



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 02:04 pm BST

PDB ID : 3QV1
Title : Crystal structure of the binary complex of photosynthetic A4 glyceraldehyde 3-phosphate dehydrogenase (GAPDH) with cp12-2, both from *Arabidopsis thaliana*.
Authors : Thumiger, A.; Fermani, S.; Falini, G.; Marri, L.; Sparla, F.; Trost, P.
Deposited on : 2011-02-25
Resolution : 2.00 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

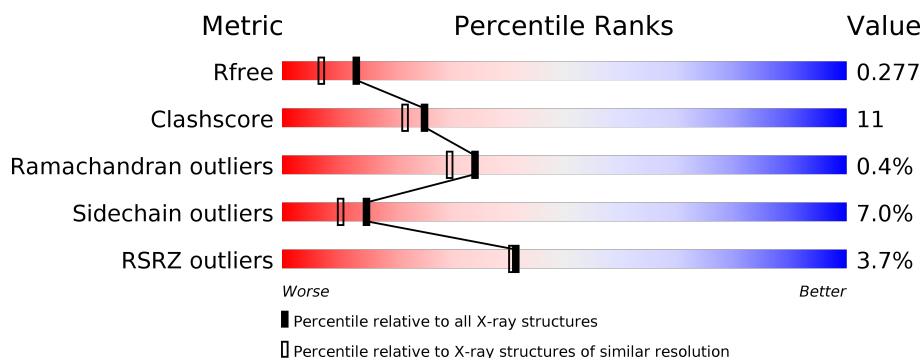
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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





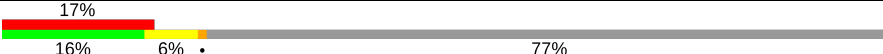
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on 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	337	
1	B	337	
1	C	337	
1	D	337	
1	E	337	
1	F	337	

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Mol	Chain	Length	Quality of chain
2	G	82	
2	H	82	
2	I	82	

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
4	PEG	A	334	-	-	X	-
4	PEG	A	336	-	-	X	-
4	PEG	C	334	-	-	X	-
6	EDO	C	336	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 16337 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 Glyceraldehyde-3-phosphate dehydrogenase A, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	1	0
			2562	1620	445	488	9			
1	B	336	Total	C	N	O	S	0	1	0
			2557	1616	444	487	10			
1	C	336	Total	C	N	O	S	0	0	0
			2552	1612	444	487	9			
1	D	337	Total	C	N	O	S	0	1	0
			2562	1620	445	488	9			
1	E	336	Total	C	N	O	S	0	0	0
			2552	1612	444	487	9			
1	F	336	Total	C	N	O	S	0	1	0
			2557	1616	444	487	10			

- Molecule 2 is a protein called CP12 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	21	Total	C	N	O	S	0	0	0
			177	104	28	43	2			
2	H	20	Total	C	N	O	S	0	0	0
			168	100	26	40	2			
2	I	19	Total	C	N	O	S	0	0	0
			160	96	25	37	2			

There are 12 discrepancies between the modelled and reference sequences:

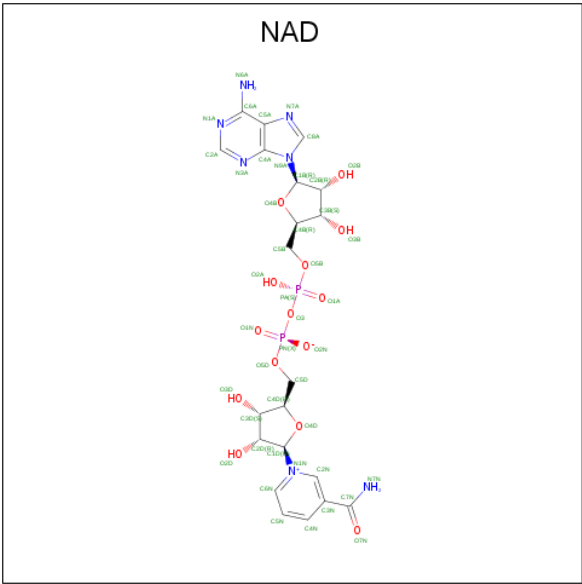
Chain	Residue	Modelled	Actual	Comment	Reference
G	-3	GLY	-	EXPRESSION TAG	UNP Q9LZP9
G	-2	SER	-	EXPRESSION TAG	UNP Q9LZP9
G	-1	HIS	-	EXPRESSION TAG	UNP Q9LZP9
G	0	MET	-	EXPRESSION TAG	UNP Q9LZP9
H	-3	GLY	-	EXPRESSION TAG	UNP Q9LZP9
H	-2	SER	-	EXPRESSION TAG	UNP Q9LZP9
H	-1	HIS	-	EXPRESSION TAG	UNP Q9LZP9

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Chain	Residue	Modelled	Actual	Comment	Reference
H	0	MET	-	EXPRESSION TAG	UNP Q9LZP9
I	-3	GLY	-	EXPRESSION TAG	UNP Q9LZP9
I	-2	SER	-	EXPRESSION TAG	UNP Q9LZP9
I	-1	HIS	-	EXPRESSION TAG	UNP Q9LZP9
I	0	MET	-	EXPRESSION TAG	UNP Q9LZP9

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



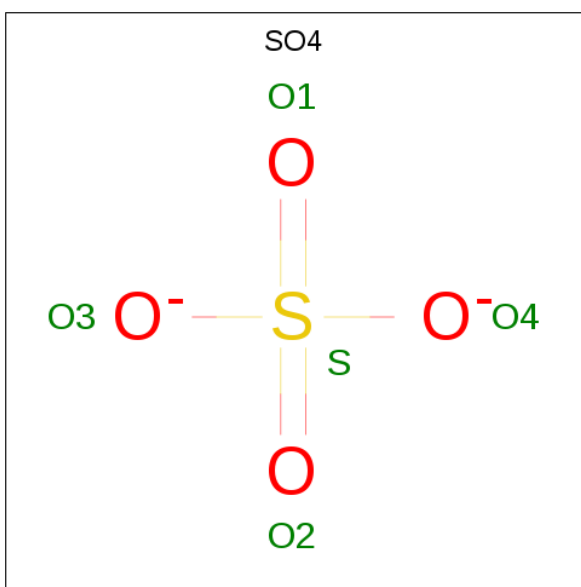
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

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

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			4	2	2		

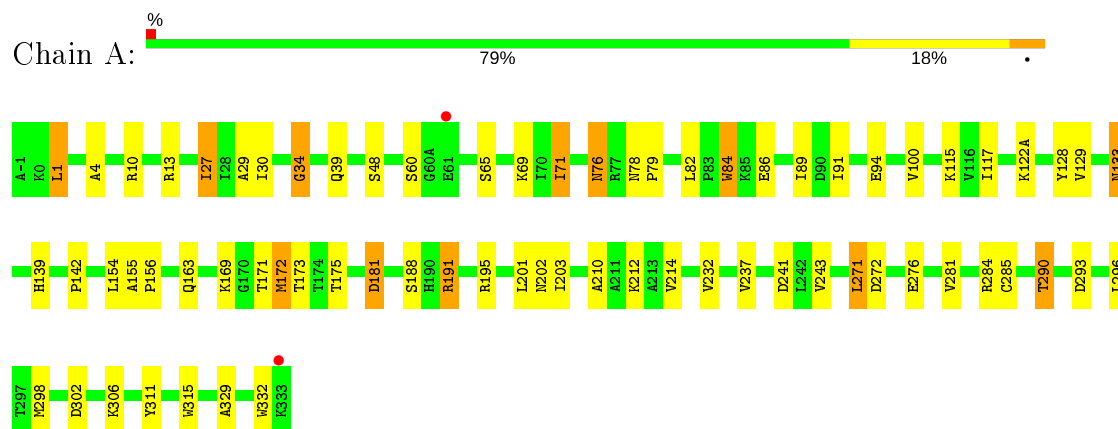
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	57	Total	O	0	0
			57	57		
7	B	33	Total	O	0	0
			33	33		
7	C	31	Total	O	0	0
			31	31		
7	D	38	Total	O	0	0
			38	38		
7	E	17	Total	O	0	0
			17	17		
7	F	14	Total	O	0	0
			14	14		
7	G	1	Total	O	0	0
			1	1		

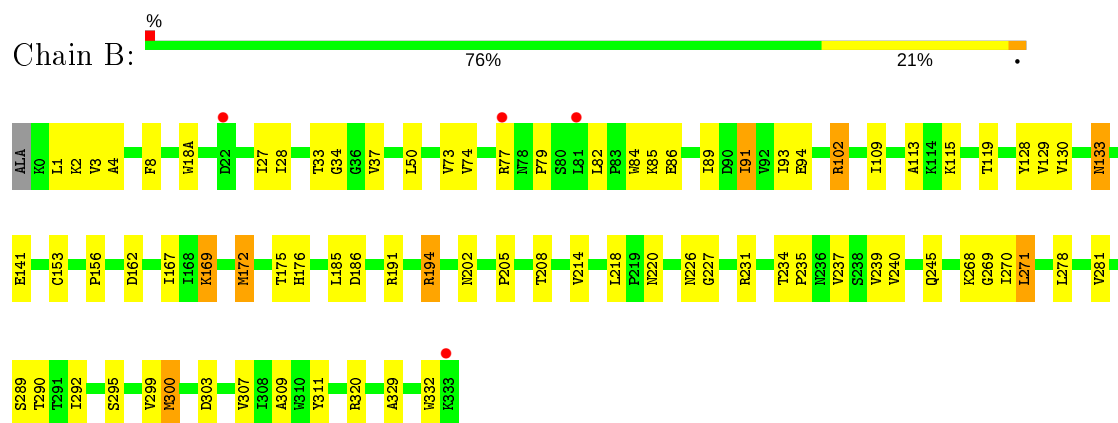
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

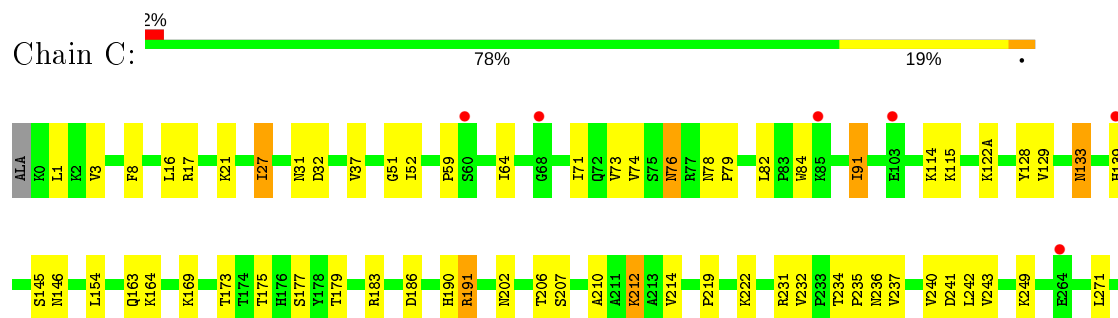
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A, chloroplastic



- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A, chloroplastic

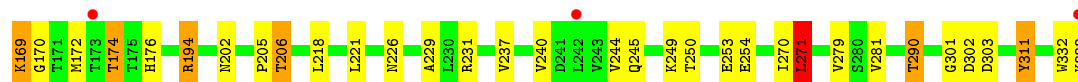
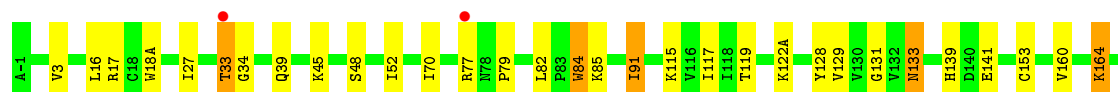
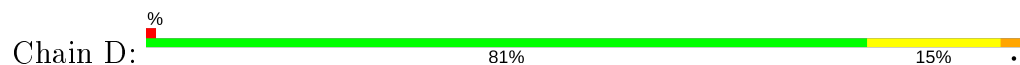


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A, chloroplastic

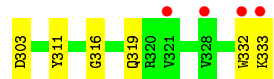
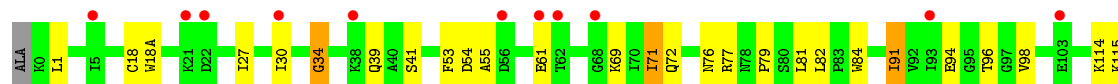
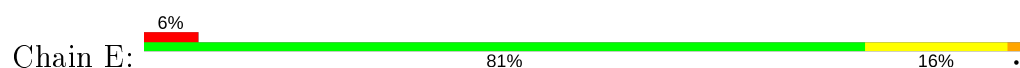




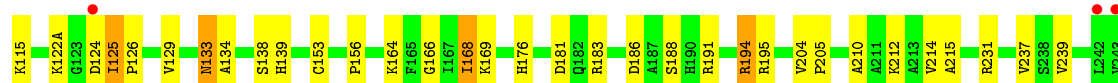
- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A, chloroplatic



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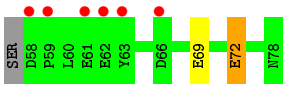


- Molecule 1: Glyceraldehyde-3-phosphate dehydrogenase A, chloroplatic



- Molecule 2: CP12 protein

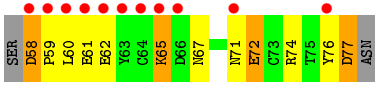




● Molecule 2: CP12 protein



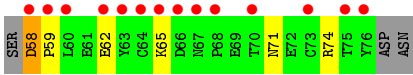
GLY	SER	HIS	MET	ALA	ALA	PRO	GLU	GLY	GLY	ILE	SER	ASP	VAL	VAL	GLU	LYS	SER	ILE	LYS	GLU	ALA	GLN	GLU	THR	CYS	ALA	GLY	ASP	PRO	VAL	SER	GLY	GLU	CYS	VAL	ALA	ALA	TRP	ASP	GLU	VAL	VAL	GLU	GLU	LEU	SER	ALA	ALA	ALA	SER	HIS	ALA	ARG	ASP	LYS	LYS	LYS	ALA	ASP	GLY
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● Molecule 2: CP12 protein



GLY	SER	HIS	MET	ALA	ALA	PRO	GLU	GLY	GLY	ILE	SER	ASP	VAL	VAL	GLU	LYS	SER	ILE	LYS	GLU	ALA	GLN	GLU	THR	CYS	ALA	GLY	ASP	PRO	VAL	SER	GLY	GLU	CYS	VAL	ALA	ALA	TRP	ASP	GLU	VAL	VAL	GLU	GLU	LEU	SER	ALA	ALA	ALA	SER	HIS	ALA	ARG	ASP	LYS	LYS	LYS	ALA	ASP	GLY
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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	142.67Å 245.99Å 138.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.20 – 2.00 34.26 – 1.98	Depositor EDS
% Data completeness (in resolution range)	99.0 (34.20-2.00) 99.1 (34.26-1.98)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.225 , 0.272 0.233 , 0.277	Depositor DCC
R_{free} test set	8373 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.672	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.077 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.078 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16337	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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: PEG, SO4, NAD, EDO

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	0.86	0/2609	0.91	5/3541 (0.1%)
1	B	0.69	0/2604	0.82	3/3533 (0.1%)
1	C	0.80	0/2596	0.88	4/3523 (0.1%)
1	D	0.79	0/2609	0.89	5/3541 (0.1%)
1	E	0.68	0/2596	0.81	2/3523 (0.1%)
1	F	0.65	0/2604	0.79	4/3533 (0.1%)
2	G	0.73	0/180	0.79	0/243
2	H	0.88	0/171	0.79	0/232
2	I	0.72	0/163	0.61	0/221
All	All	0.75	0/16132	0.85	23/21890 (0.1%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	17	ARG	NE-CZ-NH2	-8.35	116.13	120.30
1	C	186	ASP	CB-CG-OD2	-7.71	111.36	118.30
1	C	186	ASP	CB-CG-OD1	6.92	124.53	118.30
1	D	17	ARG	NE-CZ-NH1	6.75	123.67	120.30
1	D	194	ARG	NE-CZ-NH1	6.22	123.41	120.30
1	A	34	GLY	N-CA-C	6.22	128.64	113.10
1	B	271	LEU	CA-CB-CG	6.13	129.40	115.30
1	C	17	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	D	271	LEU	CA-CB-CG	5.91	128.90	115.30
1	F	186	ASP	CB-CG-OD2	-5.87	113.02	118.30
1	C	17	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	E	194	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	B	194	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	A	191	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	A	298	MET	CG-SD-CE	-5.34	91.66	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	271	LEU	CA-CB-CG	5.27	127.42	115.30
1	A	284	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	F	194	ARG	NE-CZ-NH1	5.13	122.87	120.30
1	F	186	ASP	CB-CG-OD1	5.11	122.90	118.30
1	A	302	ASP	CB-CG-OD1	5.08	122.88	118.30
1	B	102	ARG	NE-CZ-NH1	-5.06	117.77	120.30
1	D	34	GLY	N-CA-C	5.05	125.72	113.10
1	E	34	GLY	N-CA-C	5.04	125.69	113.10

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	2562	0	2609	63	0
1	B	2557	0	2602	68	0
1	C	2552	0	2593	74	0
1	D	2562	0	2609	61	0
1	E	2552	0	2593	34	1
1	F	2557	0	2602	59	0
2	G	177	0	144	2	0
2	H	168	0	138	7	0
2	I	160	0	134	7	0
3	A	44	0	26	0	0
3	B	44	0	26	1	0
3	C	44	0	26	0	0
3	D	44	0	26	1	0
3	E	44	0	26	0	0
3	F	44	0	26	1	0
4	A	14	0	20	19	0
4	C	7	0	10	15	0
5	A	10	0	0	0	0
6	C	4	0	6	7	0
7	A	57	0	0	1	0
7	B	33	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	31	0	0	0	0
7	D	38	0	0	0	0
7	E	17	0	0	0	0
7	F	14	0	0	0	0
7	G	1	0	0	0	0
All	All	16337	0	16216	345	1

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 11.

All (345) 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:203[A]:ILE:HD12	1:A:232:VAL:HG12	1.45	0.99
1:C:76:ASN:HD22	1:C:78:ASN:H	1.08	0.95
1:C:241:ASP:HB2	4:C:334:PEG:H32	1.58	0.84
1:B:102:ARG:HG2	1:B:102:ARG:HH11	1.45	0.81
1:B:34:GLY:O	2:H:60:LEU:HB2	1.83	0.79
1:B:115:LYS:NZ	1:B:141:GLU:O	2.15	0.78
1:B:194:ARG:NH1	1:B:205:PRO:HG2	1.99	0.78
1:B:33:THR:HG21	1:B:77:ARG:HH11	1.52	0.75
1:A:30:ILE:HD12	1:A:71:ILE:HD11	1.68	0.75
1:A:172:MET:HA	4:A:334:PEG:H11	1.67	0.75
2:I:58:ASP:HB3	2:I:59:PRO:HD3	1.70	0.74
1:F:129:VAL:H	1:F:133:ASN:HD21	1.35	0.74
1:A:171:THR:HG23	4:A:334:PEG:O1	1.85	0.74
1:F:194:ARG:HD2	1:F:205:PRO:O	1.88	0.73
1:C:241:ASP:CB	4:C:334:PEG:C3	2.69	0.71
1:A:156:PRO:HB2	1:A:290:THR:HG21	1.73	0.70
1:B:2:LYS:HG2	1:B:28:ILE:HD13	1.71	0.70
1:C:128:TYR:HA	1:C:133:ASN:HD21	1.57	0.70
1:F:38:LYS:HD2	1:F:38:LYS:H	1.56	0.70
1:B:202:ASN:HD21	1:D:281:VAL:HG12	1.55	0.69
1:E:176:HIS:HB3	1:E:231:ARG:HD3	1.73	0.69
1:C:241:ASP:HB2	4:C:334:PEG:C3	2.22	0.69
1:C:76:ASN:ND2	1:C:78:ASN:H	1.86	0.69
1:C:129:VAL:H	1:C:133:ASN:ND2	1.92	0.69
1:A:281:VAL:HB	1:C:202:ASN:ND2	2.08	0.69
1:F:168:ILE:HD11	1:F:247:SER:HB3	1.75	0.68
1:F:293:ASP:CG	1:F:296:LEU:HD12	2.13	0.68
1:C:241:ASP:CB	4:C:334:PEG:H32	2.22	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:334:PEG:O1	1:C:306:LYS:NZ	2.26	0.68
1:B:202:ASN:HD21	1:D:281:VAL:CG1	2.06	0.67
1:C:241:ASP:HB3	4:C:334:PEG:H31	1.75	0.67
1:D:33:THR:CG2	1:D:77:ARG:NH1	2.58	0.67
1:F:298:MET:HB3	1:F:300[A]:MET:CE	2.25	0.67
1:A:201:LEU:HD22	4:A:336:PEG:H31	1.76	0.67
1:B:172:MET:HE1	1:B:208:THR:HG21	1.78	0.65
1:C:139:HIS:HB3	1:C:333:LYS:NZ	2.11	0.65
1:D:271:LEU:HD22	1:D:290:THR:OG1	1.96	0.65
2:I:58:ASP:CB	2:I:59:PRO:HD3	2.26	0.65
1:B:129:VAL:H	1:B:133:ASN:HD21	1.45	0.65
1:B:194:ARG:HH11	1:B:205:PRO:HG2	1.59	0.65
1:F:293:ASP:CB	1:F:296:LEU:HD12	2.27	0.64
1:A:129:VAL:H	1:A:133:ASN:ND2	1.95	0.64
1:D:131:GLY:HA3	1:D:270:ILE:HD13	1.77	0.64
1:A:175:THR:OG1	4:C:334:PEG:H11	1.98	0.64
1:C:129:VAL:H	1:C:133:ASN:HD21	1.46	0.64
1:D:115:LYS:NZ	1:D:141:GLU:O	2.26	0.64
1:F:89:ILE:O	1:F:113:ALA:HA	1.97	0.64
1:F:293:ASP:HB3	1:F:296:LEU:HD12	1.78	0.63
1:C:241:ASP:CB	4:C:334:PEG:H31	2.29	0.63
1:F:210:ALA:O	1:F:214:VAL:HG23	1.99	0.63
1:E:210:ALA:O	1:E:214:VAL:HG23	1.98	0.63
1:A:76:ASN:HD22	1:A:78:ASN:H	1.44	0.62
1:F:102:ARG:HG2	1:F:102:ARG:HH11	1.64	0.61
1:B:115:LYS:HD2	1:B:332:TRP:HZ3	1.65	0.61
2:I:62:GLU:OE1	2:I:65:LYS:NZ	2.33	0.61
1:A:188:SER:HB2	1:D:39:GLN:OE1	1.99	0.61
1:C:281:VAL:HG11	1:D:48:SER:HA	1.82	0.61
1:B:33:THR:HG22	1:B:77:ARG:HD2	1.82	0.61
1:B:300[A]:MET:CE	1:D:226:ASN:HD22	2.13	0.61
1:B:28:ILE:HD11	1:B:89:ILE:HD12	1.82	0.60
1:C:154:LEU:HD23	1:C:214:VAL:CG2	2.31	0.60
1:A:175:THR:OG1	4:C:334:PEG:C1	2.49	0.60
1:E:82:LEU:HD13	1:E:84:TRP:CZ2	2.36	0.60
1:A:30:ILE:CD1	1:A:71:ILE:HD11	2.31	0.60
1:C:16:LEU:O	1:C:16:LEU:HD23	2.01	0.60
1:A:202:ASN:HD21	1:C:281:VAL:HB	1.67	0.60
1:E:122(A):LYS:HD3	1:E:123:GLY:N	2.16	0.60
1:A:281:VAL:CG1	1:C:202:ASN:HD21	2.14	0.60
1:B:33:THR:CG2	1:B:77:ARG:HD2	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:122(A):LYS:O	1:F:125:ILE:HG13	2.01	0.59
1:E:114:LYS:HD2	1:E:332:TRP:HZ2	1.68	0.59
1:A:281:VAL:HB	1:C:202:ASN:HD21	1.67	0.59
1:E:129:VAL:H	1:E:133:ASN:ND2	2.01	0.59
1:A:4:ALA:HB2	1:A:89:ILE:HD12	1.83	0.59
1:A:154:LEU:HD23	1:A:214:VAL:HG21	1.85	0.59
1:D:194:ARG:HD2	1:D:205:PRO:O	2.03	0.59
1:B:3:VAL:HG13	1:B:91:ILE:HD13	1.85	0.58
1:B:129:VAL:H	1:B:133:ASN:ND2	2.00	0.58
4:A:336:PEG:H41	1:C:235:PRO:HD2	1.84	0.58
1:F:129:VAL:H	1:F:133:ASN:ND2	2.01	0.58
1:D:170:GLY:HA3	1:D:244:VAL:HG12	1.85	0.58
1:C:139:HIS:HB3	1:C:333:LYS:HZ2	1.69	0.57
1:D:176:HIS:HB3	1:D:231:ARG:HD3	1.86	0.57
1:C:154:LEU:HD23	1:C:214:VAL:HG21	1.85	0.57
1:D:16:LEU:HD23	1:D:16:LEU:O	2.04	0.57
1:F:298:MET:HB3	1:F:300[A]:MET:HE3	1.86	0.57
4:A:336:PEG:H11	1:C:235:PRO:HD3	1.85	0.57
1:E:115:LYS:HG3	1:E:332:TRP:CZ3	2.40	0.57
1:A:181:ASP:OD2	1:A:195:ARG:NH1	2.37	0.57
1:F:298:MET:HB3	1:F:300[A]:MET:HE1	1.87	0.57
1:B:18(A):TRP:CE3	1:B:27:ILE:HD12	2.40	0.56
1:C:173:THR:HG21	4:C:334:PEG:H21	1.87	0.56
1:E:219:PRO:O	1:E:222:LYS:HB2	2.05	0.56
1:E:281:VAL:HG11	1:F:48:SER:HA	1.87	0.56
1:C:3:VAL:HG22	1:C:91:ILE:HG23	1.88	0.56
1:D:119:THR:O	3:D:335:NAD:H1D	2.05	0.56
1:F:4:ALA:HB2	1:F:89:ILE:HD13	1.86	0.56
1:F:318:SER:O	1:F:322:VAL:HG23	2.06	0.56
4:A:336:PEG:H12	4:A:336:PEG:C4	2.34	0.56
4:A:336:PEG:H12	4:A:336:PEG:H42	1.88	0.56
1:E:128:TYR:HA	1:E:133:ASN:HD21	1.71	0.56
1:A:241:ASP:CB	4:C:334:PEG:H22	2.36	0.56
1:F:168:ILE:HG22	1:F:169:LYS:HD3	1.88	0.55
1:D:33:THR:HG21	1:D:77:ARG:NH1	2.22	0.55
1:A:128:TYR:HA	1:A:133:ASN:HD21	1.72	0.55
1:A:79:PRO:HA	1:A:82:LEU:HD12	1.89	0.55
1:B:28:ILE:HD11	1:B:89:ILE:CD1	2.37	0.55
1:B:300[A]:MET:HE2	1:D:226:ASN:HD22	1.70	0.55
1:A:129:VAL:H	1:A:133:ASN:HD21	1.52	0.55
1:C:179:THR:OG1	1:C:231:ARG:NH1	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:115:LYS:HG3	1:E:332:TRP:HZ3	1.72	0.55
1:A:91:ILE:HG21	1:A:117:ILE:HD12	1.88	0.55
1:B:194:ARG:HD2	1:B:205:PRO:O	2.07	0.55
1:F:9:GLY:O	1:F:13:ARG:HG3	2.07	0.55
1:B:102:ARG:CG	1:B:102:ARG:HH11	2.12	0.54
1:C:302:ASP:OD2	1:F:102:ARG:HD2	2.08	0.54
1:A:271:LEU:HD12	1:A:272:ASP:N	2.23	0.54
1:A:241:ASP:O	4:A:334:PEG:H22	2.07	0.54
1:B:102:ARG:HG2	1:B:102:ARG:NH1	2.19	0.54
1:C:139:HIS:HB2	1:C:331:ASN:O	2.08	0.54
1:B:172:MET:CE	1:B:227:GLY:HA3	2.38	0.54
1:E:256:ASN:HD21	1:E:297:THR:CB	2.21	0.54
1:A:142:PRO:HD2	7:A:357:HOH:O	2.08	0.54
1:C:175:THR:OG1	4:C:334:PEG:H41	2.08	0.53
1:F:255:VAL:O	1:F:258:ALA:HB3	2.08	0.53
1:A:306:LYS:NZ	4:A:334:PEG:H41	2.23	0.53
1:A:48:SER:HA	1:B:281:VAL:HG11	1.90	0.53
1:C:52:ILE:HG13	6:C:336:EDO:C1	2.38	0.53
1:D:133:ASN:H	1:D:133:ASN:HD22	1.56	0.53
2:I:58:ASP:CB	2:I:59:PRO:CD	2.86	0.53
1:C:52:ILE:HG13	6:C:336:EDO:H12	1.89	0.53
1:E:129:VAL:H	1:E:133:ASN:HD21	1.54	0.53
1:F:168:ILE:CD1	1:F:168:ILE:N	2.72	0.53
1:B:82:LEU:HD13	1:B:84:TRP:CZ2	2.44	0.53
4:A:336:PEG:H11	1:C:234:THR:HA	1.91	0.53
1:B:4:ALA:HB2	1:B:89:ILE:HG12	1.91	0.53
1:E:79:PRO:HA	1:E:82:LEU:HD12	1.92	0.53
1:C:183:ARG:CZ	2:H:71:ASN:HB3	2.39	0.53
1:E:195:ARG:HH22	1:E:231:ARG:NH2	2.07	0.52
1:F:212:LYS:O	1:F:215:ALA:HB3	2.09	0.52
1:C:271:LEU:HD12	1:C:290:THR:OG1	2.08	0.52
1:E:18:CYS:HB3	1:E:319:GLN:OE1	2.10	0.52
1:A:293:ASP:HB3	1:A:296:LEU:HD12	1.91	0.52
1:F:153:CYS:O	1:F:156:PRO:HD2	2.09	0.52
1:F:74:VAL:HG21	1:F:82:LEU:HD22	1.92	0.52
4:A:336:PEG:H41	1:C:235:PRO:CD	2.39	0.52
1:E:91:ILE:HB	1:E:115:LYS:HB2	1.92	0.52
1:D:33:THR:HB	1:D:77:ARG:NH1	2.25	0.52
1:F:183:ARG:CZ	2:I:71:ASN:HB3	2.40	0.52
1:B:162:ASP:OD2	1:B:220:ASN:ND2	2.43	0.52
1:E:96:THR:OG1	1:E:98:VAL:HG22	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:133:ASN:N	1:F:133:ASN:HD22	2.06	0.51
1:F:82:LEU:HD13	1:F:84:TRP:CZ2	2.46	0.51
1:B:128:TYR:HA	1:B:133:ASN:HD21	1.74	0.51
1:B:1:LEU:HD22	1:B:329:ALA:CB	2.41	0.51
1:B:33:THR:HG22	1:B:77:ARG:CD	2.41	0.51
1:F:313:ASN:O	3:F:335:NAD:H4N	2.10	0.51
2:I:58:ASP:HB3	2:I:59:PRO:CD	2.40	0.51
1:A:76:ASN:ND2	1:A:78:ASN:H	2.09	0.51
1:D:174:THR:HB	1:D:240:VAL:HG12	1.92	0.51
1:F:285:CYS:HA	1:F:315:TRP:CD1	2.46	0.51
1:F:301:GLY:O	1:F:302:ASP:HB2	2.10	0.51
1:A:243:VAL:HG23	4:A:334:PEG:H12	1.93	0.51
1:C:51:GLY:HA2	6:C:336:EDO:H21	1.91	0.51
1:D:153:CYS:HA	1:D:290:THR:HG22	1.93	0.50
1:D:160:VAL:O	1:D:164:LYS:HB2	2.11	0.50
1:B:176:HIS:HB3	1:B:231:ARG:HD3	1.92	0.50
1:C:210:ALA:O	1:C:214:VAL:HG23	2.10	0.50
1:C:79:PRO:HA	1:C:82:LEU:CD1	2.42	0.50
1:B:115:LYS:HD2	1:B:332:TRP:CZ3	2.47	0.50
1:C:154:LEU:CD2	1:C:214:VAL:HG21	2.41	0.50
1:D:129:VAL:H	1:D:133:ASN:ND2	2.10	0.50
2:H:76:TYR:O	2:H:77:ASP:HB2	2.11	0.50
1:F:256:ASN:O	1:F:260:ARG:HG3	2.11	0.50
1:B:18(A):TRP:CD2	1:B:27:ILE:HD12	2.47	0.50
6:C:336:EDO:H22	1:D:281:VAL:O	2.12	0.50
1:D:128:TYR:HA	1:D:133:ASN:HD21	1.76	0.49
1:C:51:GLY:HA2	6:C:336:EDO:C2	2.41	0.49
1:D:153:CYS:O	1:D:290:THR:HG21	2.11	0.49
2:I:71:ASN:HA	2:I:74:ARG:NH1	2.27	0.49
1:F:102:ARG:NH1	1:F:102:ARG:HG2	2.27	0.49
1:A:172:MET:HG2	1:A:173:THR:N	2.27	0.49
1:A:203[A]:ILE:HD11	1:C:232:VAL:HG21	1.93	0.49
1:B:50:LEU:HD21	7:B:352:HOH:O	2.13	0.49
1:B:37:VAL:HG22	1:B:73:VAL:HB	1.94	0.49
2:H:58:ASP:CG	2:H:59:PRO:HD3	2.33	0.49
2:H:60:LEU:HD11	2:H:72:GLU:HB2	1.93	0.49
1:B:102:ARG:NH1	7:B:360:HOH:O	2.39	0.48
1:D:133:ASN:N	1:D:133:ASN:HD22	2.11	0.48
2:H:62:GLU:OE1	2:H:65:LYS:HE3	2.12	0.48
1:E:156:PRO:HB2	1:E:290:THR:HG21	1.95	0.48
1:D:206:THR:HG22	1:D:229:ALA:HB3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:ASN:HD22	1:B:133:ASN:H	1.61	0.48
1:E:53:PHE:O	1:E:55:ALA:N	2.46	0.48
1:A:169:LYS:NZ	1:C:303:ASP:OD1	2.26	0.48
1:A:1:LEU:HD21	1:A:332:TRP:CE3	2.49	0.48
1:C:51:GLY:HA2	6:C:336:EDO:C1	2.43	0.48
1:B:269:GLY:C	1:B:270:ILE:HD12	2.34	0.47
1:C:206:THR:OG1	1:C:207:SER:N	2.47	0.47
1:D:18(A):TRP:CE3	1:D:27:ILE:HD12	2.49	0.47
1:F:139:HIS:CD2	1:F:333:LYS:H	2.33	0.47
1:C:212:LYS:HE2	1:C:212:LYS:HB2	1.52	0.47
1:D:129:VAL:H	1:D:133:ASN:HD21	1.61	0.47
1:D:45:LYS:HD2	1:D:52:ILE:HG23	1.96	0.47
1:D:77:ARG:HH22	2:G:69:GLU:HG3	1.79	0.47
1:A:100:VAL:HG23	1:A:122(A):LYS:HG2	1.97	0.47
1:B:202:ASN:ND2	1:D:279:VAL:HG23	2.29	0.47
1:F:37:VAL:HG21	1:F:61:GLU:O	2.15	0.47
1:B:162:ASP:HB2	1:B:167:ILE:HD12	1.97	0.47
1:E:195:ARG:NH2	1:E:231:ARG:NH2	2.62	0.47
1:C:114:LYS:O	1:C:115:LYS:HG2	2.15	0.47
1:D:122(A):LYS:HD3	1:D:122(A):LYS:HA	1.83	0.47
1:A:79:PRO:HA	1:A:82:LEU:CD1	2.45	0.47
1:F:181:ASP:OD1	1:F:195:ARG:NE	2.47	0.47
1:B:133:ASN:C	1:B:133:ASN:ND2	2.66	0.47
1:E:18(A):TRP:HH2	1:E:69:LYS:HE2	1.80	0.47
1:A:241:ASP:OD2	4:C:334:PEG:H22	2.16	0.46
1:D:3:VAL:HG22	1:D:91[A]:ILE:HG23	1.96	0.46
1:B:300[A]:MET:SD	1:D:226:ASN:HB3	2.55	0.46
1:D:91[A]:ILE:HD11	1:D:117:ILE:HD12	1.98	0.46
1:C:302:ASP:CG	1:F:102:ARG:HD2	2.35	0.46
1:F:264:GLU:O	1:F:268:LYS:HE2	2.16	0.46
1:D:16:LEU:HD23	1:D:16:LEU:C	2.35	0.46
1:A:285:CYS:HA	1:A:315:TRP:CD1	2.51	0.46
1:E:30:ILE:CD1	1:E:71:ILE:HD11	2.45	0.46
1:F:133:ASN:HD22	1:F:133:ASN:H	1.64	0.46
1:A:306:LYS:HZ2	4:A:334:PEG:H41	1.81	0.46
1:B:79:PRO:HA	1:B:82:LEU:HD12	1.96	0.46
1:A:281:VAL:CB	1:C:202:ASN:HD21	2.26	0.46
1:E:319:GLN:OE1	1:E:319:GLN:HA	2.16	0.46
1:F:134:ALA:HB1	1:F:327:ILE:HD13	1.98	0.46
1:B:185:LEU:O	1:B:186:ASP:C	2.53	0.46
1:B:278:LEU:O	1:D:194:ARG:NH1	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:249:LYS:HG2	1:E:302:ASP:HB3	1.97	0.46
1:A:34:GLY:O	1:A:39:GLN:OE1	2.33	0.46
1:B:119:THR:O	3:B:335:NAD:H1D	2.15	0.46
1:C:31:ASN:OD1	1:C:74:VAL:HG23	2.15	0.46
1:D:33:THR:HB	1:D:77:ARG:HH11	1.81	0.46
1:A:241:ASP:HB3	4:A:334:PEG:H22	1.98	0.46
1:F:3:VAL:HG22	1:F:91:ILE:HG23	1.98	0.46
1:A:1:LEU:HD22	1:A:329:ALA:HA	1.98	0.45
1:C:37:VAL:HG22	1:C:73:VAL:HB	1.98	0.45
1:D:139:HIS:CG	1:D:139:HIS:O	2.69	0.45
1:A:10:ARG:HH11	1:A:13:ARG:NH2	2.14	0.45
1:B:153:CYS:O	1:B:156:PRO:HD2	2.16	0.45
1:A:271:LEU:CD1	1:A:272:ASP:N	2.79	0.45
1:C:139:HIS:CD2	1:C:333:LYS:HD2	2.51	0.45
1:D:253:GLU:H	1:D:253:GLU:CD	2.20	0.45
1:B:109:ILE:HG23	1:B:113:ALA:O	2.16	0.45
1:B:214:VAL:HG13	1:B:218:LEU:HD12	1.98	0.45
1:F:101:ASP:HB3	1:F:122(A):LYS:HB2	1.99	0.45
1:A:30:ILE:CG1	1:A:71:ILE:HD11	2.47	0.45
1:F:114:LYS:O	1:F:115:LYS:HG2	2.17	0.45
1:A:155:ALA:N	1:A:156:PRO:CD	2.79	0.44
1:D:250:THR:OG1	1:D:254:GLU:HB3	2.17	0.44
1:A:154:LEU:HD23	1:A:214:VAL:CG2	2.46	0.44
1:C:8:PHE:N	1:C:32:ASP:OD1	2.43	0.44
2:G:72:GLU:H	2:G:72:GLU:HG2	1.26	0.44
1:B:130:VAL:HB	1:B:320:ARG:HD3	1.99	0.44
1:D:218:LEU:HB3	1:D:221:LEU:HD23	2.00	0.44
1:C:190:HIS:NE2	2:H:74:ARG:O	2.47	0.44
1:B:91:ILE:HD11	1:B:93:ILE:HD11	1.98	0.44
1:A:115:LYS:HB2	1:A:115:LYS:HE2	1.85	0.44
1:B:169:LYS:NZ	1:D:301:GLY:O	2.51	0.44
1:D:169:LYS:HE2	1:D:245:GLN:OE1	2.17	0.44
1:A:10:ARG:HH11	1:A:13:ARG:HH21	1.64	0.44
1:E:34:GLY:O	1:E:39:GLN:OE1	2.36	0.43
1:F:84:TRP:CE3	1:F:84:TRP:HA	2.53	0.43
1:A:241:ASP:O	4:A:334:PEG:H11	2.18	0.43
1:B:133:ASN:HD22	1:B:133:ASN:N	2.14	0.43
4:A:334:PEG:H42	1:C:243:VAL:CG2	2.48	0.43
1:C:37:VAL:HG13	1:C:64:ILE:CG2	2.48	0.43
1:C:3:VAL:HA	1:C:91:ILE:O	2.18	0.43
1:A:210:ALA:O	1:A:214:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:HIS:O	1:A:139:HIS:CG	2.72	0.43
1:B:175:THR:HB	1:B:239:VAL:HG13	2.00	0.43
4:A:334:PEG:H42	1:C:243:VAL:HG23	2.00	0.43
1:F:293:ASP:CG	1:F:296:LEU:CD1	2.85	0.43
1:F:84:TRP:HE3	1:F:84:TRP:HA	1.84	0.43
1:D:249:LYS:CG	1:D:302:ASP:HB3	2.49	0.43
1:D:332:TRP:O	1:D:333:LYS:HD2	2.19	0.43
1:D:77:ARG:HG2	1:D:77:ARG:H	1.56	0.43
1:B:202:ASN:HD21	1:D:281:VAL:CB	2.32	0.43
1:D:153:CYS:O	1:D:290:THR:CG2	2.67	0.43
1:F:31:ASN:HB2	1:F:74:VAL:HG22	2.00	0.43
1:B:234:THR:HA	1:B:235:PRO:HD3	1.82	0.43
1:B:240:VAL:HG23	1:B:309:ALA:HB3	2.01	0.43
1:B:281:VAL:HG12	1:D:202:ASN:HD21	1.83	0.43
1:D:79:PRO:HA	1:D:82:LEU:CD1	2.48	0.42
1:A:84:TRP:CE3	1:A:84:TRP:HA	2.54	0.42
1:E:176:HIS:O	1:E:231:ARG:HA	2.19	0.42
1:B:239:VAL:HA	1:B:309:ALA:O	2.19	0.42
1:B:245:GLN:HA	1:B:303:ASP:O	2.19	0.42
1:E:139:HIS:CE1	1:E:333:LYS:H	2.37	0.42
1:B:1:LEU:HD22	1:B:329:ALA:HB2	2.01	0.42
1:A:27:ILE:HG13	1:A:27:ILE:O	2.20	0.42
1:C:146:ASN:CG	1:C:146:ASN:O	2.57	0.42
1:F:2:LYS:HB3	1:F:28:ILE:HD11	2.02	0.42
1:D:84:TRP:CE3	1:D:84:TRP:HA	2.54	0.42
1:F:204:VAL:HB	1:F:231:ARG:HB2	2.01	0.42
1:C:16:LEU:C	1:C:16:LEU:HD23	2.40	0.42
1:C:177:SER:OG	1:C:236:ASN:C	2.58	0.42
1:C:191:ARG:CZ	1:C:191:ARG:HB2	2.49	0.42
1:C:173:THR:CG2	4:C:334:PEG:H21	2.50	0.42
1:E:165:PHE:HA	1:E:248:LYS:HB2	2.01	0.42
1:B:172:MET:HE2	1:B:227:GLY:HA3	2.01	0.41
1:D:115:LYS:HG3	1:D:332:TRP:HZ3	1.85	0.41
1:E:181:ASP:OD2	1:E:195:ARG:HD3	2.20	0.41
1:E:139:HIS:CD2	1:E:333:LYS:HB2	2.55	0.41
1:A:203[A]:ILE:HD11	1:C:232:VAL:CG2	2.50	0.41
1:B:299:VAL:O	1:B:300[A]:MET:HE3	2.20	0.41
1:C:133:ASN:C	1:C:133:ASN:ND2	2.72	0.41
1:C:51:GLY:HA2	6:C:336:EDO:H12	2.01	0.41
1:E:332:TRP:O	1:E:333:LYS:HG3	2.20	0.41
1:F:109:ILE:HA	1:F:113:ALA:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:MET:O	1:B:227:GLY:HA2	2.20	0.41
1:C:175:THR:OG1	4:C:334:PEG:C4	2.68	0.41
1:B:202:ASN:ND2	1:D:281:VAL:HB	2.36	0.41
1:B:292:ILE:HD12	1:B:307:VAL:CG1	2.51	0.41
1:C:240:VAL:O	1:C:308:ILE:HA	2.20	0.41
1:A:201:LEU:O	4:A:336:PEG:H42	2.21	0.41
1:A:169:LYS:HD3	1:C:301:GLY:HA3	2.01	0.41
1:D:311:TYR:CD1	1:D:311:TYR:N	2.89	0.41
1:E:301:GLY:O	1:E:302:ASP:HB2	2.21	0.41
1:C:129:VAL:N	1:C:133:ASN:HD21	2.14	0.41
1:A:241:ASP:HB3	4:C:334:PEG:H22	2.03	0.41
1:A:29:ALA:HA	1:A:71:ILE:HD12	2.03	0.41
1:F:277:PRO:C	1:F:278:LEU:HD23	2.41	0.41
1:C:59:PRO:HA	1:C:64:ILE:HA	2.02	0.41
1:D:176:HIS:O	1:D:231:ARG:HA	2.20	0.41
1:F:278:LEU:N	1:F:278:LEU:HD23	2.35	0.41
1:D:33:THR:CB	1:D:77:ARG:NH1	2.84	0.41
1:D:245:GLN:HA	1:D:303:ASP:O	2.20	0.41
1:F:251:PHE:CZ	1:F:254:GLU:HB2	2.56	0.41
1:F:239:VAL:HA	1:F:309:ALA:O	2.20	0.41
1:C:154:LEU:HD11	1:C:242:LEU:HD13	2.03	0.40
1:C:249:LYS:HE3	1:F:124:ASP:HB2	2.03	0.40
1:F:55:ALA:HB1	1:F:67:ASP:OD1	2.21	0.40
1:F:176:HIS:HB3	1:F:231:ARG:HD3	2.02	0.40
1:D:250:THR:OG1	1:D:254:GLU:OE1	2.39	0.40
1:E:316:GLY:O	1:E:319:GLN:HB2	2.21	0.40
1:A:65:SER:HA	1:A:69:LYS:O	2.21	0.40
1:C:27:ILE:HD11	1:C:71:ILE:HD13	2.02	0.40
1:D:84:TRP:HA	1:D:84:TRP:HE3	1.85	0.40
1:F:271:LEU:HD13	1:F:272:ASP:N	2.37	0.40

All (1) 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)
1:E:169:LYS:NZ	1:E:303:ASP:OD1[4_565]	2.18	0.02

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	336/337 (100%)	323 (96%)	12 (4%)	1 (0%)	41	37
1	B	335/337 (99%)	309 (92%)	25 (8%)	1 (0%)	41	37
1	C	334/337 (99%)	317 (95%)	16 (5%)	1 (0%)	41	37
1	D	336/337 (100%)	317 (94%)	18 (5%)	1 (0%)	41	37
1	E	334/337 (99%)	307 (92%)	25 (8%)	2 (1%)	25	19
1	F	335/337 (99%)	305 (91%)	27 (8%)	3 (1%)	17	11
2	G	19/82 (23%)	19 (100%)	0	0	100	100
2	H	18/82 (22%)	16 (89%)	2 (11%)	0	100	100
2	I	17/82 (21%)	15 (88%)	2 (12%)	0	100	100
All	All	2064/2268 (91%)	1928 (93%)	127 (6%)	9 (0%)	34	30

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	237	VAL
1	B	237	VAL
1	C	237	VAL
1	D	237	VAL
1	E	237	VAL
1	F	237	VAL
1	F	188	SER
1	F	166	GLY
1	E	54	ASP

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	280/279 (100%)	262 (94%)	18 (6%)	17	13
1	B	280/279 (100%)	261 (93%)	19 (7%)	16	11
1	C	279/279 (100%)	260 (93%)	19 (7%)	16	11
1	D	280/279 (100%)	265 (95%)	15 (5%)	22	18
1	E	279/279 (100%)	259 (93%)	20 (7%)	14	9
1	F	280/279 (100%)	254 (91%)	26 (9%)	9	5
2	G	21/65 (32%)	20 (95%)	1 (5%)	25	22
2	H	20/65 (31%)	14 (70%)	6 (30%)	0	0
2	I	19/65 (29%)	18 (95%)	1 (5%)	22	18
All	All	1738/1869 (93%)	1613 (93%)	125 (7%)	15	9

All (125) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LEU
1	A	27	ILE
1	A	60	SER
1	A	71	ILE
1	A	76	ASN
1	A	84	TRP
1	A	86	GLU
1	A	94	GLU
1	A	133	ASN
1	A	163	GLN
1	A	172	MET
1	A	181	ASP
1	A	191	ARG
1	A	212	LYS
1	A	271	LEU
1	A	276	GLU
1	A	290	THR
1	A	311	TYR
1	B	8	PHE
1	B	74	VAL
1	B	85	LYS
1	B	86	GLU
1	B	91	ILE

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Mol	Chain	Res	Type
1	B	94	GLU
1	B	133	ASN
1	B	169	LYS
1	B	172	MET
1	B	191	ARG
1	B	226	ASN
1	B	268	LYS
1	B	271	LEU
1	B	289	SER
1	B	290	THR
1	B	295	SER
1	B	300[A]	MET
1	B	300[B]	MET
1	B	311	TYR
1	C	1	LEU
1	C	21	LYS
1	C	27	ILE
1	C	76	ASN
1	C	84	TRP
1	C	91	ILE
1	C	122(A)	LYS
1	C	133	ASN
1	C	145	SER
1	C	163	GLN
1	C	164	LYS
1	C	169	LYS
1	C	191	ARG
1	C	212	LYS
1	C	219	PRO
1	C	222	LYS
1	C	281	VAL
1	C	290	THR
1	C	333	LYS
1	D	33	THR
1	D	70	ILE
1	D	84	TRP
1	D	85	LYS
1	D	91[A]	ILE
1	D	91[B]	ILE
1	D	133	ASN
1	D	164	LYS
1	D	169	LYS

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Mol	Chain	Res	Type
1	D	172	MET
1	D	174	THR
1	D	206	THR
1	D	271	LEU
1	D	290	THR
1	D	311	TYR
1	E	1	LEU
1	E	27	ILE
1	E	41	SER
1	E	61	GLU
1	E	71	ILE
1	E	72	GLN
1	E	76	ASN
1	E	77	ARG
1	E	81	LEU
1	E	91	ILE
1	E	94	GLU
1	E	122(A)	LYS
1	E	133	ASN
1	E	169	LYS
1	E	191	ARG
1	E	212	LYS
1	E	248	LYS
1	E	276	GLU
1	E	290	THR
1	E	311	TYR
1	F	1	LEU
1	F	22	ASP
1	F	26	ASP
1	F	33	THR
1	F	38	LYS
1	F	39	GLN
1	F	60	SER
1	F	64	ILE
1	F	74	VAL
1	F	85	LYS
1	F	91	ILE
1	F	94	GLU
1	F	125	ILE
1	F	126	PRO
1	F	133	ASN
1	F	138	SER

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Mol	Chain	Res	Type
1	F	164	LYS
1	F	168	ILE
1	F	191	ARG
1	F	261	ASP
1	F	271	LEU
1	F	289	SER
1	F	290	THR
1	F	300[A]	MET
1	F	300[B]	MET
1	F	333	LYS
2	G	72	GLU
2	H	58	ASP
2	H	61	GLU
2	H	65	LYS
2	H	67	ASN
2	H	72	GLU
2	H	77	ASP
2	I	58	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	ASN
1	A	133	ASN
1	A	139	HIS
1	A	152	ASN
1	A	202	ASN
1	A	256	ASN
1	A	330	ASN
1	B	133	ASN
1	B	152	ASN
1	B	202	ASN
1	B	226	ASN
1	B	256	ASN
1	B	330	ASN
1	C	76	ASN
1	C	133	ASN
1	C	139	HIS
1	C	152	ASN
1	C	202	ASN
1	C	256	ASN
1	C	330	ASN

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Mol	Chain	Res	Type
1	D	133	ASN
1	D	202	ASN
1	D	226	ASN
1	D	256	ASN
1	D	330	ASN
1	E	39	GLN
1	E	133	ASN
1	E	152	ASN
1	E	202	ASN
1	E	256	ASN
1	F	18(B)	HIS
1	F	133	ASN
1	F	152	ASN
1	F	256	ASN
1	F	330	ASN
1	F	331	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

12 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
4	PEG	A	336	-	6,6,6	0.59	0	5,5,5	0.87	0
4	PEG	A	334	-	6,6,6	0.41	0	5,5,5	0.71	0
3	NAD	F	335	-	42,48,48	1.59	4 (9%)	50,73,73	1.41	5 (10%)
3	NAD	E	335	-	42,48,48	1.68	5 (11%)	50,73,73	1.34	6 (12%)
3	NAD	B	335	-	42,48,48	1.75	4 (9%)	50,73,73	1.54	4 (8%)
3	NAD	A	335	-	42,48,48	1.68	4 (9%)	50,73,73	1.52	5 (10%)
3	NAD	C	335	-	42,48,48	1.63	5 (11%)	50,73,73	1.51	5 (10%)
6	EDO	C	336	-	3,3,3	0.51	0	2,2,2	0.28	0
5	SO4	A	337	-	4,4,4	0.28	0	6,6,6	0.52	0
3	NAD	D	335	-	42,48,48	1.60	5 (11%)	50,73,73	1.53	6 (12%)
5	SO4	A	338	-	4,4,4	0.32	0	6,6,6	0.39	0
4	PEG	C	334	-	6,6,6	0.53	0	5,5,5	0.83	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
4	PEG	A	334	-	-	3/4/4/4	-
3	NAD	F	335	-	-	5/26/62/62	0/5/5/5
3	NAD	E	335	-	-	5/26/62/62	0/5/5/5
3	NAD	B	335	-	-	5/26/62/62	0/5/5/5
3	NAD	A	335	-	-	7/26/62/62	0/5/5/5
3	NAD	C	335	-	-	6/26/62/62	0/5/5/5
6	EDO	C	336	-	-	0/1/1/1	-
3	NAD	D	335	-	-	2/26/62/62	0/5/5/5
4	PEG	A	336	-	-	2/4/4/4	-
4	PEG	C	334	-	-	4/4/4/4	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	335	NAD	O7N-C7N	8.21	1.39	1.24
3	B	335	NAD	O7N-C7N	8.14	1.39	1.24
3	E	335	NAD	O7N-C7N	8.05	1.39	1.24
3	F	335	NAD	O7N-C7N	7.62	1.38	1.24
3	D	335	NAD	O7N-C7N	7.11	1.37	1.24
3	C	335	NAD	O7N-C7N	6.75	1.37	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	335	NAD	C2A-N3A	4.17	1.38	1.32
3	C	335	NAD	C2A-N3A	3.98	1.38	1.32
3	B	335	NAD	C2A-N3A	3.98	1.38	1.32
3	A	335	NAD	C2A-N3A	3.88	1.38	1.32
3	F	335	NAD	C2A-N3A	3.45	1.37	1.32
3	C	335	NAD	C2N-N1N	3.08	1.38	1.35
3	D	335	NAD	C2A-N3A	3.07	1.37	1.32
3	E	335	NAD	C2A-N1A	2.86	1.39	1.33
3	C	335	NAD	C2A-N1A	2.81	1.39	1.33
3	A	335	NAD	C2A-N1A	2.80	1.39	1.33
3	D	335	NAD	C2N-N1N	2.78	1.38	1.35
3	B	335	NAD	C2N-N1N	2.75	1.38	1.35
3	F	335	NAD	C2A-N1A	2.61	1.38	1.33
3	B	335	NAD	C2A-N1A	2.52	1.38	1.33
3	E	335	NAD	C2N-N1N	2.39	1.37	1.35
3	F	335	NAD	O4B-C4B	-2.38	1.39	1.45
3	D	335	NAD	O2D-C2D	-2.37	1.37	1.43
3	A	335	NAD	C2N-N1N	2.09	1.37	1.35
3	E	335	NAD	O4B-C1B	2.07	1.44	1.41
3	C	335	NAD	PA-O1A	-2.01	1.43	1.50
3	D	335	NAD	C2A-N1A	2.01	1.37	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	335	NAD	N3A-C2A-N1A	-6.28	118.87	128.68
3	C	335	NAD	N3A-C2A-N1A	-5.77	119.65	128.68
3	F	335	NAD	N3A-C2A-N1A	-5.74	119.70	128.68
3	D	335	NAD	N3A-C2A-N1A	-5.59	119.94	128.68
3	B	335	NAD	C1B-N9A-C4A	-5.19	117.51	126.64
3	A	335	NAD	O7N-C7N-C3N	-4.97	113.68	119.63
3	E	335	NAD	N3A-C2A-N1A	-4.92	120.98	128.68
3	A	335	NAD	C3N-C7N-N7N	4.72	123.41	117.75
3	D	335	NAD	C3N-C7N-N7N	4.50	123.15	117.75
3	A	335	NAD	N3A-C2A-N1A	-4.50	121.65	128.68
3	F	335	NAD	O7N-C7N-C3N	-4.26	114.53	119.63
3	D	335	NAD	C1B-N9A-C4A	-3.65	120.22	126.64
3	B	335	NAD	O7N-C7N-C3N	-3.46	115.50	119.63
3	C	335	NAD	C3N-C7N-N7N	3.36	121.78	117.75
3	C	335	NAD	O7N-C7N-C3N	-3.14	115.88	119.63
3	F	335	NAD	C3N-C7N-N7N	3.03	121.38	117.75
3	A	335	NAD	C1B-N9A-C4A	-2.95	121.46	126.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	335	NAD	O7N-C7N-C3N	-2.94	116.11	119.63
3	E	335	NAD	C3N-C7N-N7N	2.85	121.17	117.75
3	F	335	NAD	C1B-N9A-C4A	-2.84	121.65	126.64
3	C	335	NAD	C1B-N9A-C4A	-2.80	121.72	126.64
3	E	335	NAD	C1B-N9A-C4A	-2.79	121.73	126.64
3	C	335	NAD	O3B-C3B-C4B	-2.58	103.58	111.05
3	D	335	NAD	O2N-PN-O1N	2.37	123.97	112.24
3	D	335	NAD	O7N-C7N-N7N	-2.36	119.22	122.58
3	B	335	NAD	C3N-C7N-N7N	2.21	120.40	117.75
3	F	335	NAD	O2N-PN-O1N	2.16	122.92	112.24
3	D	335	NAD	PN-O3-PA	-2.15	125.45	132.83
3	E	335	NAD	C6N-N1N-C2N	-2.11	120.06	121.97
3	A	335	NAD	C2N-C3N-C4N	2.03	120.56	118.26
3	E	335	NAD	O4B-C4B-C3B	2.01	109.09	105.11

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	335	NAD	O4D-C1D-N1N-C2N
3	F	335	NAD	O4D-C1D-N1N-C6N
3	F	335	NAD	C2D-C1D-N1N-C2N
3	F	335	NAD	C2D-C1D-N1N-C6N
3	E	335	NAD	O4D-C1D-N1N-C2N
3	E	335	NAD	O4D-C1D-N1N-C6N
3	E	335	NAD	C2D-C1D-N1N-C6N
3	B	335	NAD	O4D-C1D-N1N-C2N
3	B	335	NAD	O4D-C1D-N1N-C6N
3	B	335	NAD	C2D-C1D-N1N-C2N
3	B	335	NAD	C2D-C1D-N1N-C6N
3	A	335	NAD	O4D-C1D-N1N-C2N
3	A	335	NAD	O4D-C1D-N1N-C6N
3	A	335	NAD	C2D-C1D-N1N-C2N
3	A	335	NAD	C2D-C1D-N1N-C6N
3	C	335	NAD	O4D-C1D-N1N-C2N
3	C	335	NAD	O4D-C1D-N1N-C6N
3	C	335	NAD	C2D-C1D-N1N-C2N
3	C	335	NAD	C2D-C1D-N1N-C6N
3	D	335	NAD	O4D-C1D-N1N-C2N
4	A	334	PEG	O2-C3-C4-O4
4	C	334	PEG	O1-C1-C2-O2
4	A	336	PEG	O2-C3-C4-O4

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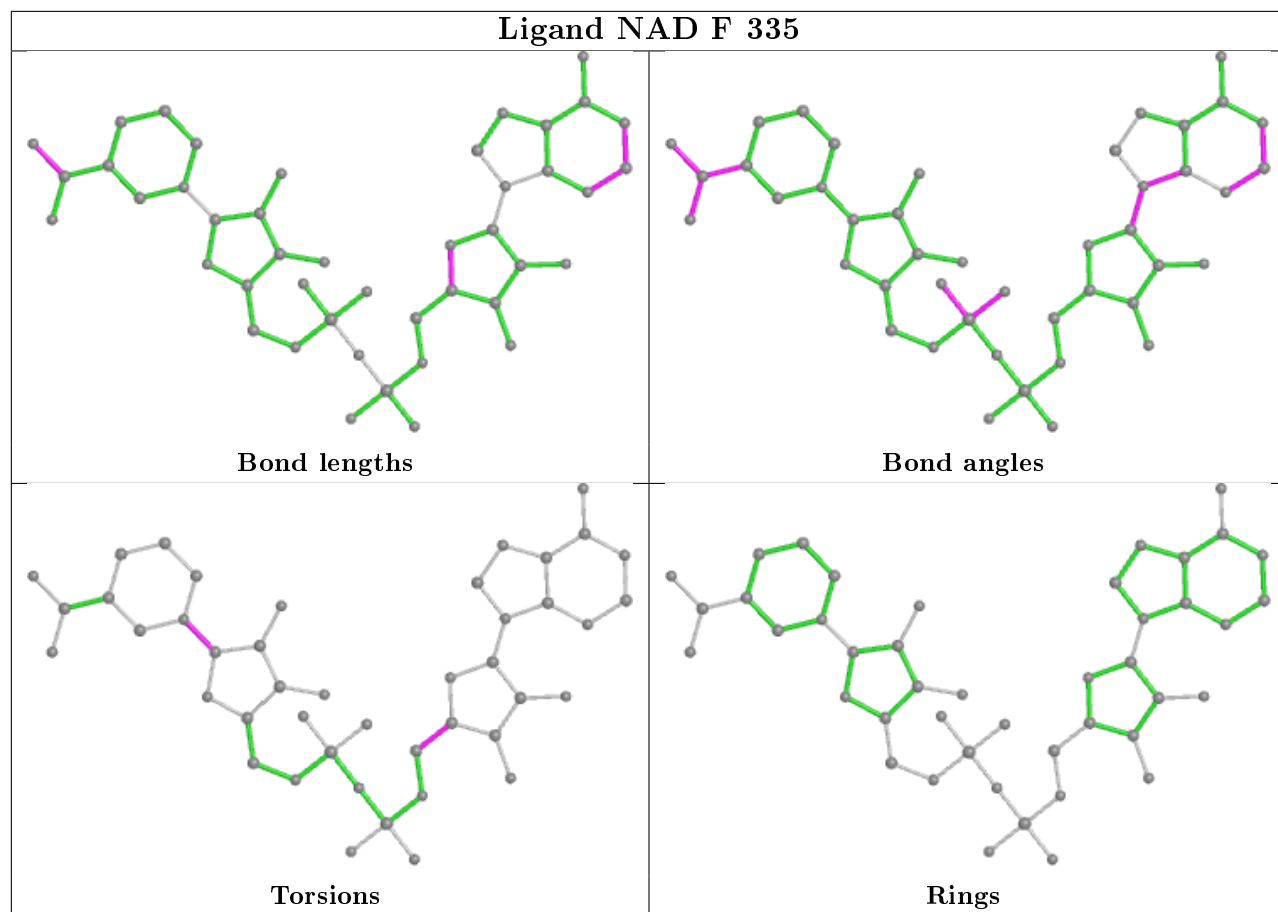
Mol	Chain	Res	Type	Atoms
4	C	334	PEG	O2-C3-C4-O4
4	C	334	PEG	C4-C3-O2-C2
3	E	335	NAD	O4B-C4B-C5B-O5B
4	A	334	PEG	C4-C3-O2-C2
4	C	334	PEG	C1-C2-O2-C3
3	D	335	NAD	O4B-C4B-C5B-O5B
3	C	335	NAD	PN-O3-PA-O1A
4	A	334	PEG	O1-C1-C2-O2
4	A	336	PEG	C1-C2-O2-C3
3	A	335	NAD	C4N-C3N-C7N-N7N
3	A	335	NAD	C4N-C3N-C7N-O7N
3	E	335	NAD	C2D-C1D-N1N-C2N
3	F	335	NAD	O4B-C4B-C5B-O5B
3	B	335	NAD	O4B-C4B-C5B-O5B
3	A	335	NAD	O4B-C4B-C5B-O5B
3	C	335	NAD	O4B-C4B-C5B-O5B

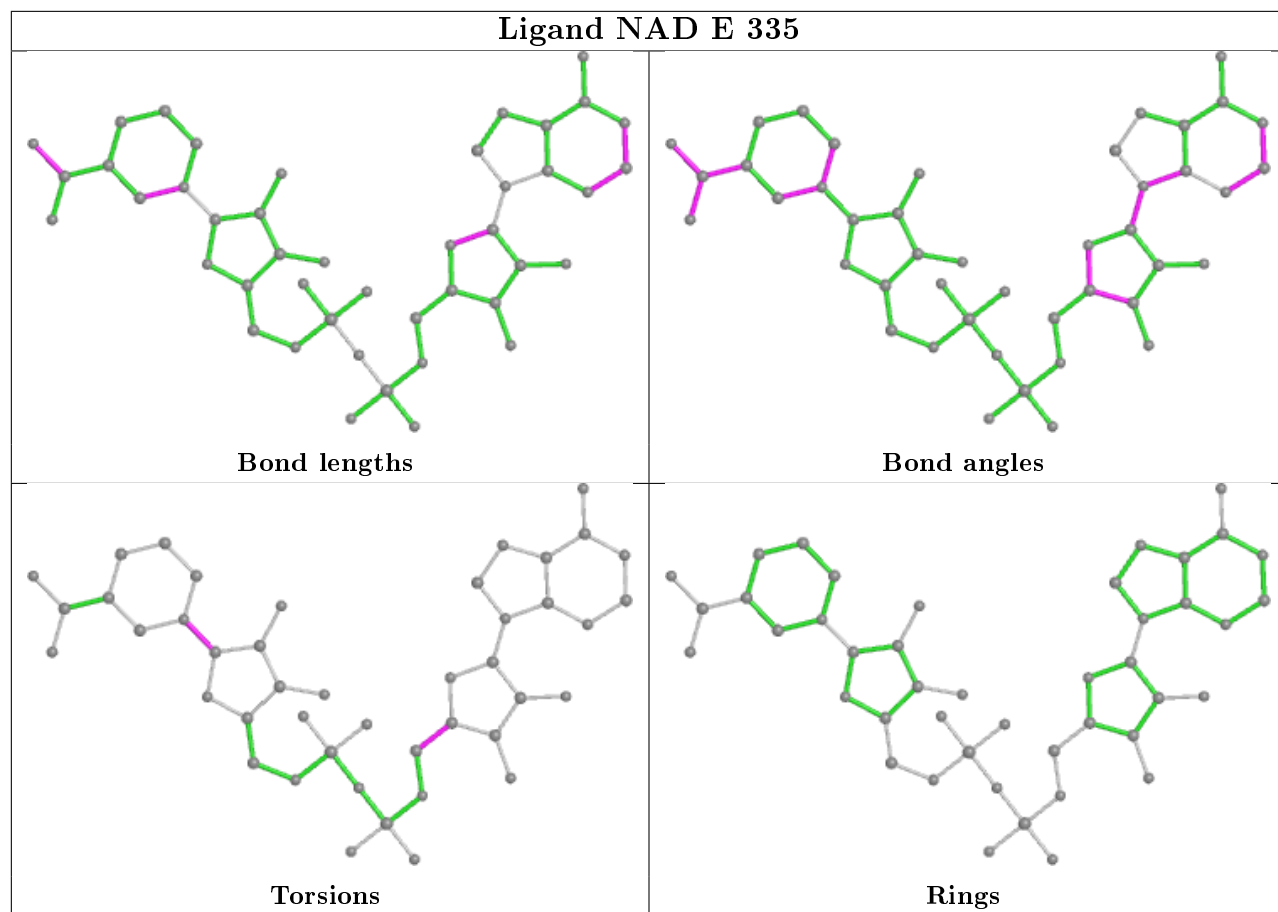
There are no ring outliers.

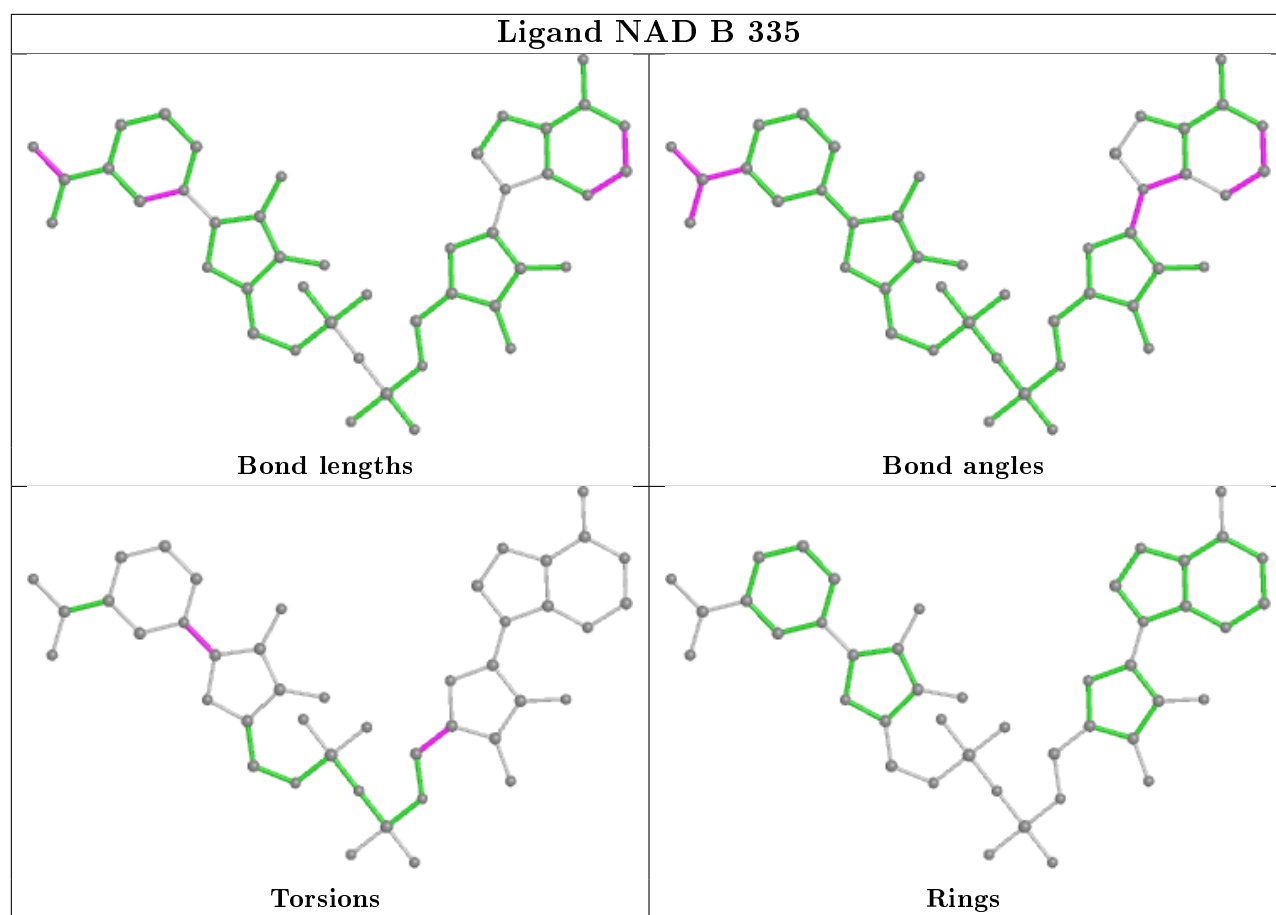
7 monomers are involved in 44 short contacts:

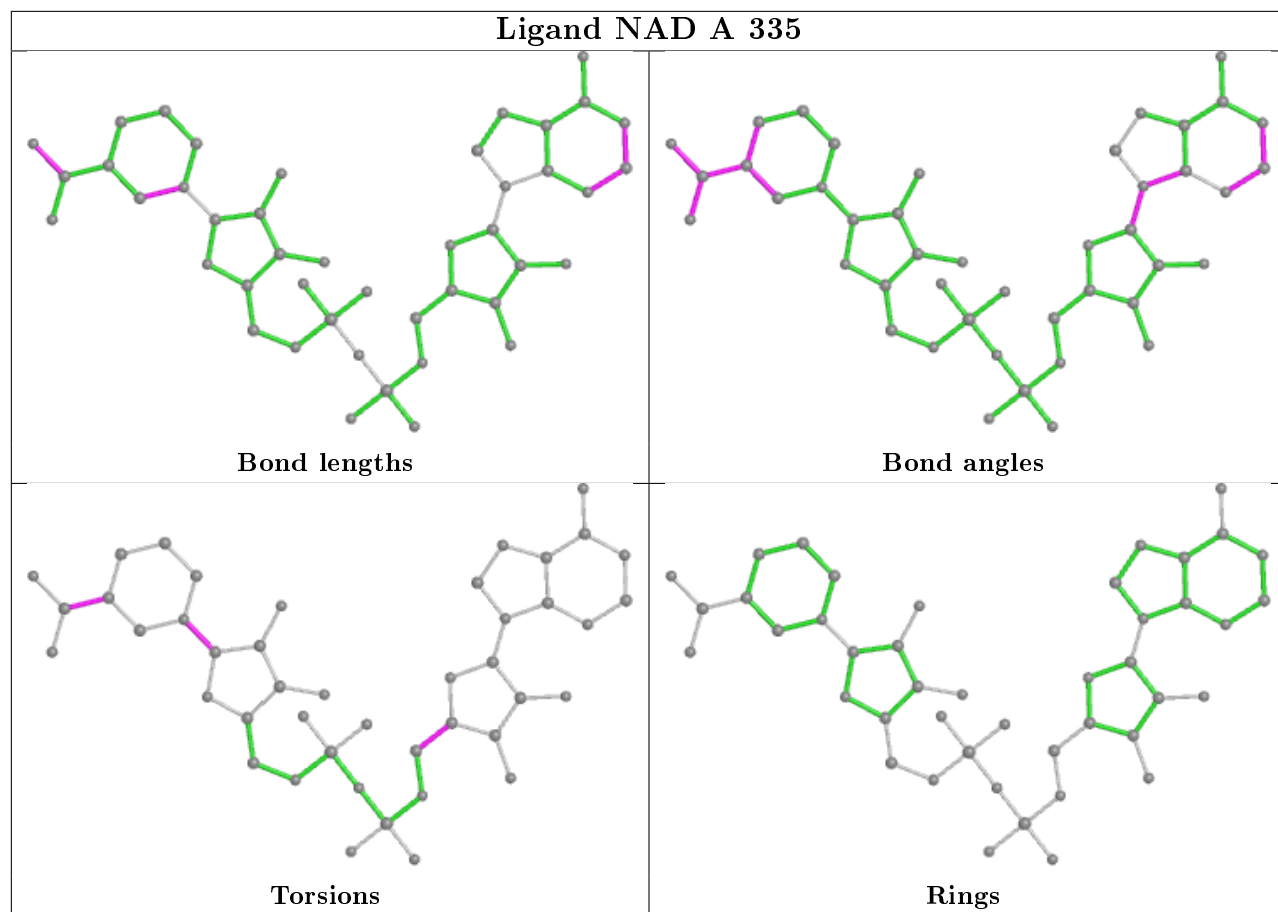
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	336	PEG	8	0
4	A	334	PEG	11	0
3	F	335	NAD	1	0
3	B	335	NAD	1	0
6	C	336	EDO	7	0
3	D	335	NAD	1	0
4	C	334	PEG	15	0

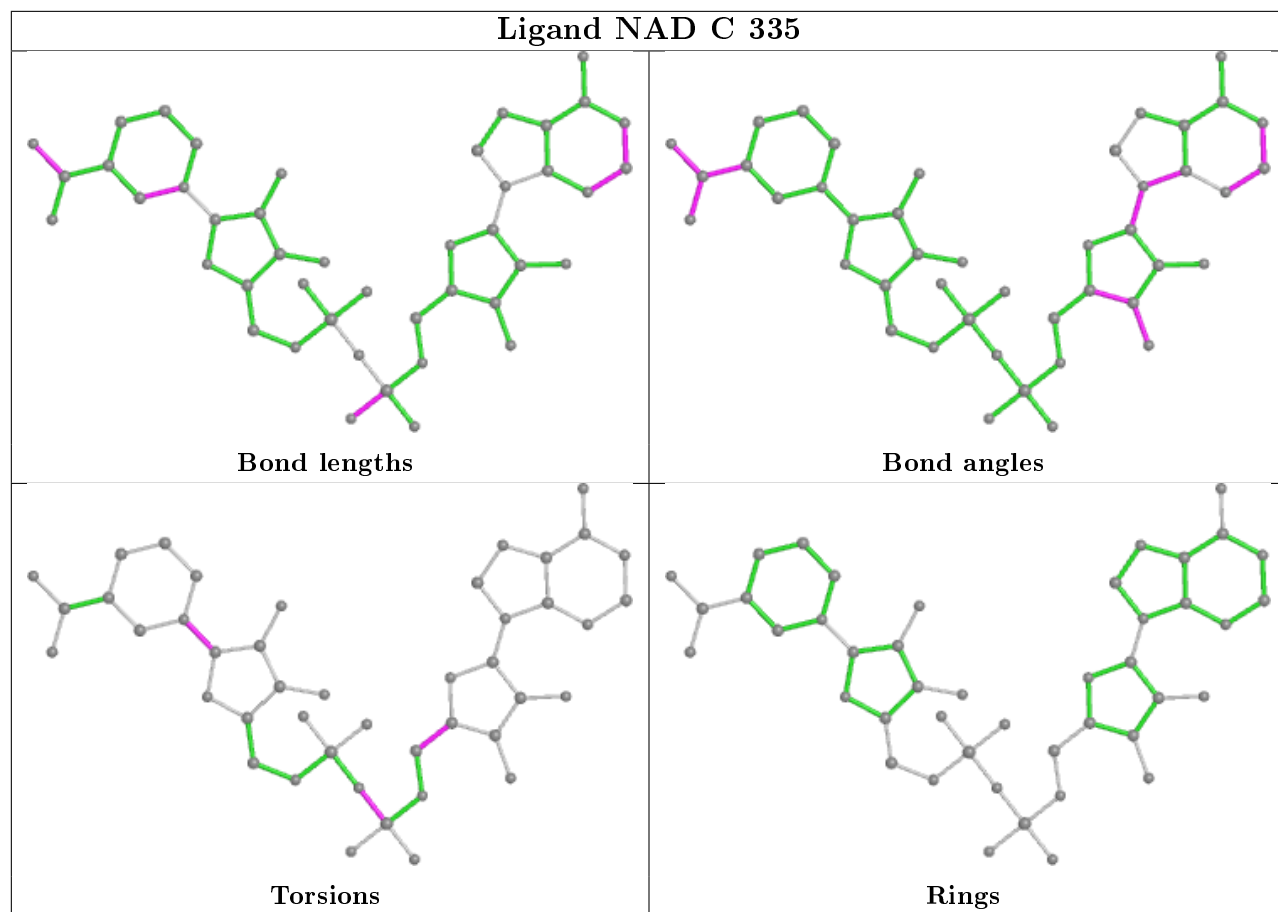
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

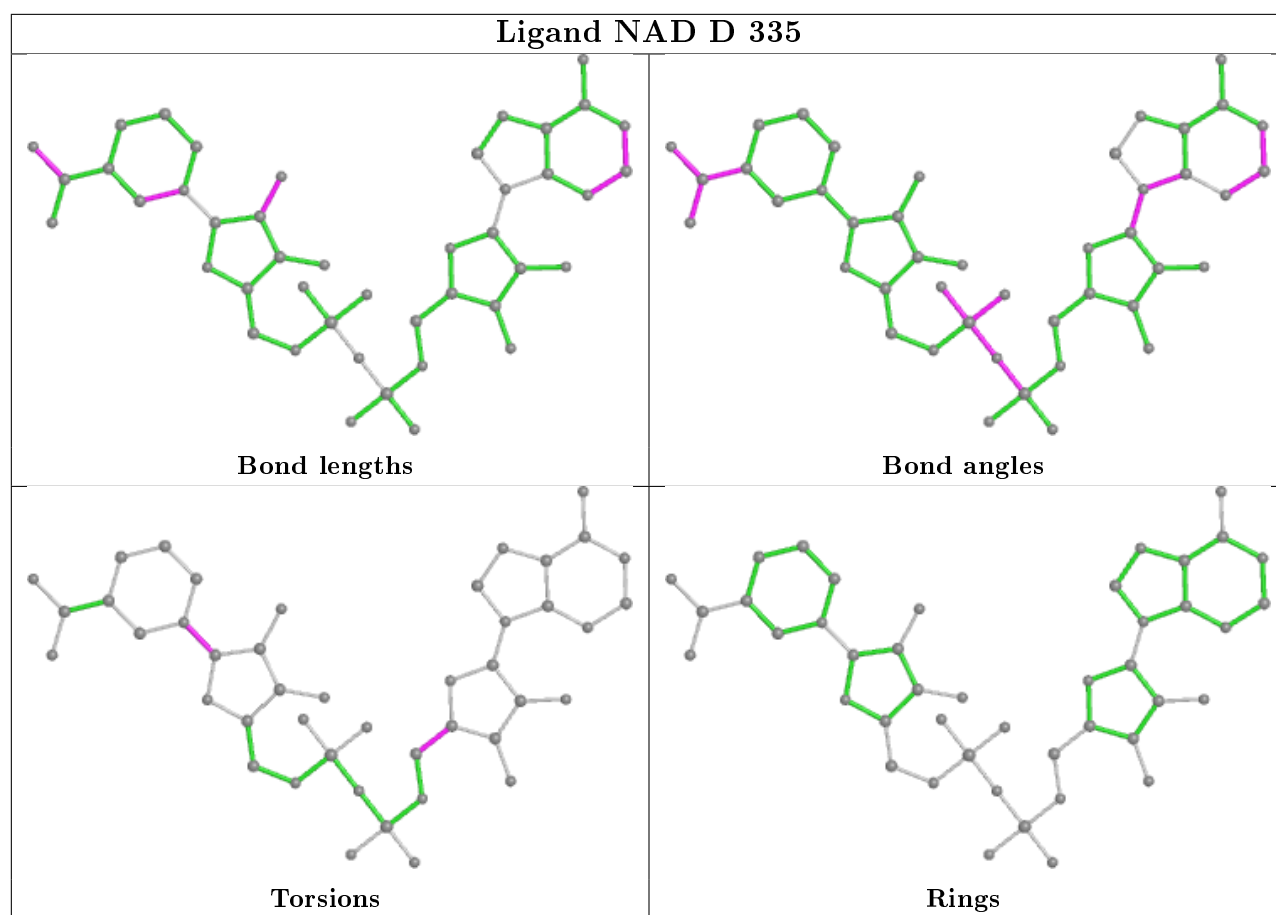












5.7 Other polymers [i](#)

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	337/337 (100%)	-0.11	2 (0%) 89 88	9, 23, 37, 48	0
1	B	336/337 (99%)	0.22	4 (1%) 79 78	13, 34, 48, 62	0
1	C	336/337 (99%)	0.06	7 (2%) 63 62	10, 27, 43, 55	0
1	D	337/337 (100%)	0.01	5 (1%) 73 72	13, 28, 41, 51	0
1	E	336/337 (99%)	0.40	19 (5%) 23 23	17, 35, 54, 61	0
1	F	336/337 (99%)	0.44	9 (2%) 54 53	22, 37, 52, 67	0
2	G	21/82 (25%)	1.50	6 (28%) 0 0	32, 42, 55, 56	0
2	H	20/82 (24%)	2.39	11 (55%) 0 0	41, 53, 62, 63	0
2	I	19/82 (23%)	3.00	14 (73%) 0 0	55, 61, 67, 68	0
All	All	2078/2268 (91%)	0.23	77 (3%) 41 41	9, 31, 51, 68	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	66	ASP	6.0
2	I	64	CYS	5.9
2	I	63	TYR	4.9
2	I	68	PRO	4.6
2	H	58	ASP	4.4
2	I	59	PRO	4.4
2	H	63	TYR	4.3
2	H	62	GLU	4.2
2	G	62	GLU	4.2
2	I	58	ASP	4.1
1	F	68	GLY	3.7
2	I	67	ASN	3.7
2	I	76	TYR	3.6
1	E	333	LYS	3.5
2	H	59	PRO	3.5

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Mol	Chain	Res	Type	RSRZ
2	H	61	GLU	3.4
2	I	75	THR	3.2
2	H	65	LYS	3.2
2	I	65	LYS	3.2
1	E	21	LYS	3.1
1	A	333	LYS	3.1
1	B	333	LYS	2.9
1	F	22	ASP	2.9
1	E	68	GLY	2.9
1	E	93	ILE	2.9
1	B	77	ARG	2.8
1	F	333	LYS	2.8
1	F	93	ILE	2.7
1	E	321	VAL	2.7
2	I	70	THR	2.7
1	E	5	ILE	2.7
1	B	81	LEU	2.7
1	E	328	VAL	2.6
1	F	243	VAL	2.6
2	I	73	CYS	2.6
2	H	71	ASN	2.6
2	G	63	TYR	2.6
1	E	22	ASP	2.6
2	G	58	ASP	2.5
2	I	62	GLU	2.5
1	D	77	ARG	2.5
1	F	124	ASP	2.5
1	B	22	ASP	2.4
1	E	117	ILE	2.4
2	I	60	LEU	2.4
2	G	61	GLU	2.4
1	F	242	LEU	2.4
2	G	66	ASP	2.4
1	C	68	GLY	2.4
2	I	66	ASP	2.4
2	G	59	PRO	2.3
1	E	264	GLU	2.3
1	E	62	THR	2.3
1	C	333	LYS	2.3
1	E	61	GLU	2.3
1	C	139	HIS	2.3
1	F	62	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	333	LYS	2.2
1	D	173	THR	2.2
1	D	242	LEU	2.2
1	E	138	SER	2.2
2	H	64	CYS	2.2
1	E	103	GLU	2.2
1	D	33	THR	2.2
1	E	332	TRP	2.1
2	H	60	LEU	2.1
1	A	61	GLU	2.1
1	C	103	GLU	2.1
1	E	56	ASP	2.1
1	C	85	LYS	2.1
1	C	60	SER	2.1
1	E	30	ILE	2.1
1	E	38	LYS	2.1
1	E	139	HIS	2.0
2	H	76	TYR	2.0
1	C	264	GLU	2.0
1	F	301	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 carbohydrates 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.

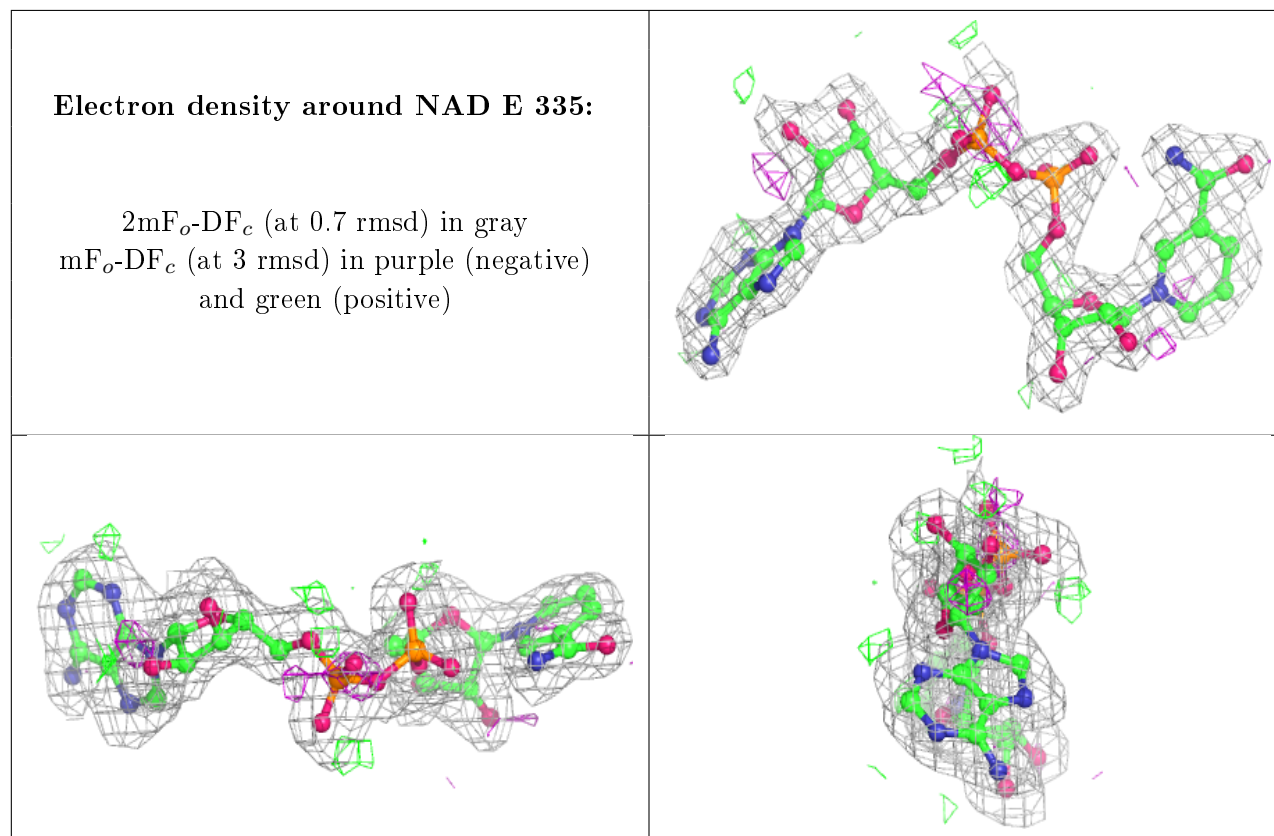
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	SO4	A	337	5/5	0.77	0.25	74,74,75,75	0
5	SO4	A	338	5/5	0.84	0.22	71,71,72,72	0
4	PEG	A	334	7/7	0.91	0.23	27,28,30,32	0

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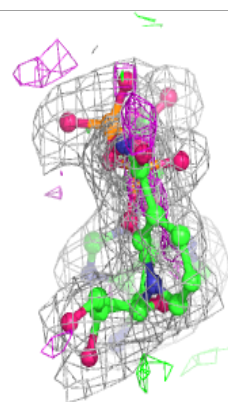
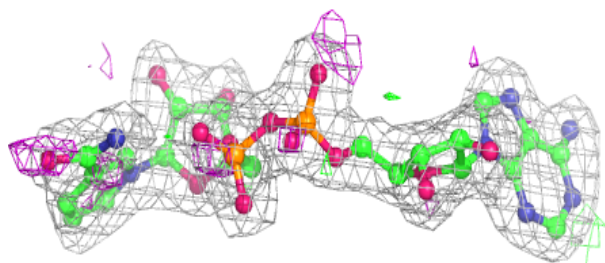
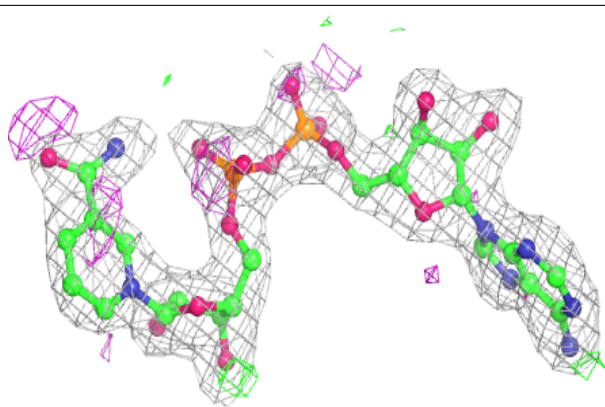
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	PEG	C	334	7/7	0.92	0.18	30,32,33,34	0
4	PEG	A	336	7/7	0.93	0.22	24,28,32,34	0
3	NAD	E	335	44/44	0.95	0.12	21,27,31,31	0
6	EDO	C	336	4/4	0.95	0.17	27,29,29,29	0
3	NAD	F	335	44/44	0.96	0.10	25,28,33,34	0
3	NAD	A	335	44/44	0.97	0.11	13,17,19,19	0
3	NAD	D	335	44/44	0.97	0.08	14,20,22,23	0
3	NAD	C	335	44/44	0.97	0.12	15,21,23,24	0
3	NAD	B	335	44/44	0.97	0.10	19,24,26,26	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.

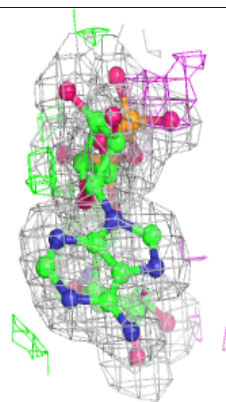
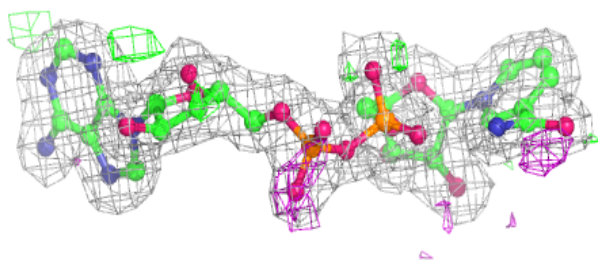
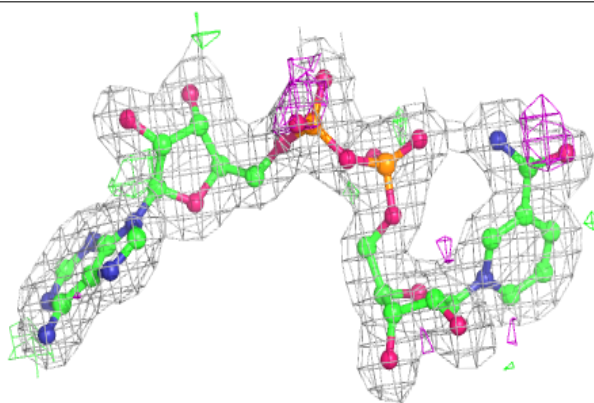


Electron density around NAD F 335:

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

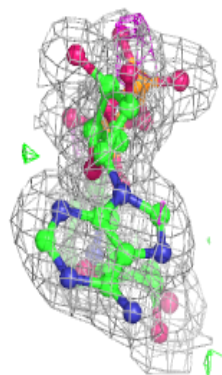
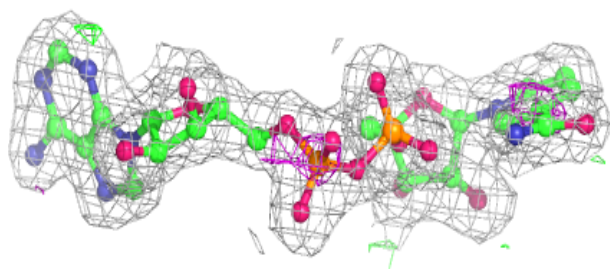
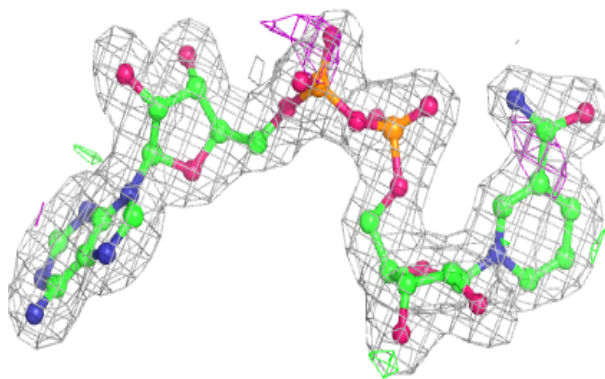
**Electron density around NAD A 335:**

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

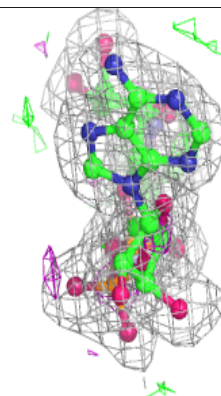
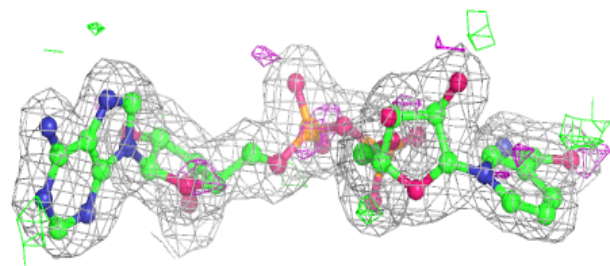
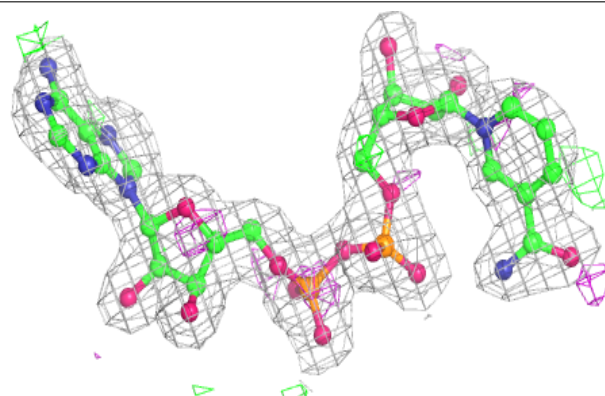


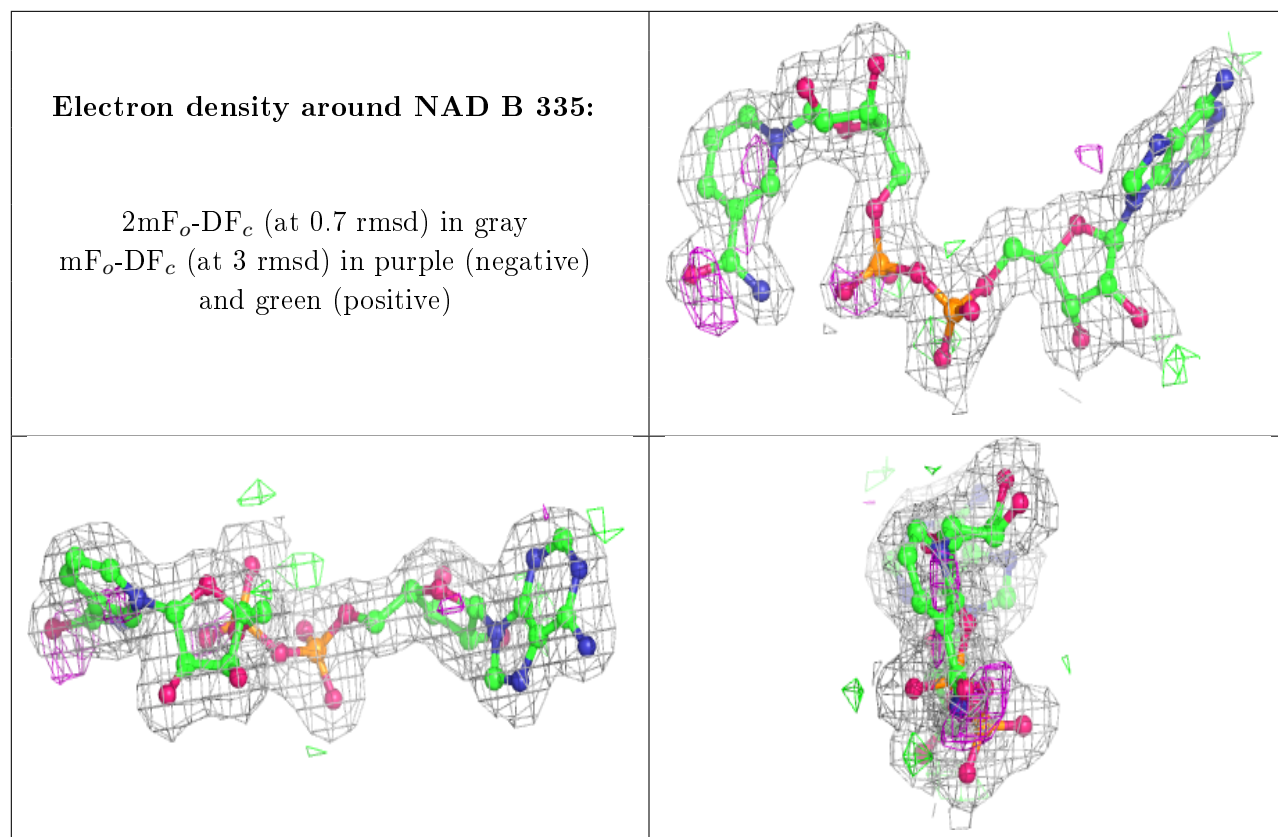
Electron density around NAD D 335:

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

**Electron density around NAD C 335:**

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





6.5 Other polymers ⓘ

There are no such residues in this entry.