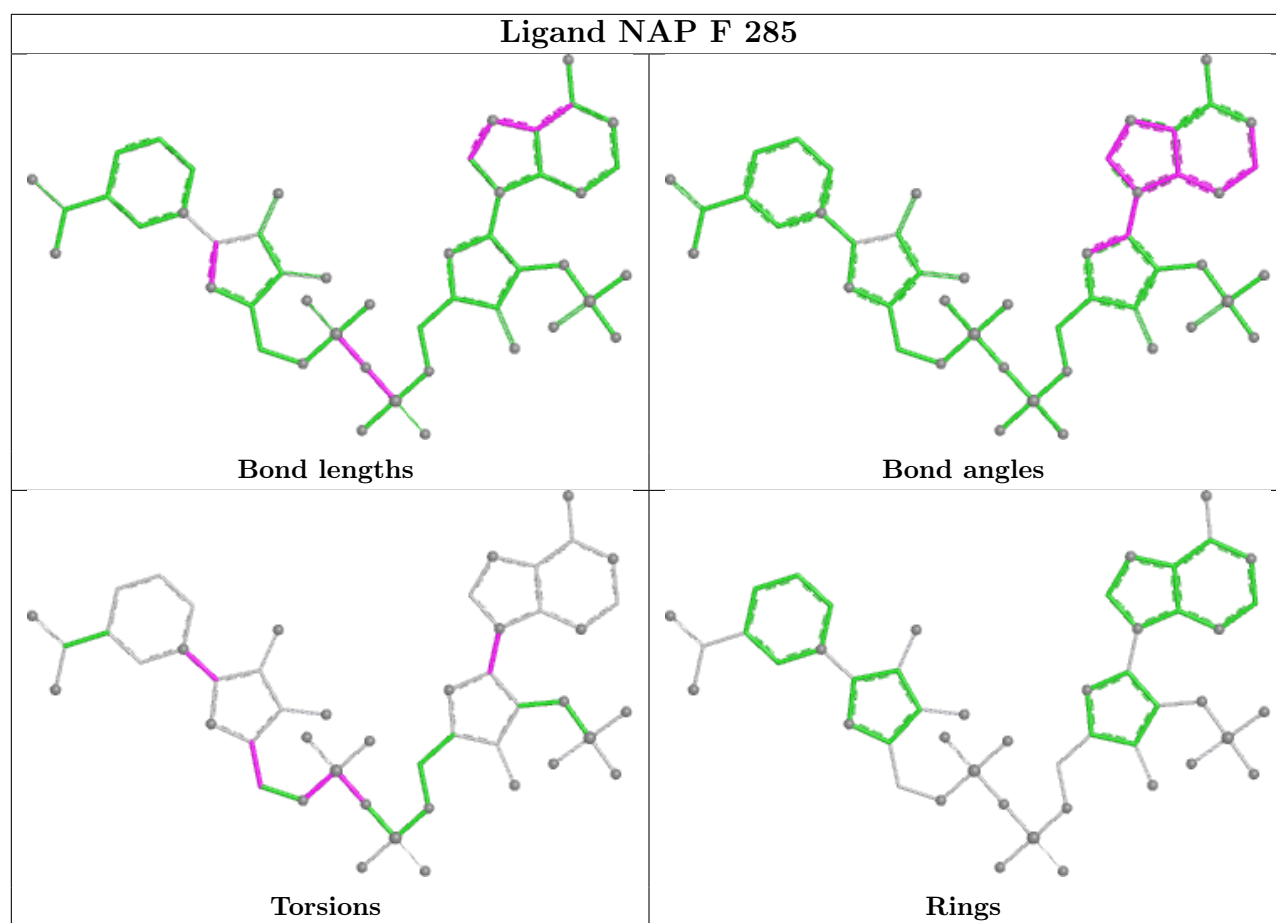
Ligand NAP B 285 (A)

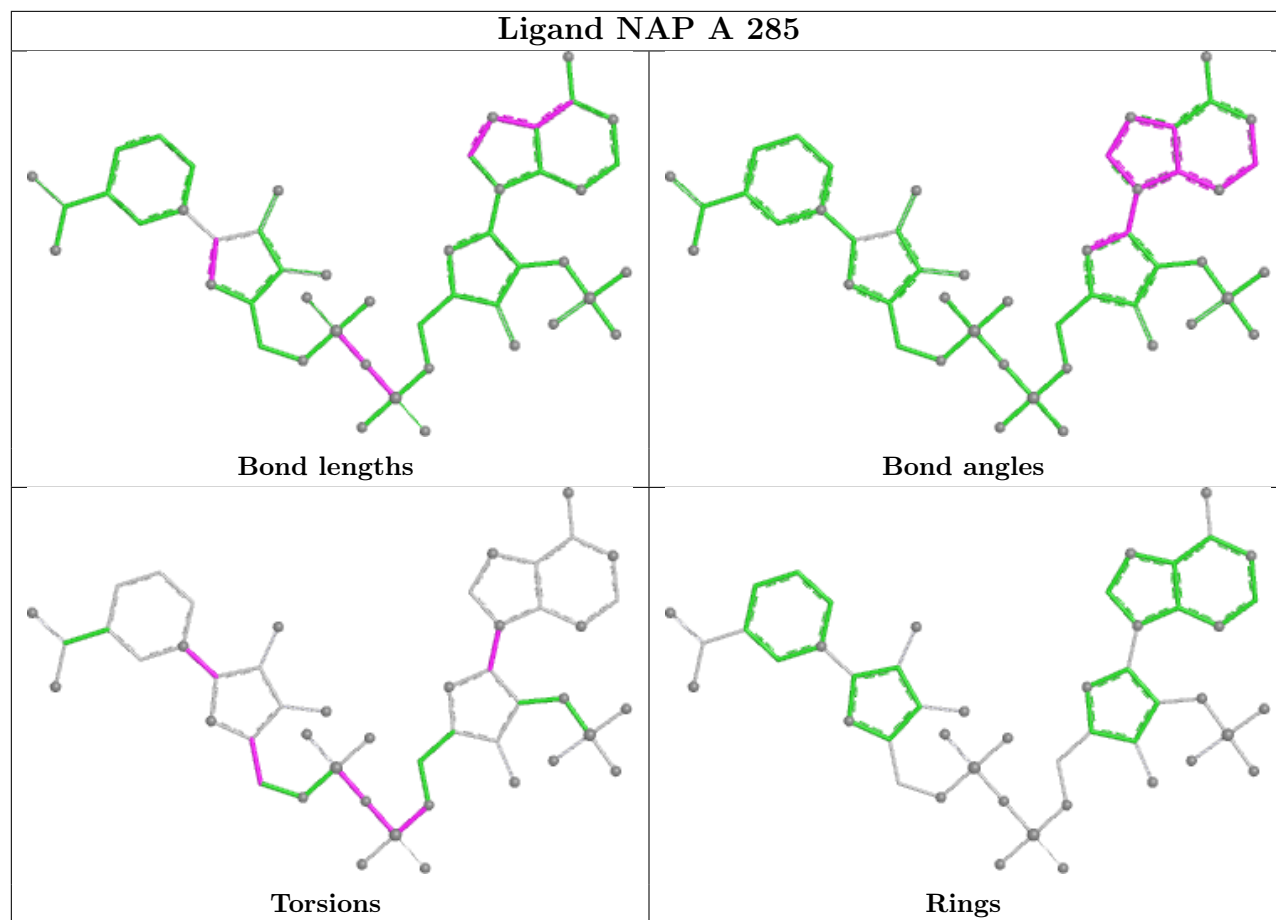












5.7 Other polymers [i](#)

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	281/287 (97%)	0.28	7 (2%) 58 52	21, 63, 89, 109	3 (1%)
1	B	280/287 (97%)	0.04	2 (0%) 84 82	26, 49, 67, 75	1 (0%)
1	C	281/287 (97%)	0.07	3 (1%) 78 75	28, 53, 71, 85	3 (1%)
1	D	283/287 (98%)	0.24	6 (2%) 63 57	26, 59, 83, 98	2 (0%)
1	E	278/287 (96%)	0.20	5 (1%) 67 63	32, 57, 80, 95	1 (0%)
1	F	281/287 (97%)	0.10	3 (1%) 78 75	30, 51, 69, 89	1 (0%)
All	All	1684/1722 (97%)	0.15	26 (1%) 72 67	21, 55, 80, 109	11 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	245	ALA	3.6
1	D	-1	ASN	3.4
1	E	119[A]	HIS	3.4
1	E	170	LYS	3.3
1	F	119[A]	HIS	3.2
1	A	119[A]	HIS	3.1
1	D	119[A]	HIS	3.0
1	E	282	LYS	2.8
1	C	172[A]	ARG	2.7
1	D	244	ALA	2.6
1	C	265	LEU	2.5
1	D	14	LEU	2.5
1	B	119[A]	HIS	2.4
1	C	240	GLU	2.4
1	E	15	GLY	2.4
1	F	172	ARG	2.2
1	A	244	ALA	2.2
1	A	14	LEU	2.2
1	A	247	ALA	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	281	THR	2.2
1	B	245	ALA	2.1
1	D	170	LYS	2.1
1	A	210	SER	2.1
1	F	14	LEU	2.1
1	E	242	PHE	2.1
1	A	243	GLY	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
3	SO4	F	288	5/5	0.22	0.17	104,104,105,105	5
3	SO4	B	289	5/5	0.45	0.18	95,95,95,95	5
3	SO4	A	287	5/5	0.50	0.12	100,100,100,100	5
3	SO4	F	286	5/5	0.55	0.13	77,78,79,79	5
3	SO4	E	287	5/5	0.58	0.13	90,90,91,91	5
3	SO4	E	288	5/5	0.62	0.15	89,89,90,90	5
3	SO4	F	287	5/5	0.71	0.13	78,79,79,80	5
3	SO4	E	289	5/5	0.71	0.13	73,74,75,75	5
3	SO4	D	286	5/5	0.78	0.11	78,79,79,80	5
3	SO4	C	287	5/5	0.80	0.10	67,69,69,70	5
3	SO4	E	286	5/5	0.80	0.09	73,73,75,75	5
3	SO4	B	288	5/5	0.81	0.13	73,73,75,75	5
3	SO4	D	287	5/5	0.83	0.13	67,69,70,71	5
3	SO4	A	286	5/5	0.87	0.08	80,80,80,81	5
3	SO4	B	286	5/5	0.90	0.18	69,71,72,74	0
3	SO4	B	287	5/5	0.90	0.07	69,71,72,72	5

Continued on next page...

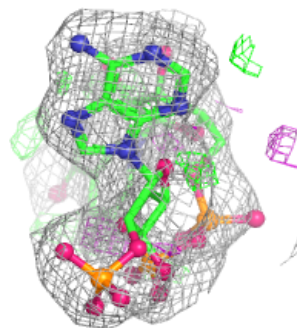
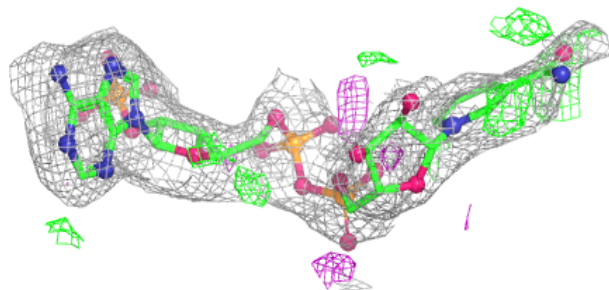
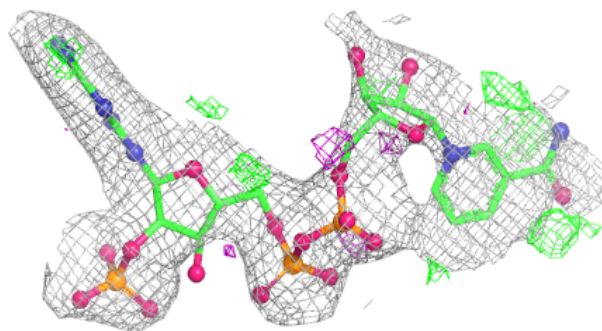
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	C	286	5/5	0.91	0.11	52,54,55,56	5
2	NAP	B	285[B]	48/48	0.93	0.11	65,68,82,83	48
2	NAP	B	285[A]	48/48	0.93	0.11	23,36,53,56	48
2	NAP	D	285	48/48	0.94	0.09	49,54,68,70	0
2	NAP	A	285	48/48	0.94	0.09	48,59,76,77	0
2	NAP	F	285	48/48	0.95	0.08	47,51,66,68	0
2	NAP	C	285	48/48	0.96	0.07	42,49,61,63	0
2	NAP	E	285	48/48	0.96	0.08	46,55,70,72	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.

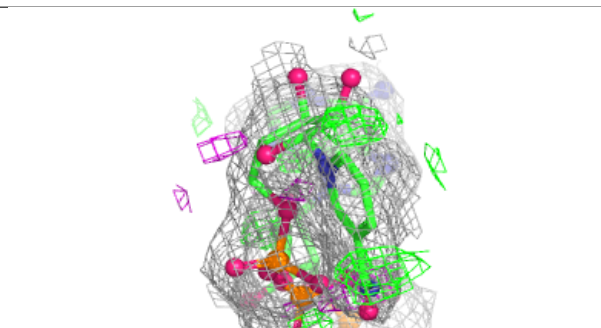
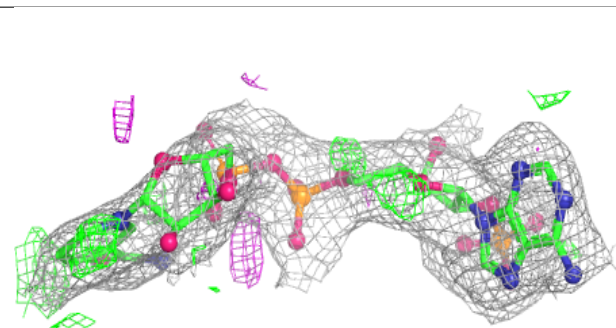
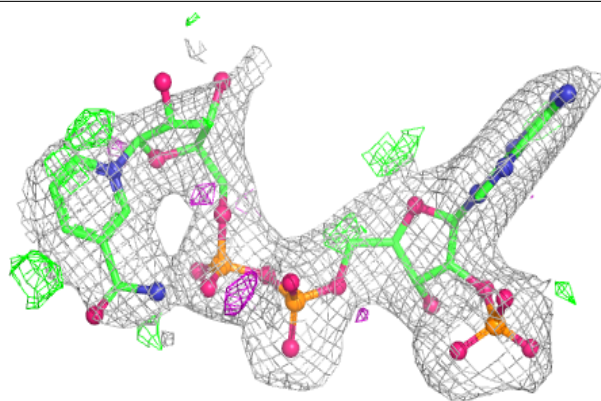
Electron density around NAP B 285 (B):

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

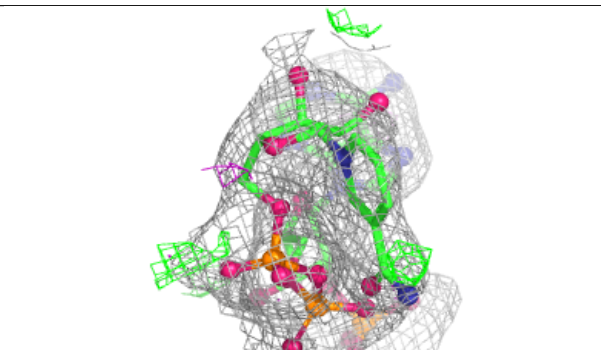
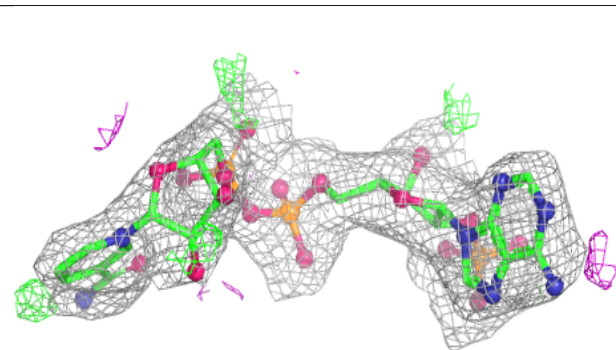
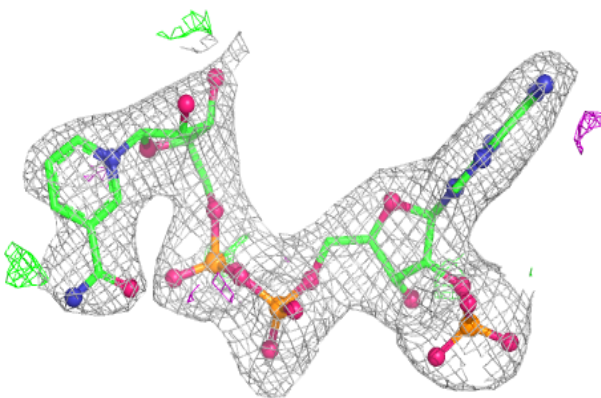


Electron density around NAP B 285 (A):

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

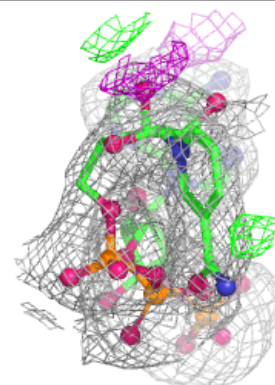
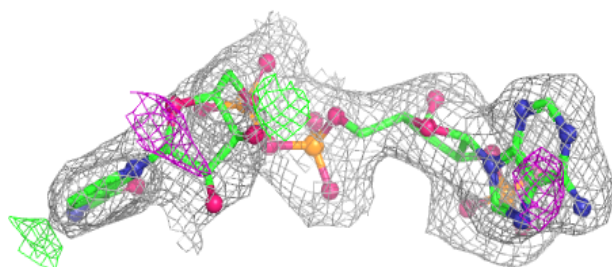
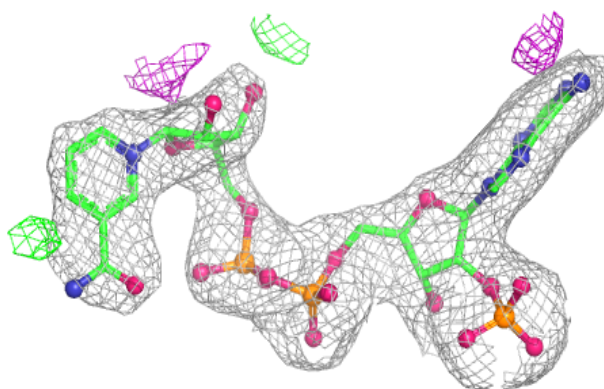
**Electron density around NAP D 285:**

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

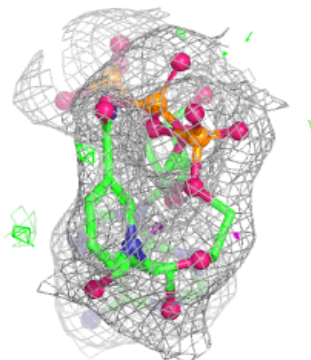
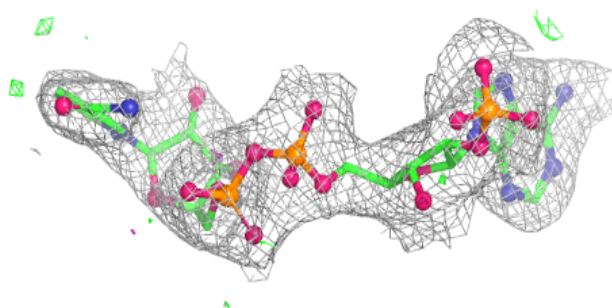
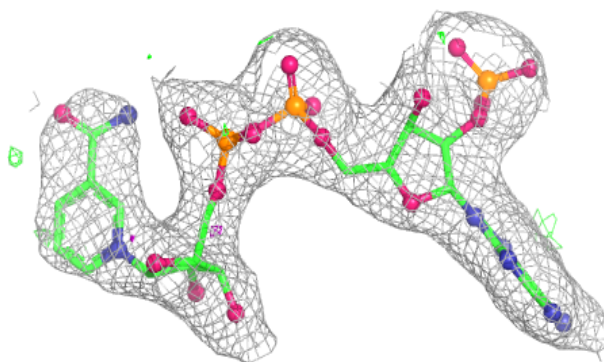


Electron density around NAP A 285:

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

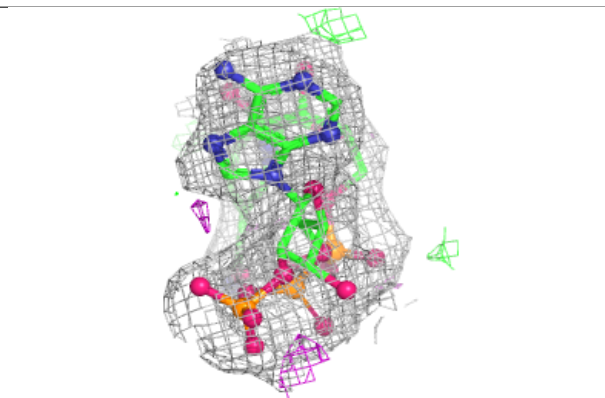
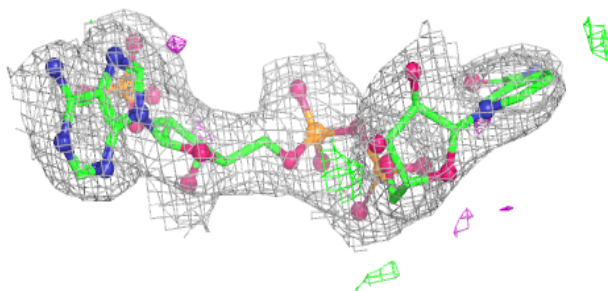
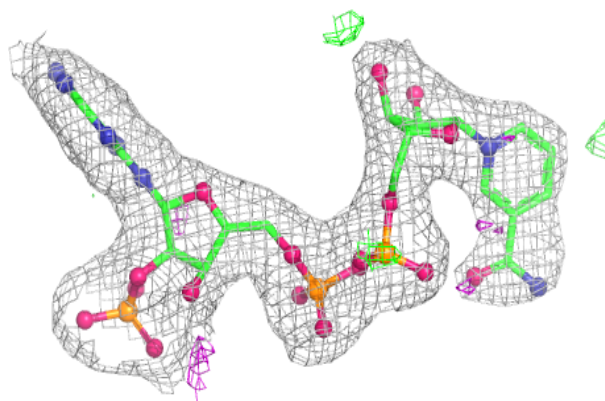
**Electron density around NAP F 285:**

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

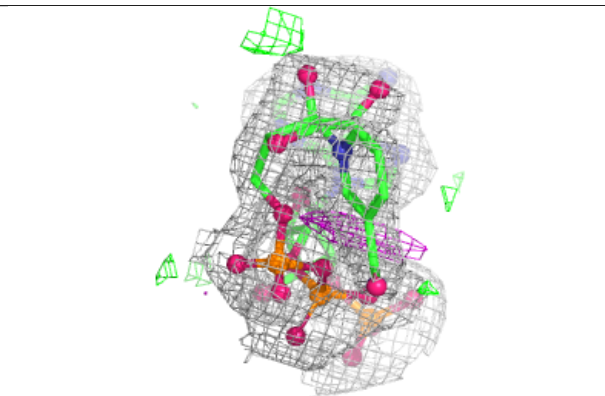
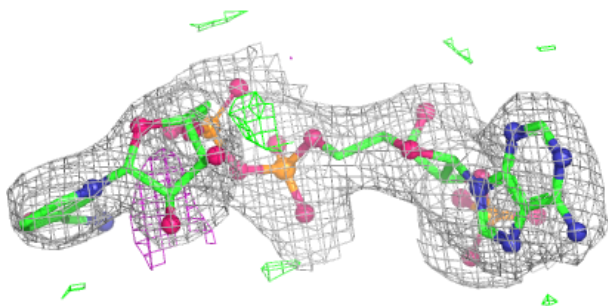
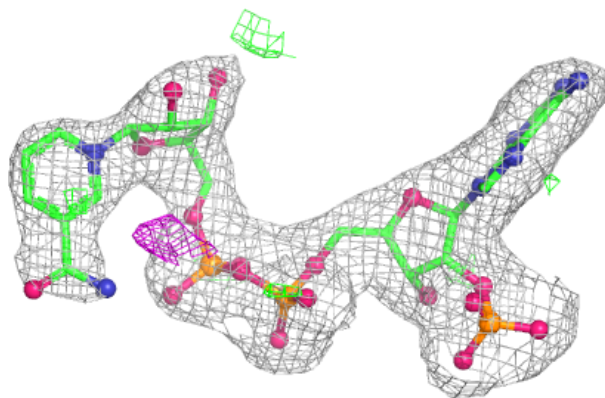


Electron density around NAP C 285:

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

**Electron density around NAP E 285:**

$2mF_o-DF_c$ (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.