Full wwPDB X-ray Structure Validation Report (i)

Oct 16, 2023 – 06:25 AM EDT

PDB ID	:	1NVE
Title	:	Crystal structure of 3-dehydroquinate synthase (DHQS) in complex with ZN2+ and NAD
Authors	:	Nichols, C.E.; Ren, J.; Lamb, H.K.; Hawkins, A.R.; Stammers, D.K.
Deposited on	:	2003-02-03
Resolution	:	2.58 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 (i) 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 (i)) were used in the production of this report:

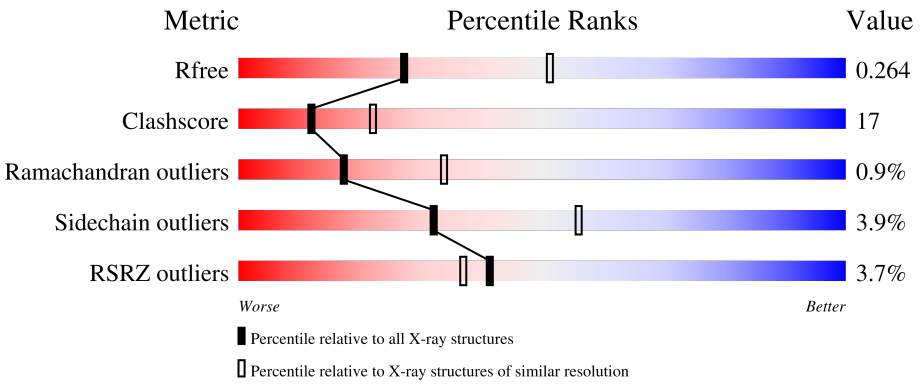
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.58 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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 <=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	393	3% 63% 30% • •
1	B	393	4% 68% 27% • •
1	C	393	4% 66% 28% • •
1	D	393	4% 62% 31% • •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 12493 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 3-DEHYDROQUINATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	378	2887	1834	499	542	12	0	0	0
1	B	383	2933	1862	508	551	12	0	0	0
1	C	377	2886	1831	499	544	12	0	0	0
1	D	381	2915	1851	505	547	12	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

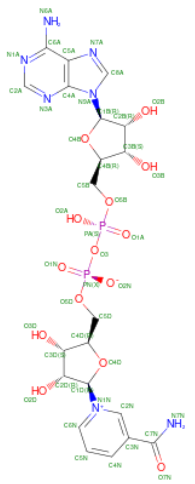
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0
2	D	1	1	1	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
3	A	1	1	1	0	0
3	C	1	1	1	0	0

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).

NAD



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

- Molecule 5 is water.

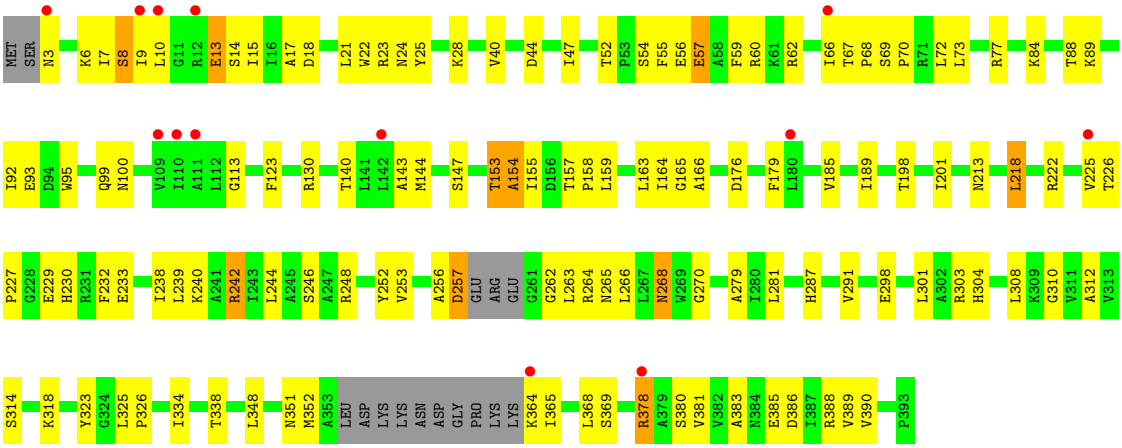
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	173	173	173	0	0
5	B	162	162	162	0	0
5	C	176	176	176	0	0
5	D	179	179	179	0	0

3 Residue-property plots (i)

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.

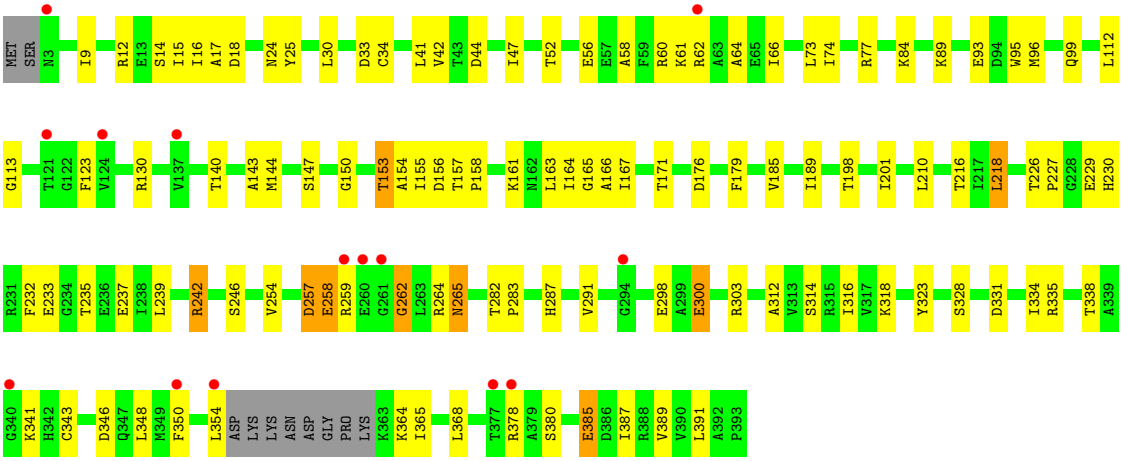
• Molecule 1: 3-DEHYDROQUINATE SYNTHASE

Chain A:



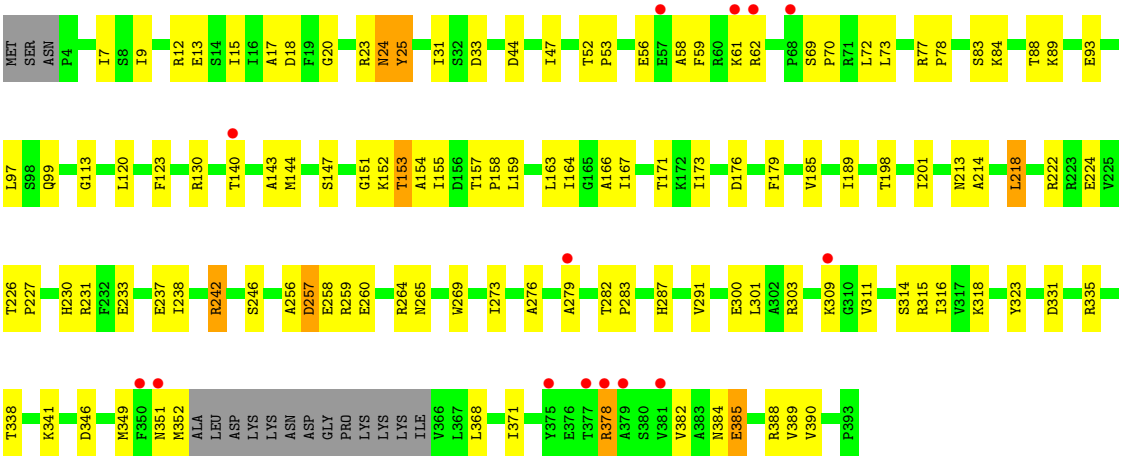
• Molecule 1: 3-DEHYDROQUINATE SYNTHASE

Chain B:



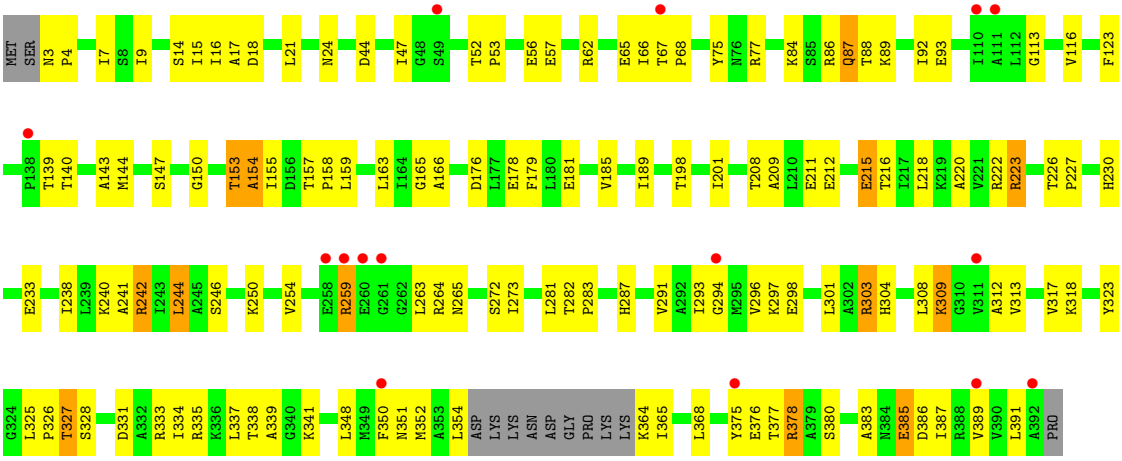
• Molecule 1: 3-DEHYDROQUINATE SYNTHASE

Chain C:



- Molecule 1: 3-DEHYDROQUINATE SYNTHASE

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	83.64Å 70.35Å 144.32Å 90.00° 90.78° 90.00°	Depositor
Resolution (Å)	25.57 – 2.58 25.56 – 2.58	Depositor EDS
% Data completeness (in resolution range)	98.4 (25.57-2.58) 98.5 (25.56-2.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.21 (at 2.57Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.208 , 0.268 0.202 , 0.264	Depositor DCC
R_{free} test set	5121 reflections (9.82%)	wwPDB-VP
Wilson B-factor (Å^2)	51.8	Xtriage
Anisotropy	0.428	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 61.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	12493	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.58 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6086e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.34	0/2935	0.58	0/3980
1	B	0.34	0/2982	0.59	0/4043
1	C	0.33	0/2935	0.59	0/3980
1	D	0.34	0/2963	0.59	0/4019
All	All	0.34	0/11815	0.59	0/16022

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	2887	0	2964	109	0
1	B	2933	0	3014	93	0
1	C	2886	0	2956	90	0
1	D	2915	0	2994	124	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1	0	0	0	0
4	A	44	0	25	0	0
4	B	44	0	25	0	0
4	C	44	0	25	1	0
4	D	44	0	25	1	0
5	A	173	0	0	4	0
5	B	162	0	0	1	0
5	C	176	0	0	5	0
5	D	179	0	0	7	0
All	All	12493	0	12028	409	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 17.

All (409) 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:D:87:GLN:H	1:D:87:GLN:NE2	1.66	0.93
1:D:157:THR:HG22	1:D:159:LEU:H	1.34	0.93
1:D:341:LYS:HA	1:D:341:LYS:HE2	1.52	0.89
1:A:9:ILE:HD11	1:A:15:ILE:HD11	1.55	0.88
1:C:73:LEU:HD13	1:C:99:GLN:HG3	1.56	0.87
1:D:201:ILE:HG22	1:D:368:LEU:HD21	1.54	0.87
1:B:258:GLU:O	1:B:258:GLU:HG3	1.75	0.85
1:C:230:HIS:HB2	1:C:233:GLU:HG3	1.60	0.84
1:D:87:GLN:H	1:D:87:GLN:HE21	1.25	0.83
1:A:89:LYS:O	1:A:93:GLU:HG3	1.79	0.82
1:D:123:PHE:HB2	1:D:153:THR:HG23	1.60	0.82
1:D:84:LYS:HG3	1:D:116:VAL:HG13	1.61	0.81
1:D:226:THR:HG23	1:D:227:PRO:HD2	1.62	0.80
1:D:259:ARG:HA	1:D:259:ARG:HH11	1.47	0.80
1:A:253:VAL:HG13	1:A:263:LEU:HG	1.63	0.80
1:C:201:ILE:HG22	1:C:368:LEU:HD21	1.64	0.79
1:C:143:ALA:HA	1:C:147:SER:OG	1.84	0.78
1:C:84:LYS:HG2	1:C:155:ILE:HA	1.67	0.77
1:D:89:LYS:O	1:D:93:GLU:HG3	1.84	0.77
1:C:61:LYS:HG3	1:C:62:ARG:HD3	1.66	0.76
1:D:157:THR:HB	1:D:163:LEU:HD12	1.66	0.75
1:B:230:HIS:HB2	1:B:233:GLU:HG3	1.68	0.75
1:A:308:LEU:HD11	1:A:312:ALA:HB3	1.70	0.74
1:C:385:GLU:O	1:C:389:VAL:HG23	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:209:ALA:HB1	1:D:242:ARG:NH1	2.03	0.74
1:B:189:ILE:HD12	1:B:323:TYR:O	1.88	0.73
1:D:254:VAL:HG13	1:D:264:ARG:HH12	1.53	0.73
1:D:84:LYS:HD2	1:D:116:VAL:HA	1.70	0.73
1:C:157:THR:HG22	1:C:159:LEU:H	1.52	0.73
1:D:67:THR:HG23	5:D:1724:HOH:O	1.89	0.73
1:A:155:ILE:HD11	1:A:164:ILE:HB	1.69	0.72
1:C:47:ILE:HD13	1:C:113:GLY:HA2	1.70	0.72
1:C:301:LEU:HA	1:C:390:VAL:HG21	1.72	0.72
1:A:73:LEU:HD13	1:A:99:GLN:HG3	1.71	0.71
1:A:201:ILE:HG22	1:A:368:LEU:HD21	1.72	0.71
1:C:153:THR:HG22	1:C:166:ALA:O	1.90	0.71
1:A:238:ILE:HD12	1:A:238:ILE:H	1.56	0.70
1:D:150:GLY:HA2	1:D:254:VAL:HG11	1.73	0.70
1:A:7:ILE:HB	1:A:15:ILE:HB	1.72	0.69
1:A:225:VAL:HG13	1:A:229:GLU:HG3	1.74	0.69
1:A:226:THR:O	1:A:229:GLU:HG2	1.92	0.68
1:C:198:THR:HG21	1:C:246:SER:HA	1.75	0.68
1:D:351:ASN:O	1:D:354:LEU:HB2	1.93	0.68
1:D:387:ILE:O	1:D:391:LEU:HD13	1.94	0.68
1:B:262:GLY:O	1:B:265:ASN:HB2	1.92	0.68
1:D:303:ARG:HH11	1:D:303:ARG:HB3	1.58	0.68
1:B:47:ILE:HD13	1:B:113:GLY:HA2	1.75	0.68
1:B:264:ARG:HD2	1:B:264:ARG:H	1.59	0.67
1:A:143:ALA:HA	1:A:147:SER:OG	1.94	0.67
1:D:348:LEU:O	1:D:352:MET:HG3	1.95	0.67
1:C:273:ILE:HD13	1:C:352:MET:SD	2.35	0.66
1:A:57:GLU:HG2	1:A:60:ARG:HH11	1.60	0.66
1:A:257:ASP:OD2	1:A:264:ARG:HD2	1.96	0.66
1:B:73:LEU:HD13	1:B:99:GLN:HG3	1.76	0.66
1:A:189:ILE:HG22	1:A:325:LEU:HD23	1.78	0.66
1:D:62:ARG:O	1:D:65:GLU:HG2	1.96	0.65
1:A:238:ILE:HD12	1:A:238:ILE:N	2.11	0.65
1:B:84:LYS:HG2	1:B:155:ILE:HA	1.78	0.65
1:B:350:PHE:O	1:B:354:LEU:HD13	1.95	0.65
1:D:242:ARG:CA	1:D:242:ARG:HE	2.10	0.65
1:A:165:GLY:O	1:B:130:ARG:HA	1.96	0.65
1:D:385:GLU:H	1:D:385:GLU:CD	1.98	0.65
1:C:378:ARG:HD3	1:C:378:ARG:N	2.11	0.64
1:B:123:PHE:HB2	1:B:153:THR:HG23	1.79	0.64
1:C:12:ARG:HD2	1:C:13:GLU:O	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:HIS:HB2	1:A:233:GLU:HG3	1.79	0.64
1:A:383:ALA:HB1	1:A:385:GLU:OE1	1.98	0.63
1:B:189:ILE:HD11	1:B:218:LEU:HD11	1.78	0.63
1:B:226:THR:HG23	1:B:227:PRO:HD2	1.80	0.63
1:A:364:LYS:HG2	1:A:381:VAL:HG22	1.80	0.63
1:C:189:ILE:HD12	1:C:323:TYR:O	1.98	0.63
1:D:143:ALA:HA	1:D:147:SER:OG	1.98	0.63
1:B:155:ILE:HD11	1:B:164:ILE:HB	1.80	0.63
1:C:382:VAL:HA	5:C:1626:HOH:O	1.98	0.63
1:B:58:ALA:O	1:B:61:LYS:HG2	1.98	0.63
1:B:52:THR:O	1:B:56:GLU:HG3	1.99	0.62
1:B:226:THR:CG2	1:B:227:PRO:HD2	2.29	0.62
1:D:189:ILE:HD12	1:D:323:TYR:O	1.99	0.62
1:D:293:ILE:O	1:D:297:LYS:HG2	1.99	0.62
1:B:303:ARG:HG2	1:B:303:ARG:HH11	1.63	0.62
1:B:198:THR:HG21	1:B:246:SER:HA	1.81	0.62
1:C:58:ALA:O	1:C:61:LYS:HG2	1.99	0.62
1:D:123:PHE:HA	1:D:153:THR:HG21	1.82	0.62
1:A:157:THR:HG22	1:A:159:LEU:H	1.64	0.62
1:C:23:ARG:HB2	5:C:1613:HOH:O	2.00	0.62
1:C:157:THR:HG23	1:C:158:PRO:HD2	1.81	0.61
1:C:24:ASN:O	1:C:25:TYR:HB2	1.98	0.61
1:B:15:ILE:O	1:B:16:ILE:HD13	2.00	0.61
1:C:20:GLY:HA2	5:C:1613:HOH:O	1.99	0.61
1:C:349:MET:HE2	1:C:384:ASN:O	1.99	0.61
1:A:47:ILE:HD13	1:A:113:GLY:HA2	1.83	0.61
1:B:41:LEU:HB3	1:B:74:ILE:HD13	1.81	0.61
1:A:198:THR:HG21	1:A:246:SER:HA	1.81	0.61
1:D:313:VAL:O	1:D:317:VAL:HG23	1.99	0.60
1:B:237:GLU:CD	1:B:237:GLU:H	2.03	0.60
1:B:312:ALA:O	1:B:316:ILE:HG13	2.02	0.60
1:A:189:ILE:HG22	1:A:325:LEU:CD2	2.31	0.60
1:D:220:ALA:HA	1:D:223:ARG:HG2	1.82	0.60
1:B:335:ARG:CB	1:B:335:ARG:HH11	2.15	0.60
1:D:242:ARG:HE	1:D:242:ARG:N	2.00	0.59
1:D:153:THR:O	1:D:154:ALA:HB2	2.03	0.59
1:B:143:ALA:HA	1:B:147:SER:OG	2.02	0.59
1:C:97:LEU:HD13	1:D:163:LEU:HD13	1.83	0.59
1:C:269:TRP:HH2	1:C:316:ILE:HD13	1.68	0.59
1:C:385:GLU:CD	1:C:385:GLU:H	2.06	0.59
1:D:3:ASN:CG	1:D:4:PRO:HD2	2.23	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:ARG:HD2	1:B:264:ARG:N	2.17	0.58
1:A:189:ILE:HD12	1:A:323:TYR:O	2.03	0.58
1:D:272:SER:HB2	1:D:298:GLU:OE2	2.02	0.58
1:A:123:PHE:HA	1:A:153:THR:HG21	1.85	0.58
1:A:157:THR:HB	1:A:163:LEU:HD12	1.85	0.58
1:A:348:LEU:O	1:A:352:MET:HG3	2.04	0.58
1:D:7:ILE:HB	1:D:15:ILE:HB	1.84	0.58
1:D:378:ARG:HG2	1:D:378:ARG:HH11	1.68	0.58
1:A:9:ILE:HG22	1:A:10:LEU:HD23	1.84	0.58
1:A:6:LYS:O	1:A:7:ILE:HD13	2.03	0.58
1:D:383:ALA:HB3	1:D:386:ASP:OD2	2.04	0.58
1:D:230:HIS:HB2	1:D:233:GLU:HG3	1.86	0.57
1:D:52:THR:O	1:D:56:GLU:HG3	2.03	0.57
1:A:185:VAL:O	1:A:189:ILE:HG12	2.04	0.57
1:A:238:ILE:H	1:A:238:ILE:CD1	2.17	0.57
1:A:298:GLU:HG2	1:A:365:ILE:HG21	1.86	0.57
1:B:331:ASP:HB3	1:B:334:ILE:HD12	1.85	0.57
1:C:231:ARG:HG3	1:C:231:ARG:HH11	1.69	0.57
1:A:365:ILE:N	1:A:365:ILE:HD12	2.19	0.57
1:B:150:GLY:HA2	1:B:254:VAL:HG11	1.87	0.57
1:B:201:ILE:HG22	1:B:368:LEU:HD21	1.85	0.57
1:C:238:ILE:N	1:C:238:ILE:HD12	2.20	0.57
1:B:257:ASP:O	1:B:258:GLU:C	2.43	0.56
1:A:153:THR:O	1:A:154:ALA:HB2	2.05	0.56
1:B:153:THR:HG22	1:B:166:ALA:H	1.70	0.56
1:D:157:THR:HG23	1:D:158:PRO:HD2	1.86	0.56
1:D:281:LEU:HD22	1:D:338:THR:HG21	1.87	0.56
1:A:303:ARG:HE	1:A:310:GLY:CA	2.18	0.55
1:D:189:ILE:HD11	1:D:218:LEU:HD21	1.88	0.55
1:A:308:LEU:HD11	1:A:312:ALA:CB	2.36	0.55
1:C:185:VAL:O	1:C:189:ILE:HG12	2.07	0.55
1:D:185:VAL:O	1:D:189:ILE:HG12	2.07	0.55
1:D:308:LEU:HD11	1:D:312:ALA:HB3	1.88	0.55
1:B:343:CYS:HB3	1:B:348:LEU:HD11	1.88	0.55
1:D:123:PHE:CB	1:D:153:THR:HG23	2.35	0.55
1:D:294:GLY:O	1:D:298:GLU:HG3	2.06	0.55
1:B:60:ARG:HD2	1:B:60:ARG:C	2.27	0.55
1:D:238:ILE:HD12	1:D:238:ILE:N	2.22	0.55
1:A:67:THR:HA	1:A:68:PRO:C	2.27	0.54
1:B:185:VAL:O	1:B:189:ILE:HG12	2.07	0.54
1:D:341:LYS:HA	1:D:341:LYS:CE	2.31	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:PHE:HA	1:B:153:THR:HG21	1.89	0.54
1:C:163:LEU:O	1:C:164:ILE:HD13	2.07	0.54
1:A:263:LEU:HD12	1:A:266:LEU:HD12	1.89	0.54
1:B:335:ARG:HB3	1:B:335:ARG:NH1	2.22	0.54
1:D:351:ASN:HA	1:D:354:LEU:HD13	1.90	0.54
1:A:66:ILE:N	1:A:66:ILE:HD12	2.23	0.54
1:A:3:ASN:N	5:A:678:HOH:O	2.40	0.53
1:A:73:LEU:CD1	1:A:99:GLN:HG3	2.38	0.53
1:A:279:ALA:HB3	1:A:351:ASN:HD22	1.73	0.53
1:C:89:LYS:HE3	1:C:123:PHE:CE2	2.43	0.53
1:B:364:LYS:HA	1:B:380:SER:O	2.08	0.53
1:B:58:ALA:HA	1:B:61:LYS:HD3	1.89	0.53
1:D:68:PRO:HA	5:D:1724:HOH:O	2.08	0.53
1:D:308:LEU:HD11	1:D:312:ALA:CB	2.38	0.53
1:B:198:THR:HG22	5:B:683:HOH:O	2.09	0.53
1:B:265:ASN:ND2	1:B:378:ARG:NH1	2.56	0.53
1:D:230:HIS:HB2	1:D:233:GLU:CG	2.38	0.53
1:A:287:HIS:O	1:A:291:VAL:HG23	2.09	0.53
1:D:215:GLU:HG3	1:D:216:THR:N	2.24	0.53
1:C:257:ASP:OD2	1:C:264:ARG:HG3	2.08	0.53
1:D:47:ILE:HD13	1:D:113:GLY:HA2	1.91	0.53
1:A:44:ASP:HA	1:A:77:ARG:O	2.09	0.52
1:A:52:THR:O	1:A:56:GLU:HG3	2.09	0.52
1:D:364:LYS:C	1:D:365:ILE:HD12	2.30	0.52
1:B:230:HIS:HB2	1:B:233:GLU:CG	2.39	0.52
1:A:364:LYS:HA	1:A:380:SER:O	2.10	0.52
1:D:189:ILE:HG23	1:D:323:TYR:O	2.09	0.52
1:D:350:PHE:O	1:D:354:LEU:HD13	2.10	0.52
1:C:237:GLU:CD	1:C:237:GLU:H	2.13	0.52
1:C:9:ILE:HD11	1:C:15:ILE:HD11	1.91	0.51
1:A:213:ASN:OD1	1:A:238:ILE:HG23	2.10	0.51
1:B:387:ILE:O	1:B:391:LEU:HD23	2.10	0.51
1:A:9:ILE:HG22	1:A:10:LEU:CD2	2.40	0.51
1:A:88:THR:O	1:A:92:ILE:HG13	2.11	0.51
1:D:287:HIS:O	1:D:291:VAL:HG23	2.11	0.51
1:D:87:GLN:NE2	1:D:87:GLN:N	2.47	0.51
1:B:303:ARG:HG2	1:B:303:ARG:NH1	2.26	0.51
1:B:73:LEU:CD1	1:B:99:GLN:HG3	2.40	0.51
1:C:287:HIS:O	1:C:291:VAL:HG23	2.11	0.51
1:B:153:THR:O	1:B:154:ALA:HB2	2.11	0.50
1:B:155:ILE:C	1:B:155:ILE:HD12	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:218:LEU:HD12	1:C:222:ARG:CZ	2.41	0.50
1:D:242:ARG:HE	1:D:242:ARG:HA	1.76	0.50
1:A:57:GLU:HG2	1:A:60:ARG:NH1	2.25	0.50
1:B:41:LEU:HD11	1:B:112:LEU:HB3	1.93	0.50
1:C:349:MET:HE2	1:C:388:ARG:HG3	1.92	0.50
1:B:385:GLU:H	1:B:385:GLU:CD	2.15	0.50
1:B:64:ALA:C	1:B:66:ILE:H	2.14	0.50
1:D:65:GLU:O	1:D:66:ILE:HD13	2.11	0.49
1:A:9:ILE:CD1	1:A:15:ILE:HD11	2.35	0.49
1:A:386:ASP:O	1:A:389:VAL:HG22	2.12	0.49
1:B:44:ASP:HA	1:B:77:ARG:O	2.12	0.49
1:B:140:THR:O	1:B:144:MET:HG3	2.11	0.49
1:C:303:ARG:HH11	1:C:303:ARG:HG2	1.77	0.49
1:D:52:THR:N	1:D:53:PRO:CD	2.74	0.49
1:A:240:LYS:O	1:A:244:LEU:HB2	2.12	0.49
1:C:230:HIS:HB2	1:C:233:GLU:CG	2.39	0.49
1:D:335:ARG:HD2	5:D:1629:HOH:O	2.12	0.49
1:C:231:ARG:HG3	1:C:231:ARG:NH1	2.28	0.49
1:C:44:ASP:HA	1:C:77:ARG:O	2.12	0.49
1:C:153:THR:O	1:C:154:ALA:HB2	2.12	0.48
1:C:18:ASP:O	1:C:176:ASP:HA	2.13	0.48
1:C:58:ALA:HA	1:C:61:LYS:HE2	1.96	0.48
1:D:226:THR:CG2	1:D:227:PRO:HD2	2.40	0.48
1:A:7:ILE:HG22	1:A:8:SER:N	2.28	0.48
1:A:153:THR:HG22	1:A:166:ALA:O	2.12	0.48
1:C:269:TRP:CH2	1:C:316:ILE:HD13	2.49	0.48
1:A:130:ARG:HD3	1:B:167:ILE:HD12	1.94	0.48
1:D:3:ASN:N	5:D:1639:HOH:O	2.45	0.48
1:D:88:THR:O	1:D:92:ILE:HG13	2.14	0.48
1:A:7:ILE:O	1:A:13:GLU:HA	2.13	0.48
1:A:189:ILE:O	1:A:325:LEU:HD21	2.14	0.48
1:D:17:ALA:O	1:D:18:ASP:HB2	2.14	0.48
1:A:59:PHE:CE2	1:A:72:LEU:HB2	2.49	0.48
1:B:338:THR:HB	1:B:341:LYS:HB2	1.95	0.48
1:C:140:THR:O	1:C:144:MET:HG3	2.14	0.48
1:D:44:ASP:HA	1:D:77:ARG:O	2.14	0.48
1:A:9:ILE:HD11	1:A:15:ILE:CD1	2.35	0.47
1:B:156:ASP:OD1	1:B:161:LYS:HA	2.14	0.47
1:B:287:HIS:O	1:B:291:VAL:HG23	2.13	0.47
1:D:230:HIS:H	1:D:233:GLU:CD	2.17	0.47
1:B:258:GLU:O	1:B:258:GLU:CG	2.54	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:LYS:HE3	1:C:123:PHE:CD2	2.49	0.47
1:C:130:ARG:HA	1:D:165:GLY:O	2.14	0.47
1:A:225:VAL:CG1	1:A:229:GLU:HG3	2.42	0.47
1:B:12:ARG:HH22	1:B:171:THR:HA	1.79	0.47
1:B:12:ARG:NH2	1:B:171:THR:HA	2.29	0.47
1:B:314:SER:O	1:B:318:LYS:HG2	2.14	0.47
1:C:300:GLU:HB3	1:C:390:VAL:HG13	1.95	0.47
1:D:3:ASN:OD1	1:D:4:PRO:HD2	2.13	0.47
1:A:123:PHE:N	1:A:153:THR:OG1	2.47	0.47
1:B:258:GLU:O	1:B:259:ARG:HB3	2.15	0.47
1:D:87:GLN:HE21	1:D:87:GLN:N	2.04	0.47
1:A:21:LEU:HD22	1:A:25:TYR:CD1	2.49	0.47
1:A:326:PRO:HG3	1:A:334:ILE:HD11	1.96	0.47
1:B:300:GLU:OE1	1:B:300:GLU:HA	2.15	0.47
1:A:230:HIS:HB2	1:A:233:GLU:CG	2.45	0.47
1:A:378:ARG:HH11	1:A:378:ARG:HG2	1.80	0.47
1:D:208:THR:HA	1:D:211:GLU:OE1	2.14	0.47
1:A:130:ARG:HA	1:B:165:GLY:O	2.15	0.47
1:B:265:ASN:ND2	1:B:378:ARG:HH12	2.13	0.47
1:C:314:SER:O	1:C:318:LYS:HG2	2.15	0.47
1:B:41:LEU:HD23	1:B:74:ILE:CD1	2.45	0.46
1:A:84:LYS:HG2	1:A:155:ILE:HA	1.97	0.46
1:C:279:ALA:HB3	1:C:351:ASN:HD22	1.81	0.46
1:D:140:THR:O	1:D:144:MET:HG3	2.16	0.46
1:A:388:ARG:HD3	5:A:715:HOH:O	2.14	0.46
1:C:9:ILE:HG21	5:C:1640:HOH:O	2.15	0.46
1:D:338:THR:O	1:D:339:ALA:C	2.53	0.46
1:A:130:ARG:CD	1:B:167:ILE:HD12	2.45	0.46
1:C:311:VAL:O	1:C:315:ARG:HG3	2.15	0.46
1:A:155:ILE:C	1:A:155:ILE:HD12	2.35	0.46
1:D:375:TYR:CD2	1:D:376:GLU:HG3	2.51	0.46
1:A:301:LEU:O	1:A:304:HIS:HB3	2.16	0.46
1:C:52:THR:N	1:C:53:PRO:CD	2.79	0.46
1:D:241:ALA:HB1	1:D:242:ARG:HH21	1.80	0.45
1:A:303:ARG:HE	1:A:310:GLY:HA3	1.81	0.45
1:D:254:VAL:HG13	1:D:264:ARG:NH1	2.27	0.45
1:D:272:SER:CB	1:D:365:ILE:HG13	2.46	0.45
1:A:252:TYR:HE2	1:A:263:LEU:HD21	1.81	0.45
1:C:238:ILE:N	1:C:238:ILE:CD1	2.79	0.45
1:C:12:ARG:NH1	1:C:171:THR:O	2.50	0.45
1:B:157:THR:HB	1:B:163:LEU:HD12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:17:ALA:CB	1:D:244:LEU:HD11	2.46	0.45
1:A:314:SER:O	1:A:318:LYS:HG2	2.16	0.45
1:C:7:ILE:HB	1:C:15:ILE:HB	1.97	0.45
1:D:155:ILE:C	1:D:155:ILE:HD12	2.36	0.45
1:A:252:TYR:CE2	1:A:263:LEU:HD21	2.51	0.45
1:D:263:LEU:HD13	1:D:377:THR:O	2.16	0.45
1:D:296:VAL:HG21	1:D:327:THR:HG22	1.98	0.45
1:D:208:THR:O	1:D:212:GLU:HG2	2.17	0.45
1:A:28:LYS:HB2	1:A:62:ARG:NE	2.31	0.45
1:B:385:GLU:O	1:B:389:VAL:HG23	2.16	0.45
1:B:226:THR:HB	1:B:229:GLU:CD	2.37	0.45
1:C:58:ALA:C	1:C:61:LYS:HG2	2.37	0.45
1:B:15:ILE:C	1:B:16:ILE:HD13	2.37	0.44
1:C:226:THR:CG2	1:C:227:PRO:HD2	2.47	0.44
1:B:61:LYS:NZ	1:B:62:ARG:CZ	2.80	0.44
1:B:335:ARG:CB	1:B:335:ARG:NH1	2.78	0.44
1:D:157:THR:HG23	1:D:158:PRO:CD	2.47	0.44
1:C:276:ALA:HB2	1:C:352:MET:HG3	1.99	0.44
1:D:176:ASP:HB3	1:D:179:PHE:CD2	2.53	0.44
1:A:22:TRP:CZ2	1:A:55:PHE:HB2	2.52	0.44
1:A:388:ARG:C	1:A:390:VAL:H	2.19	0.44
1:B:41:LEU:HD12	1:B:42:VAL:N	2.31	0.44
1:B:123:PHE:CB	1:B:153:THR:HG23	2.45	0.44
1:C:78:PRO:HG2	1:C:83:SER:OG	2.18	0.44
1:D:17:ALA:HB3	1:D:244:LEU:HD21	1.99	0.44
1:D:240:LYS:O	1:D:244:LEU:HB2	2.17	0.44
1:A:232:PHE:CD1	1:A:239:LEU:HD22	2.52	0.44
1:D:326:PRO:HD3	5:D:1641:HOH:O	2.17	0.44
1:D:333:ARG:HD2	1:D:337:LEU:HD11	1.99	0.44
1:C:151:GLY:O	1:C:167:ILE:HA	2.18	0.44
1:C:338:THR:HB	1:C:341:LYS:HB2	1.99	0.44
1:B:89:LYS:O	1:B:93:GLU:HG3	2.18	0.44
1:D:259:ARG:HH11	1:D:259:ARG:CA	2.25	0.44
1:C:214:ALA:HA	1:C:323:TYR:CZ	2.53	0.44
1:D:7:ILE:HG22	1:D:9:ILE:CD1	2.47	0.44
1:D:198:THR:HG21	1:D:246:SER:HA	1.99	0.44
1:A:17:ALA:O	1:A:18:ASP:HB2	2.16	0.43
1:B:89:LYS:HE3	1:B:123:PHE:CE2	2.53	0.43
1:D:242:ARG:CA	1:D:242:ARG:NE	2.80	0.43
1:D:293:ILE:HD11	1:D:334:ILE:CD1	2.48	0.43
1:A:364:LYS:C	1:A:365:ILE:HD12	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:ALA:O	1:B:18:ASP:HB2	2.18	0.43
1:C:256:ALA:O	1:C:258:GLU:N	2.52	0.43
1:D:386:ASP:O	1:D:389:VAL:HG22	2.18	0.43
1:A:218:LEU:HD22	1:A:222:ARG:HD2	2.00	0.43
1:C:226:THR:HG23	1:C:227:PRO:HD2	2.00	0.43
1:D:364:LYS:HA	1:D:380:SER:O	2.18	0.43
1:C:157:THR:HB	1:C:163:LEU:HD12	1.99	0.43
1:D:16:ILE:HG22	1:D:21:LEU:HD11	2.00	0.43
1:D:318:LYS:NZ	5:D:1717:HOH:O	2.51	0.43
1:C:331:ASP:O	1:C:335:ARG:HG3	2.19	0.43
1:C:93:GLU:OE1	1:D:86:ARG:NH2	2.44	0.43
1:A:18:ASP:O	1:A:176:ASP:HA	2.19	0.43
1:B:9:ILE:HD11	1:B:15:ILE:HD11	1.99	0.43
1:C:59:PHE:CE2	1:C:72:LEU:HB2	2.54	0.43
1:C:258:GLU:HG3	1:C:259:ARG:N	2.32	0.43
1:C:309:LYS:HB2	5:C:1704:HOH:O	2.19	0.43
1:B:18:ASP:O	1:B:176:ASP:HA	2.18	0.43
1:B:153:THR:HG22	1:B:166:ALA:N	2.34	0.43
1:B:335:ARG:HH11	1:B:335:ARG:HB2	1.82	0.43
1:A:123:PHE:HB2	1:A:153:THR:HG23	2.01	0.42
1:B:30:LEU:O	1:B:34:CYS:N	2.49	0.42
1:B:58:ALA:O	1:B:62:ARG:HG2	2.19	0.42
1:B:298:GLU:HG2	1:B:365:ILE:HG21	1.99	0.42
1:C:316:ILE:HG13	1:C:371:ILE:HD11	2.01	0.42
1:A:240:LYS:HD2	5:A:758:HOH:O	2.20	0.42
1:A:268:ASN:HD22	1:A:268:ASN:HA	1.59	0.42
1:B:95:TRP:CE3	1:B:96:MET:HG3	2.55	0.42
1:C:15:ILE:HG12	1:C:173:ILE:HB	2.02	0.42
1:B:60:ARG:HD2	1:B:60:ARG:O	2.19	0.42
1:D:212:GLU:HG2	1:D:212:GLU:H	1.67	0.42
1:A:155:ILE:CD1	1:A:164:ILE:HB	2.44	0.42
1:A:262:GLY:O	1:A:265:ASN:HB2	2.20	0.42
1:C:282:THR:OG1	1:C:283:PRO:HA	2.19	0.42
1:B:216:THR:CG2	1:B:235:THR:HG21	2.48	0.42
1:B:232:PHE:CD1	1:B:239:LEU:HD22	2.55	0.42
1:D:218:LEU:HD13	1:D:222:ARG:CZ	2.50	0.42
1:D:282:THR:OG1	1:D:283:PRO:HA	2.20	0.42
1:A:44:ASP:C	1:A:44:ASP:OD1	2.57	0.42
1:B:176:ASP:HB3	1:B:179:PHE:CD2	2.55	0.42
1:D:123:PHE:HA	1:D:153:THR:CG2	2.49	0.42
1:D:123:PHE:CA	1:D:153:THR:CG2	2.98	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:241:ALA:HB1	1:D:242:ARG:NH2	2.35	0.42
1:D:301:LEU:O	1:D:304:HIS:HB3	2.20	0.42
1:A:253:VAL:HG13	1:A:263:LEU:CG	2.41	0.42
1:B:157:THR:CG2	1:B:158:PRO:HD2	2.49	0.42
1:C:213:ASN:HD22	1:C:242:ARG:HD3	1.84	0.42
1:A:176:ASP:HB3	1:A:179:PHE:CD2	2.54	0.42
1:B:264:ARG:H	1:B:264:ARG:CD	2.29	0.42
1:D:309:LYS:HE3	1:D:309:LYS:HA	2.02	0.42
1:A:385:GLU:O	1:A:389:VAL:HG13	2.20	0.42
1:C:152:LYS:HB2	4:C:402:NAD:H72N	1.84	0.42
1:D:153:THR:O	1:D:154:ALA:CB	2.68	0.42
1:A:157:THR:HG23	1:A:158:PRO:HD2	2.02	0.41
1:B:282:THR:OG1	1:B:283:PRO:HA	2.20	0.41
1:C:84:LYS:CG	1:C:155:ILE:HA	2.42	0.41
1:A:54:SER:OG	1:A:55:PHE:N	2.53	0.41
1:A:242:ARG:HE	1:A:242:ARG:HA	1.85	0.41
1:B:41:LEU:HD11	1:B:112:LEU:CB	2.50	0.41
1:C:176:ASP:HB3	1:C:179:PHE:CD2	2.55	0.41
1:A:303:ARG:HE	1:A:310:GLY:HA2	1.84	0.41
1:C:61:LYS:HG3	1:C:62:ARG:CD	2.45	0.41
1:D:238:ILE:N	1:D:238:ILE:CD1	2.84	0.41
1:A:23:ARG:HG2	1:A:23:ARG:HH11	1.84	0.41
1:A:281:LEU:HD22	1:A:338:THR:HG21	2.02	0.41
1:C:31:ILE:HD12	1:C:62:ARG:HB3	2.01	0.41
1:D:328:SER:O	1:D:331:ASP:HB2	2.21	0.41
1:A:40:VAL:HG13	1:A:95:TRP:HZ3	1.85	0.41
1:B:210:LEU:HD23	1:B:242:ARG:CG	2.51	0.41
1:C:143:ALA:HA	1:C:147:SER:HG	1.81	0.41
1:D:218:LEU:HA	1:D:218:LEU:HD23	1.85	0.41
1:A:7:ILE:HD12	1:A:248:ARG:HE	1.85	0.41
1:C:88:THR:HG22	1:C:120:LEU:HD11	2.03	0.41
1:C:301:LEU:CA	1:C:390:VAL:HG21	2.47	0.41
1:D:139:THR:HB	4:D:1403:NAD:H61A	1.86	0.41
1:A:140:THR:O	1:A:144:MET:HG3	2.21	0.41
1:A:155:ILE:HD12	1:A:155:ILE:O	2.21	0.41
1:A:364:LYS:HE3	1:A:381:VAL:HG23	2.02	0.41
1:C:52:THR:O	1:C:56:GLU:HG3	2.20	0.41
1:D:325:LEU:HA	5:D:1641:HOH:O	2.20	0.41
1:C:157:THR:CG2	1:C:158:PRO:HD2	2.50	0.40
1:D:385:GLU:CD	1:D:385:GLU:N	2.70	0.40
1:A:69:SER:HA	1:A:70:PRO:HD3	1.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:17:ALA:O	1:C:18:ASP:HB2	2.21	0.40
1:C:58:ALA:HA	1:C:61:LYS:HG2	2.02	0.40
1:D:273:ILE:HD12	1:D:391:LEU:HD11	2.03	0.40
1:D:75:TYR:CD1	1:D:75:TYR:C	2.95	0.40
1:D:351:ASN:HA	1:D:354:LEU:HD22	2.04	0.40
1:D:378:ARG:HH11	1:D:378:ARG:CG	2.34	0.40
1:A:8:SER:HB3	1:A:13:GLU:N	2.37	0.40
1:C:69:SER:HA	1:C:70:PRO:HD3	1.88	0.40
1:A:264:ARG:HD3	5:A:635:HOH:O	2.21	0.40
1:A:270:GLY:N	1:A:298:GLU:OE1	2.39	0.40
1:C:24:ASN:O	1:C:24:ASN:ND2	2.48	0.40
1:D:153:THR:HB	1:D:166:ALA:O	2.20	0.40
1:D:178:GLU:O	1:D:181:GLU:HG3	2.22	0.40
1:D:189:ILE:CD1	1:D:218:LEU:HD21	2.50	0.40
1:D:352:MET:C	1:D:354:LEU:H	2.25	0.40
1:D:365:ILE:HD12	1:D:365:ILE:N	2.36	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	372/393 (95%)	351 (94%)	16 (4%)	5 (1%)	12 23
1	B	379/393 (96%)	355 (94%)	20 (5%)	4 (1%)	14 28
1	C	373/393 (95%)	353 (95%)	17 (5%)	3 (1%)	19 37
1	D	377/393 (96%)	361 (96%)	15 (4%)	1 (0%)	41 62
All	All	1501/1572 (96%)	1420 (95%)	68 (4%)	13 (1%)	17 34

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	256	ALA
1	C	257	ASP
1	A	13	GLU
1	B	25	TYR
1	C	25	TYR
1	A	14	SER
1	B	258	GLU
1	A	154	ALA
1	B	14	SER
1	C	260	GLU
1	D	154	ALA
1	B	262	GLY
1	A	227	PRO

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	311/326 (95%)	300 (96%)	11 (4%)	36	59
1	B	316/326 (97%)	305 (96%)	11 (4%)	36	59
1	C	311/326 (95%)	301 (97%)	10 (3%)	39	63
1	D	314/326 (96%)	297 (95%)	17 (5%)	22	42
All	All	1252/1304 (96%)	1203 (96%)	49 (4%)	32	56

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	SER
1	A	24	ASN
1	A	57	GLU
1	A	100	ASN
1	A	153	THR
1	A	218	LEU
1	A	242	ARG
1	A	257	ASP
1	A	268	ASN

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Mol	Chain	Res	Type
1	A	369	SER
1	A	378	ARG
1	B	24	ASN
1	B	33	ASP
1	B	153	THR
1	B	218	LEU
1	B	242	ARG
1	B	257	ASP
1	B	265	ASN
1	B	300	GLU
1	B	328	SER
1	B	346	ASP
1	B	385	GLU
1	C	24	ASN
1	C	33	ASP
1	C	153	THR
1	C	218	LEU
1	C	224	GLU
1	C	242	ARG
1	C	265	ASN
1	C	346	ASP
1	C	378	ARG
1	C	385	GLU
1	D	14	SER
1	D	24	ASN
1	D	57	GLU
1	D	87	GLN
1	D	153	THR
1	D	215	GLU
1	D	223	ARG
1	D	242	ARG
1	D	244	LEU
1	D	250	LYS
1	D	259	ARG
1	D	265	ASN
1	D	303	ARG
1	D	309	LYS
1	D	327	THR
1	D	378	ARG
1	D	385	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	76	ASN
1	A	100	ASN
1	A	268	ASN
1	B	24	ASN
1	B	265	ASN
1	B	268	ASN
1	C	87	GLN
1	C	213	ASN
1	C	271	HIS
1	D	24	ASN
1	D	87	GLN
1	D	268	ASN
1	D	271	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	NAD	A	400	-	42,48,48	2.34	11 (26%)	50,73,73	1.73	8 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	NAD	D	1403	-	42,48,48	2.50	11 (26%)	50,73,73	1.71	8 (16%)
4	NAD	B	401	-	42,48,48	2.46	12 (28%)	50,73,73	1.77	10 (20%)
4	NAD	C	402	-	42,48,48	2.43	11 (26%)	50,73,73	1.74	9 (18%)

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	NAD	A	400	-	-	8/26/62/62	0/5/5/5
4	NAD	D	1403	-	-	6/26/62/62	0/5/5/5
4	NAD	B	401	-	-	2/26/62/62	0/5/5/5
4	NAD	C	402	-	-	1/26/62/62	0/5/5/5

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	400	NAD	O3B-C3B	-8.62	1.22	1.43
4	C	402	NAD	O3B-C3B	-8.57	1.22	1.43
4	D	1403	NAD	O3B-C3B	-8.50	1.23	1.43
4	B	401	NAD	O3B-C3B	-8.46	1.23	1.43
4	D	1403	NAD	O4B-C1B	5.66	1.49	1.41
4	B	401	NAD	C2B-C1B	-5.31	1.45	1.53
4	A	400	NAD	O4B-C1B	5.21	1.48	1.41
4	C	402	NAD	O4B-C1B	4.92	1.47	1.41
4	C	402	NAD	O4D-C1D	4.55	1.47	1.41
4	D	1403	NAD	O4D-C1D	4.45	1.47	1.41
4	D	1403	NAD	C6N-N1N	4.36	1.46	1.35
4	D	1403	NAD	C4A-N3A	4.27	1.41	1.35
4	B	401	NAD	C6N-N1N	4.26	1.45	1.35
4	B	401	NAD	O4D-C1D	4.24	1.47	1.41
4	D	1403	NAD	C2B-C1B	-4.19	1.47	1.53
4	A	400	NAD	C4A-N3A	4.12	1.41	1.35
4	B	401	NAD	C4A-N3A	4.10	1.41	1.35
4	B	401	NAD	C4N-C3N	4.01	1.46	1.39
4	C	402	NAD	C6N-N1N	3.91	1.45	1.35
4	C	402	NAD	C2N-N1N	3.90	1.39	1.35
4	C	402	NAD	C3N-C7N	3.87	1.56	1.50
4	C	402	NAD	C2B-C1B	-3.81	1.48	1.53
4	B	401	NAD	C2N-N1N	3.80	1.39	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	400	NAD	C6N-N1N	3.79	1.44	1.35
4	C	402	NAD	C4A-N3A	3.69	1.40	1.35
4	C	402	NAD	C4N-C3N	3.67	1.45	1.39
4	D	1403	NAD	C4N-C3N	3.64	1.45	1.39
4	A	400	NAD	O4D-C1D	3.63	1.46	1.41
4	B	401	NAD	O4B-C1B	3.60	1.46	1.41
4	D	1403	NAD	C2N-N1N	3.55	1.39	1.35
4	B	401	NAD	C3N-C7N	3.51	1.55	1.50
4	A	400	NAD	C4N-C3N	3.45	1.45	1.39
4	D	1403	NAD	C3N-C7N	3.39	1.55	1.50
4	A	400	NAD	C2B-C1B	-3.26	1.48	1.53
4	A	400	NAD	C2N-N1N	3.10	1.38	1.35
4	A	400	NAD	C3N-C7N	3.08	1.55	1.50
4	C	402	NAD	C2A-N1A	2.83	1.39	1.33
4	A	400	NAD	C2A-N1A	2.57	1.38	1.33
4	D	1403	NAD	C2A-N1A	2.56	1.38	1.33
4	D	1403	NAD	C5N-C4N	2.52	1.44	1.38
4	B	401	NAD	C2A-N1A	2.52	1.38	1.33
4	B	401	NAD	C5N-C4N	2.35	1.43	1.38
4	B	401	NAD	C7N-N7N	2.31	1.37	1.33
4	A	400	NAD	C5N-C4N	2.23	1.43	1.38
4	C	402	NAD	C5N-C4N	2.17	1.43	1.38

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	401	NAD	C5N-C4N-C3N	-6.73	112.38	120.34
4	A	400	NAD	C5N-C4N-C3N	-6.11	113.11	120.34
4	C	402	NAD	C5N-C4N-C3N	-5.97	113.28	120.34
4	D	1403	NAD	C5N-C4N-C3N	-5.75	113.53	120.34
4	C	402	NAD	C6N-C5N-C4N	5.03	126.75	119.44
4	A	400	NAD	C6N-C5N-C4N	4.82	126.44	119.44
4	D	1403	NAD	C6N-C5N-C4N	4.73	126.32	119.44
4	B	401	NAD	C6N-C5N-C4N	4.64	126.19	119.44
4	D	1403	NAD	C2N-C3N-C4N	4.05	122.85	118.26
4	C	402	NAD	C2N-C3N-C4N	3.99	122.78	118.26
4	A	400	NAD	C2N-C3N-C4N	3.89	122.67	118.26
4	A	400	NAD	C5N-C6N-N1N	-3.87	114.85	120.40
4	D	1403	NAD	C5N-C6N-N1N	-3.62	115.22	120.40
4	B	401	NAD	C2N-C3N-C4N	3.57	122.31	118.26
4	B	401	NAD	C5N-C6N-N1N	-3.52	115.35	120.40
4	A	400	NAD	C2N-N1N-C1D	-3.36	111.66	119.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	402	NAD	C5N-C6N-N1N	-3.30	115.67	120.40
4	D	1403	NAD	C2N-N1N-C1D	-3.23	111.93	119.14
4	C	402	NAD	C2N-N1N-C1D	-3.13	112.16	119.14
4	C	402	NAD	C5A-C6A-N6A	2.90	124.76	120.35
4	B	401	NAD	C2N-N1N-C1D	-2.87	112.74	119.14
4	B	401	NAD	O4D-C1D-C2D	-2.84	102.78	106.93
4	A	400	NAD	C5A-C6A-N6A	2.80	124.60	120.35
4	B	401	NAD	C5A-C6A-N6A	2.77	124.57	120.35
4	D	1403	NAD	C5A-C6A-N6A	2.71	124.48	120.35
4	C	402	NAD	C4N-C3N-C7N	-2.30	114.88	121.04
4	C	402	NAD	O4D-C1D-C2D	-2.30	103.57	106.93
4	D	1403	NAD	C4N-C3N-C7N	-2.25	115.03	121.04
4	A	400	NAD	C4N-C3N-C7N	-2.20	115.16	121.04
4	B	401	NAD	O2N-PN-O1N	2.19	123.07	112.24
4	B	401	NAD	C4N-C3N-C7N	-2.18	115.19	121.04
4	D	1403	NAD	O2N-PN-O1N	2.14	122.81	112.24
4	C	402	NAD	O2N-PN-O1N	2.11	122.66	112.24
4	A	400	NAD	O2N-PN-O1N	2.10	122.61	112.24
4	B	401	NAD	C3D-C2D-C1D	2.03	104.03	100.98

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	400	NAD	C5D-O5D-PN-O2N
4	A	400	NAD	O4D-C1D-N1N-C2N
4	A	400	NAD	O4D-C1D-N1N-C6N
4	A	400	NAD	C2D-C1D-N1N-C2N
4	A	400	NAD	C2D-C1D-N1N-C6N
4	C	402	NAD	O4D-C1D-N1N-C6N
4	D	1403	NAD	O4D-C1D-N1N-C2N
4	D	1403	NAD	O4D-C1D-N1N-C6N
4	D	1403	NAD	C2D-C1D-N1N-C2N
4	A	400	NAD	PA-O3-PN-O5D
4	A	400	NAD	C5D-O5D-PN-O3
4	D	1403	NAD	PN-O3-PA-O1A
4	B	401	NAD	PN-O3-PA-O1A
4	D	1403	NAD	C2D-C1D-N1N-C6N
4	B	401	NAD	PN-O3-PA-O2A
4	D	1403	NAD	PN-O3-PA-O2A
4	A	400	NAD	C5D-O5D-PN-O1N

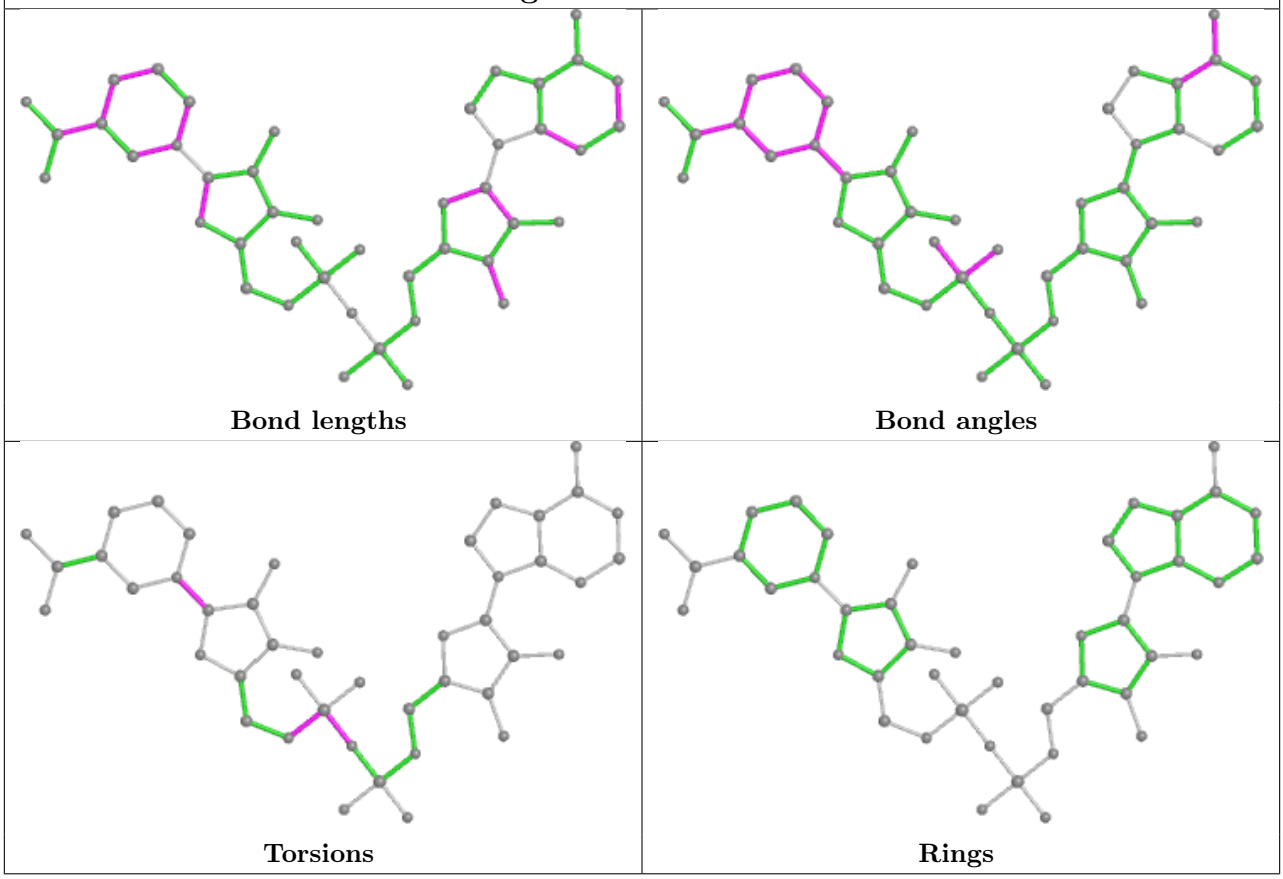
There are no ring outliers.

2 monomers are involved in 2 short contacts:

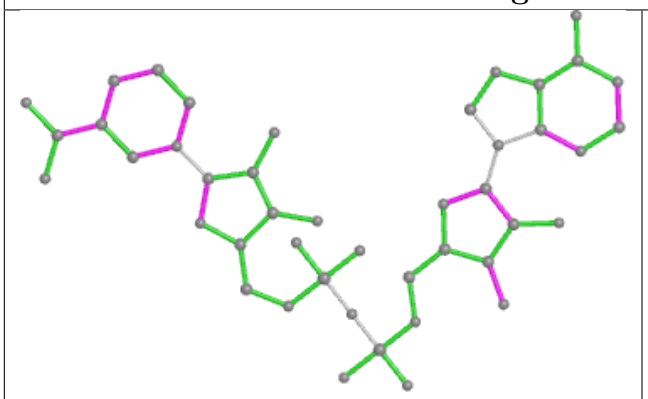
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1403	NAD	1	0
4	C	402	NAD	1	0

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

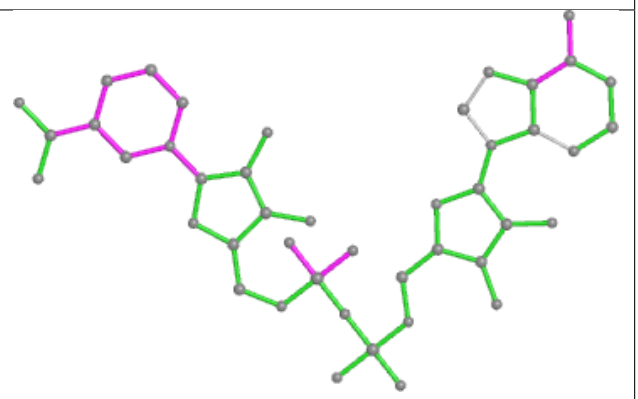
Ligand NAD A 400



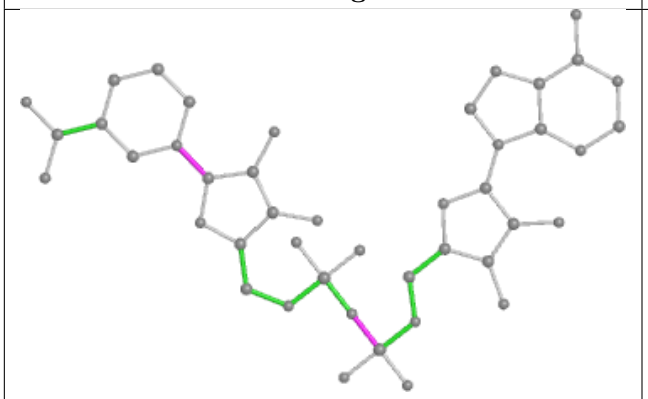
Ligand NAD D 1403



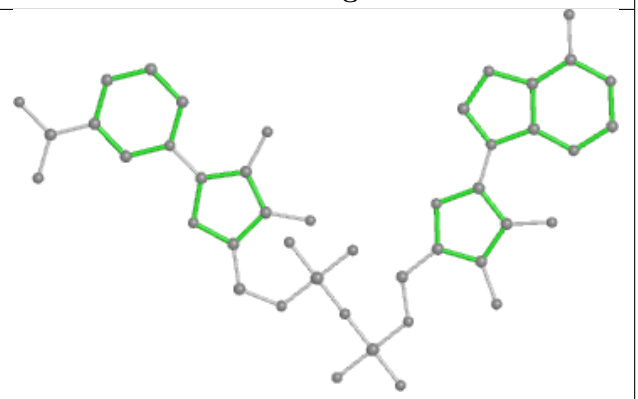
Bond lengths



Bond angles

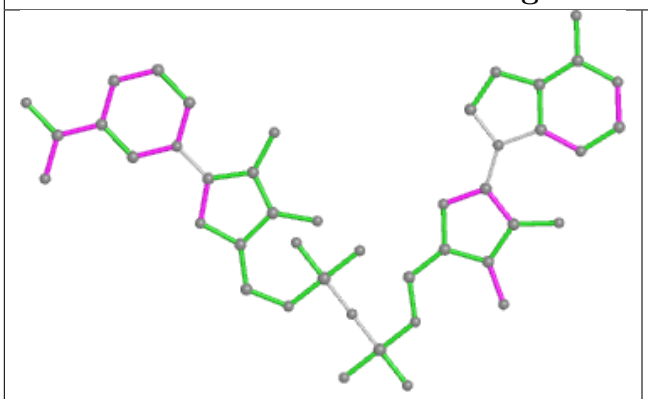


Torsions

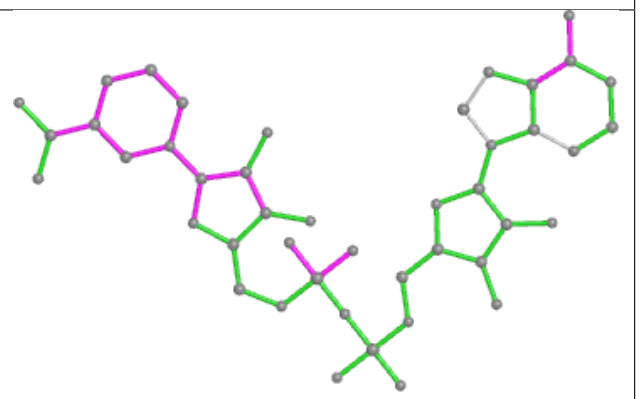


Rings

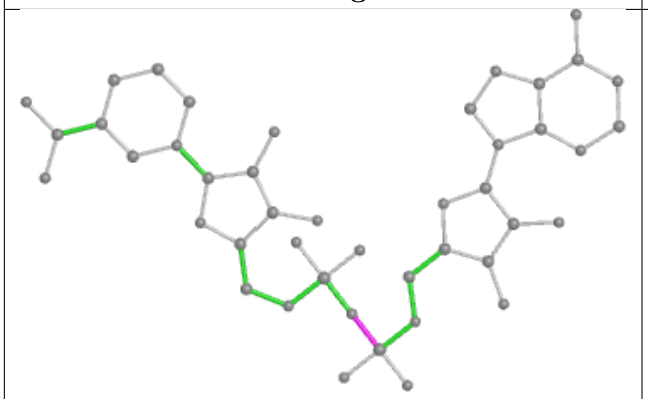
Ligand NAD B 401



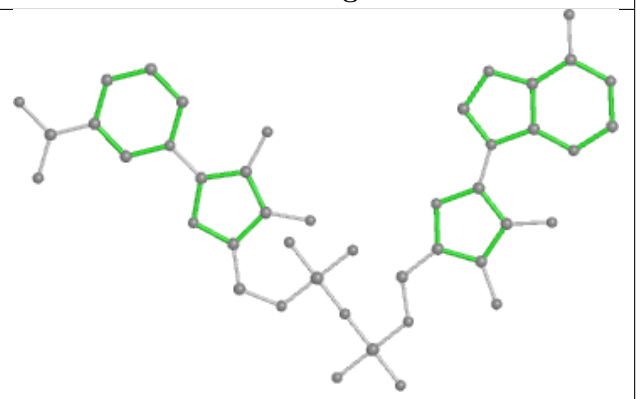
Bond lengths



Bond angles

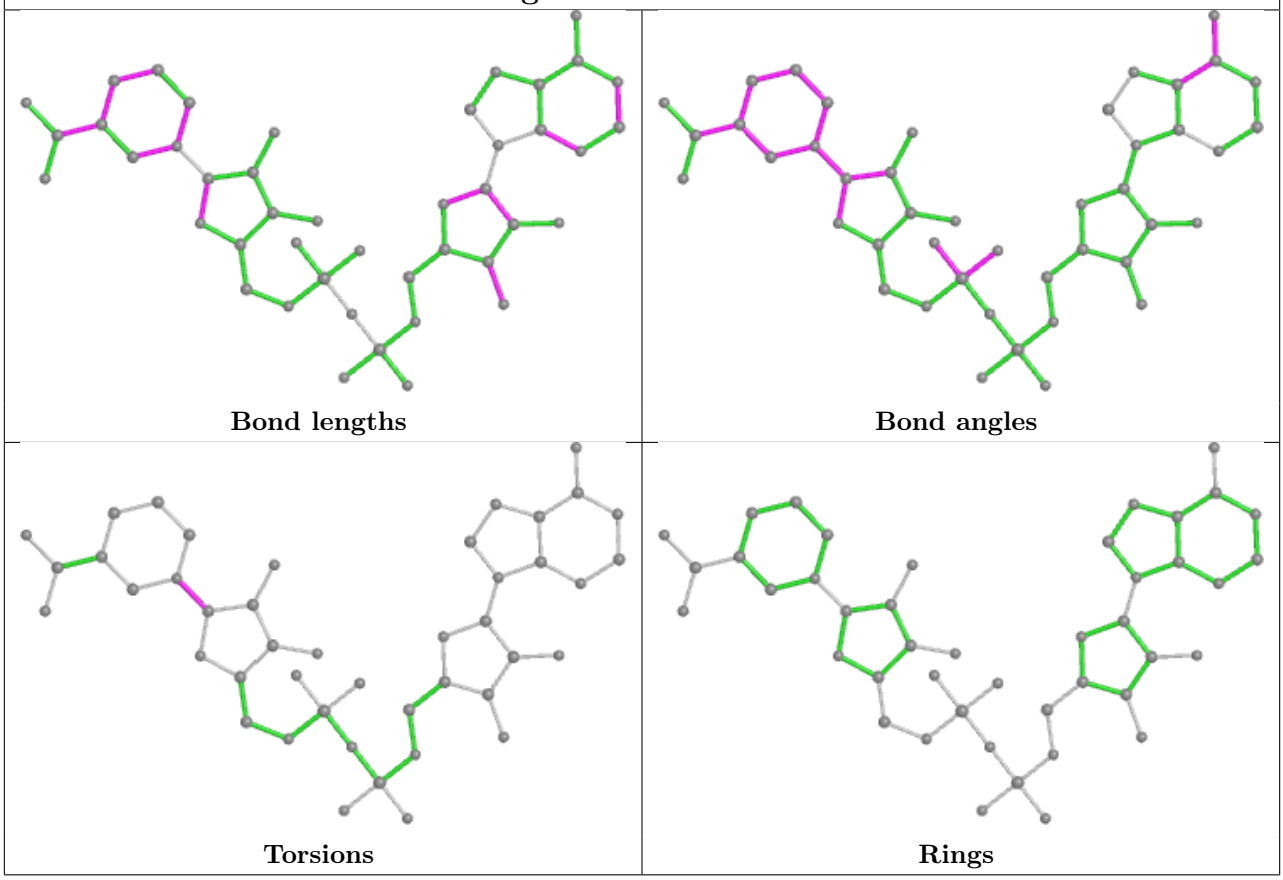


Torsions



Rings

Ligand NAD C 402



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	378/393 (96%)	-0.03	13 (3%)	45	41	32, 51, 89, 118	0
1	B	383/393 (97%)	-0.06	14 (3%)	41	37	31, 49, 89, 121	0
1	C	377/393 (95%)	0.01	14 (3%)	41	37	32, 53, 96, 120	0
1	D	381/393 (96%)	-0.05	15 (3%)	39	34	29, 49, 85, 113	0
All	All	1519/1572 (96%)	-0.03	56 (3%)	41	37	29, 51, 90, 121	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	259	ARG	6.2
1	C	377	THR	5.4
1	C	378	ARG	4.7
1	C	350	PHE	4.1
1	B	261	GLY	3.9
1	D	67	THR	3.8
1	B	354	LEU	3.7
1	C	379	ALA	3.7
1	B	377	THR	3.6
1	A	10	LEU	3.5
1	A	378	ARG	3.4
1	B	3	ASN	3.2
1	A	364	LYS	3.2
1	A	9	ILE	3.2
1	C	61	LYS	3.2
1	A	225	VAL	3.2
1	C	279	ALA	3.1
1	D	260	GLU	3.1
1	A	3	ASN	3.1
1	D	110	ILE	3.0
1	B	137	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	375	TYR	3.0
1	D	258	GLU	3.0
1	A	111	ALA	3.0
1	C	351	ASN	2.9
1	D	261	GLY	2.8
1	D	259	ARG	2.8
1	A	109	VAL	2.7
1	B	121	THR	2.7
1	B	378	ARG	2.7
1	D	311	VAL	2.7
1	A	12	ARG	2.7
1	A	110	ILE	2.6
1	B	350	PHE	2.6
1	D	392	ALA	2.6
1	A	66	ILE	2.6
1	D	375	TYR	2.5
1	B	340	GLY	2.5
1	B	294	GLY	2.5
1	C	381	VAL	2.4
1	B	62	ARG	2.4
1	C	68	PRO	2.3
1	B	260	GLU	2.3
1	D	138	PRO	2.3
1	A	180	LEU	2.3
1	C	309	LYS	2.2
1	C	62	ARG	2.2
1	D	49	SER	2.2
1	D	389	VAL	2.2
1	D	294	GLY	2.1
1	D	350	PHE	2.1
1	D	111	ALA	2.1
1	B	124	VAL	2.1
1	A	142	LEU	2.1
1	C	57	GLU	2.1
1	C	140	THR	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 monosaccharides in this entry.

6.4 Ligands (i)

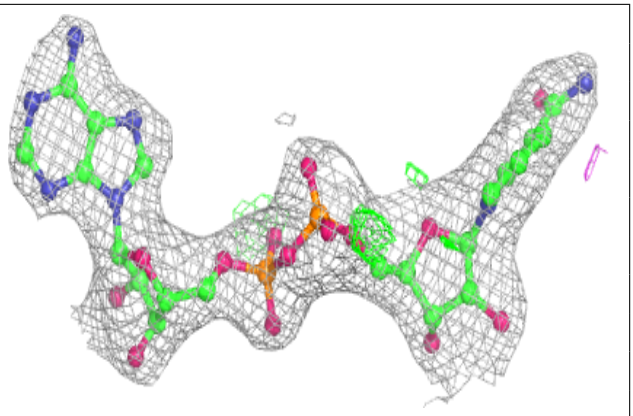
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

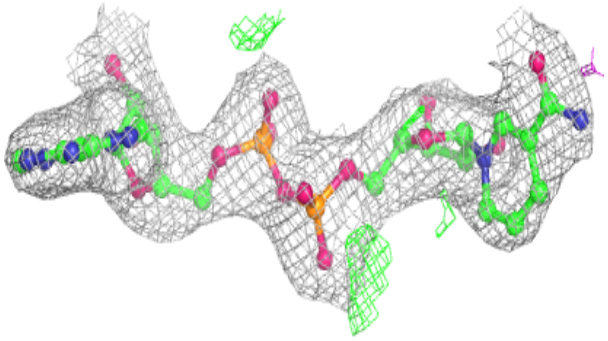
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ZN	A	600	1/1	0.74	0.10	116,116,116,116	0
2	ZN	D	1603	1/1	0.80	0.12	102,102,102,102	0
3	CL	A	604	1/1	0.89	0.21	61,61,61,61	0
3	CL	C	1605	1/1	0.94	0.20	59,59,59,59	0
2	ZN	B	601	1/1	0.95	0.04	100,100,100,100	0
2	ZN	C	602	1/1	0.96	0.09	117,117,117,117	0
4	NAD	A	400	44/44	0.97	0.12	28,44,56,62	0
4	NAD	C	402	44/44	0.97	0.10	31,45,52,54	0
4	NAD	D	1403	44/44	0.97	0.11	25,39,49,56	0
4	NAD	B	401	44/44	0.98	0.10	23,40,47,50	0

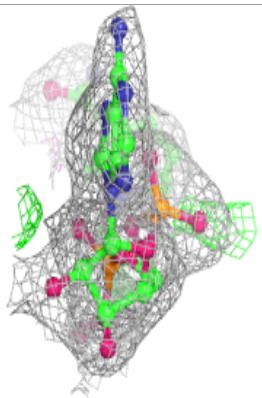
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAD A 400:

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

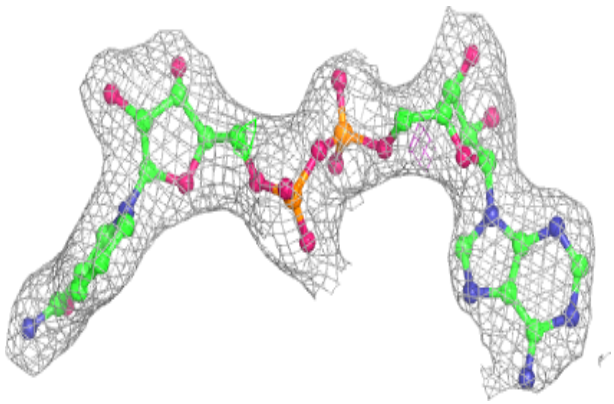


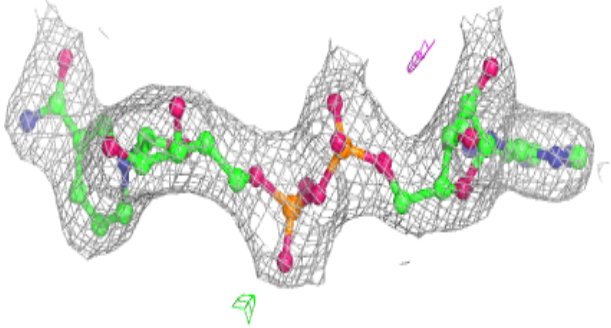


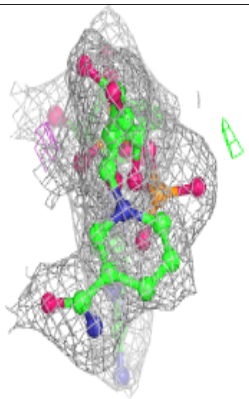


Electron density around NAD C 402:

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

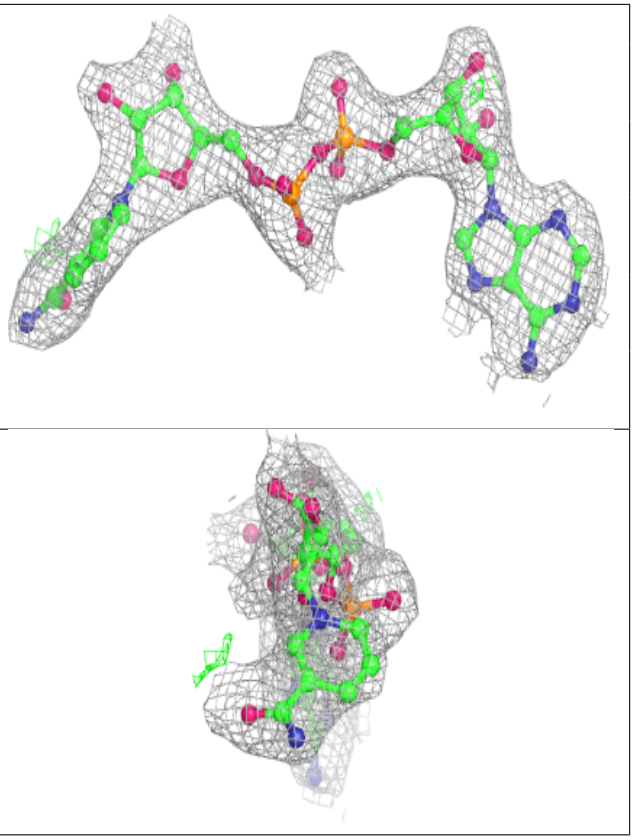






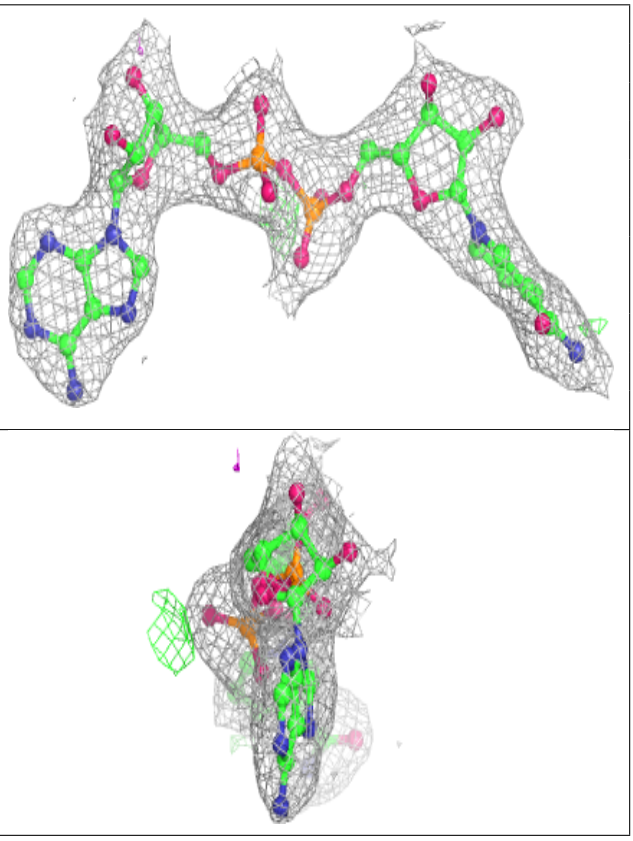
Electron density around NAD D 1403:

$2mF_o$-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 B 401:

$2mF_o$-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.