



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:32 PM UTC

PDB ID : 1QXS / pdb_00001qxs
Title : CRYSTAL STRUCTURE OF Trypanosoma cruzi GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE COMPLEXED WITH AN ANALOGUE OF 1,3- BisPHOSPHO-D-GLYCERIC ACID
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Deposited on : 2003-09-08
Resolution : 2.75 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

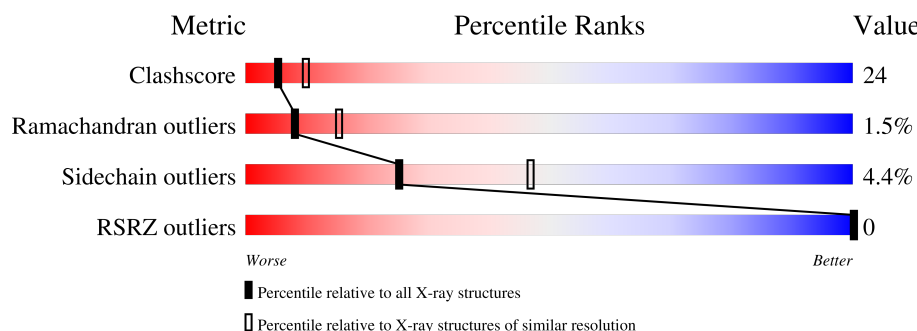
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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))
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

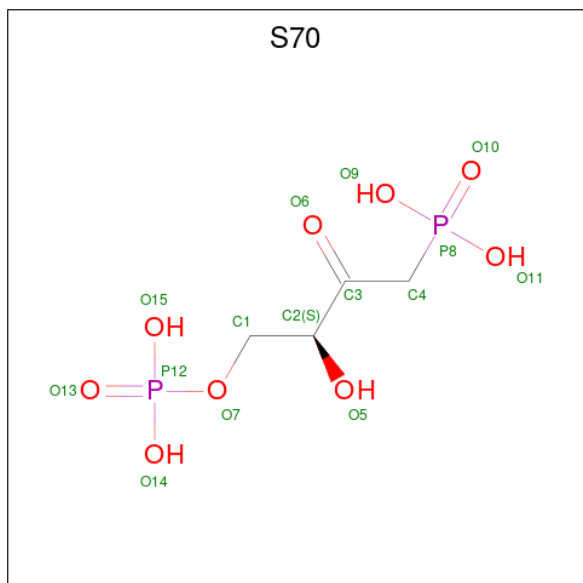
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	359	58% 38% .
1	B	359	56% 39% 5%
1	C	359	52% 45% ..
1	D	359	61% 35% .

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

- Molecule 3 is 3-HYDROXY-2-OXO-4-PHOPHONOXY- BUTYL)-PHOSPHONIC ACID (CCD ID: S70) (formula: C₄H₁₀O₉P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			15	4	9	2		

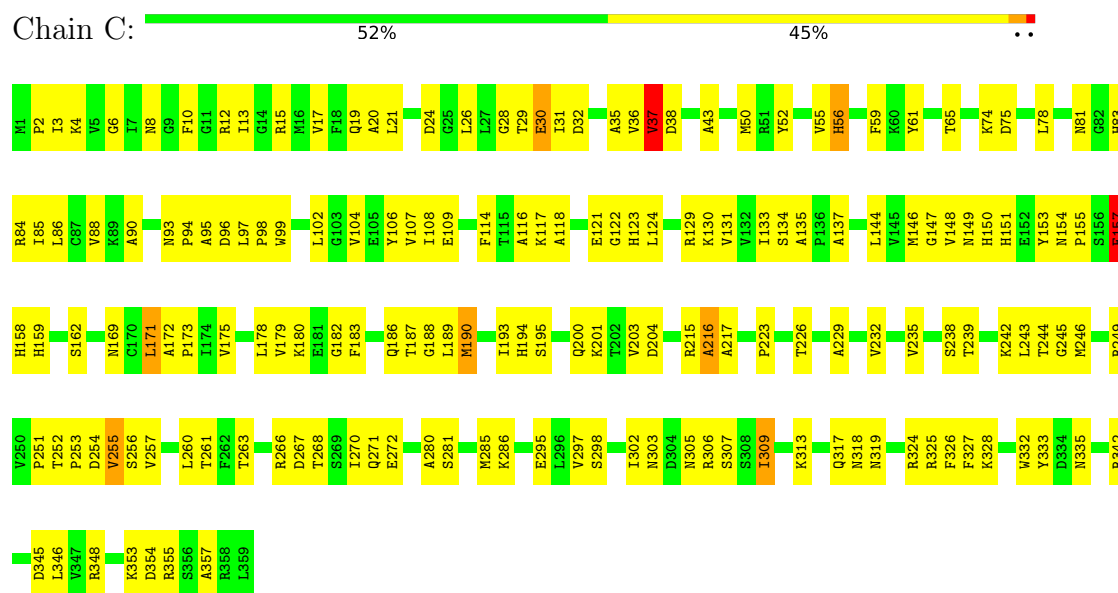
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	76	Total	O	0	0
			76	76		
4	D	127	Total	O	0	0
			127	127		
4	A	122	Total	O	0	0
			122	122		
4	B	128	Total	O	0	0
			128	128		

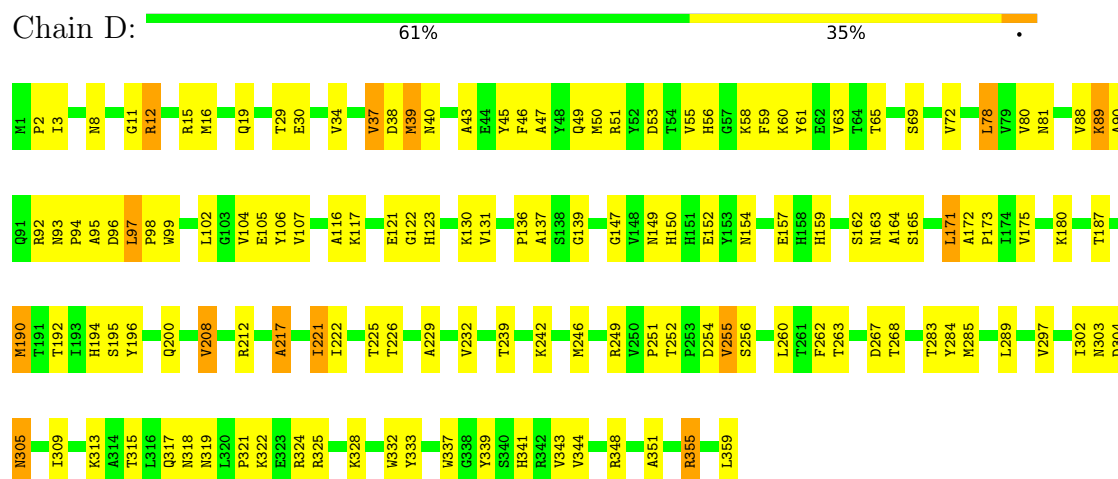
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase, glycosomal

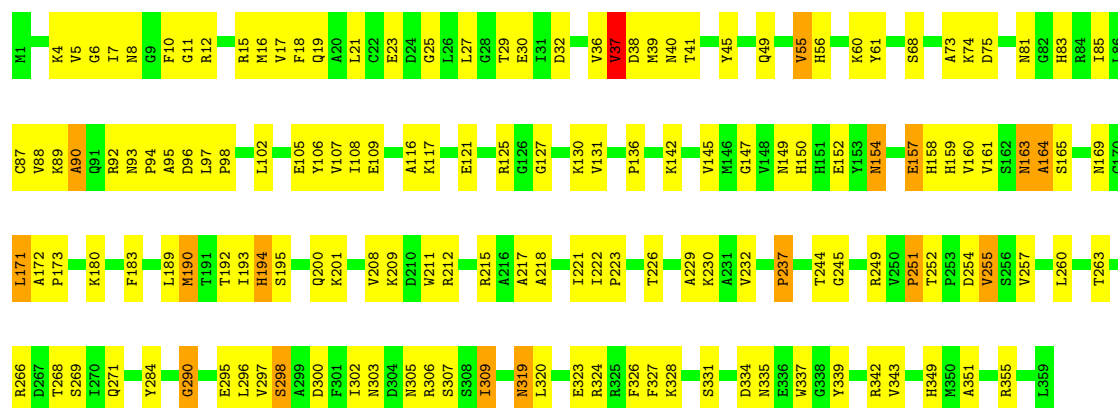


- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase, glycosomal



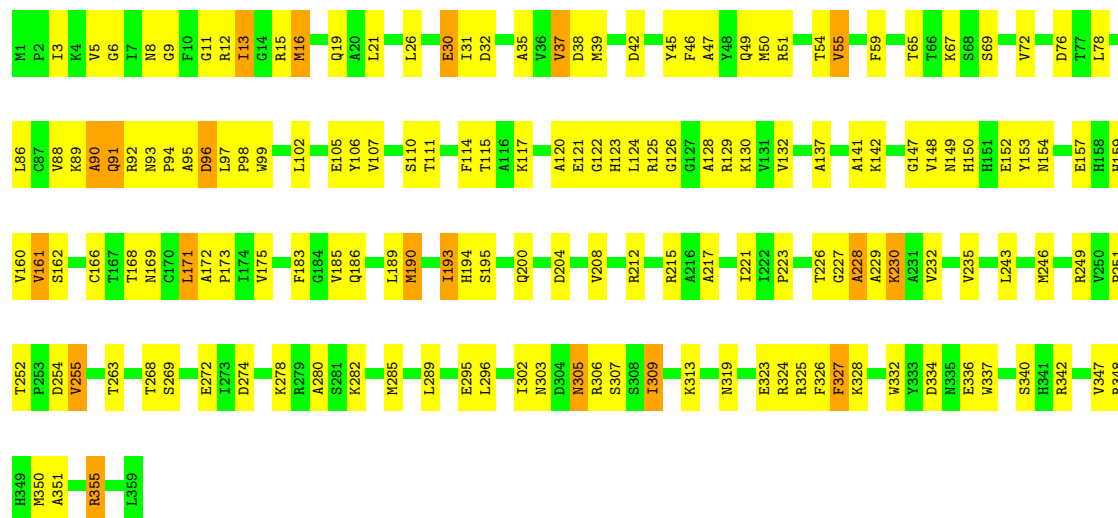
- Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase, glycosomal

Chain A:  58% 38% .



• Molecule 1: Glyceraldehyde 3-phosphate dehydrogenase, glycosomal

Chain B:  56% 39% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.76Å 85.19Å 106.42Å 90.00° 96.74° 90.00°	Depositor
Resolution (Å)	8.00 – 2.75 8.00 – 2.75	Depositor EDS
% Data completeness (in resolution range)	92.4 (8.00-2.75) 82.7 (8.00-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 2.74Å)	Xtriage
Refinement program	REFMAC 5.1.24, CNS 1.0	Depositor
R, R_{free}	0.198 , 0.276 0.189 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.640	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 82.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11576	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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: NAD, S70

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	0.46	0/2786	0.96	12/3778 (0.3%)
1	B	0.44	0/2786	0.94	10/3778 (0.3%)
1	C	0.41	0/2786	0.89	3/3778 (0.1%)
1	D	0.43	0/2786	0.96	9/3778 (0.2%)
All	All	0.43	0/11144	0.94	34/15112 (0.2%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	337	TRP	N-CA-C	8.26	121.09	111.02
1	A	217	ALA	N-CA-C	7.18	118.75	111.07
1	B	217	ALA	N-CA-C	6.98	118.68	111.14
1	A	222	ILE	N-CA-C	6.88	114.36	107.55
1	C	217	ALA	N-CA-C	6.70	118.66	111.36
1	B	208	VAL	N-CA-C	6.46	119.17	111.09
1	A	334	ASP	N-CA-C	-6.38	98.78	108.67
1	D	222	ILE	N-CA-C	6.34	113.83	107.55
1	A	154	ASN	CA-C-N	6.29	126.49	119.32
1	A	154	ASN	C-N-CA	6.29	126.49	119.32
1	B	152	GLU	N-CA-C	-6.25	105.59	112.72
1	D	260	LEU	N-CA-C	6.20	118.81	108.96
1	A	92	ARG	N-CA-C	-6.09	106.17	113.97
1	A	163	ASN	N-CA-C	-5.96	105.76	113.16
1	B	55	VAL	N-CA-C	5.94	116.92	111.81
1	B	337	TRP	N-CA-C	5.73	117.33	111.14
1	B	193	ILE	N-CA-C	-5.70	99.28	107.37
1	C	193	ILE	N-CA-C	-5.63	99.37	107.37
1	A	251	PRO	N-CA-C	5.61	124.03	112.47
1	D	318	ASN	N-CA-C	5.57	119.26	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	327	PHE	N-CA-C	5.55	117.62	109.24
1	D	196	TYR	N-CA-C	5.50	117.42	110.33
1	A	290	GLY	N-CA-C	-5.39	103.09	111.12
1	D	217	ALA	N-CA-C	5.34	118.96	112.23
1	C	56	HIS	N-CA-C	5.30	119.38	113.02
1	A	55	VAL	N-CA-C	5.29	116.36	111.81
1	A	194	HIS	N-CA-C	5.26	117.53	109.07
1	D	221	ILE	N-CA-C	-5.25	98.90	106.88
1	D	304	ASP	N-CA-C	-5.24	101.44	109.76
1	B	161	VAL	N-CA-C	5.23	116.39	108.96
1	D	208	VAL	N-CA-C	5.22	115.99	110.72
1	B	296	LEU	N-CA-C	5.14	118.49	110.32
1	B	334	ASP	N-CA-C	-5.05	99.55	108.23
1	A	56	HIS	N-CA-C	5.04	119.06	113.02

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	2733	0	2722	135	0
1	B	2733	0	2722	150	0
1	C	2733	0	2722	165	0
1	D	2733	0	2722	110	0
2	A	44	0	26	2	0
2	B	44	0	26	6	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
3	A	15	0	6	1	0
4	A	122	0	0	13	0
4	B	128	0	0	9	0
4	C	76	0	0	7	0
4	D	127	0	0	12	0
All	All	11576	0	10998	533	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 24.

All (533) 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:154:ASN:HD22	1:A:157:GLU:HB2	1.21	1.05
1:C:319:ASN:HD21	1:C:328:LYS:H	1.07	0.97
1:C:147:GLY:H	1:C:150:HIS:HD2	1.12	0.96
1:D:195:SER:HB2	1:D:252:THR:O	1.70	0.92
1:A:125:ARG:HG2	1:A:125:ARG:HH11	1.38	0.89
1:A:195:SER:HB2	1:A:252:THR:O	1.72	0.88
1:A:15:ARG:O	1:A:19:GLN:HG3	1.72	0.87
1:D:37:VAL:HB	1:D:88:VAL:HG22	1.58	0.85
1:D:319:ASN:HD21	1:D:328:LYS:H	1.22	0.85
1:B:195:SER:HB3	1:B:252:THR:OG1	1.75	0.85
1:D:8:ASN:HA	1:D:37:VAL:HG13	1.56	0.85
1:A:154:ASN:ND2	1:A:157:GLU:HB2	1.91	0.85
1:D:190:MET:HG2	1:D:229:ALA:HB2	1.59	0.85
1:C:147:GLY:H	1:C:150:HIS:CD2	1.96	0.83
1:D:19:GLN:HG3	1:D:59:PHE:HE1	1.41	0.83
1:D:117:LYS:O	1:D:121:GLU:HG3	1.80	0.82
1:B:117:LYS:O	1:B:121:GLU:HG3	1.79	0.81
1:D:37:VAL:HB	1:D:88:VAL:CG2	2.12	0.80
1:D:63:VAL:HG22	1:D:80:VAL:HG22	1.63	0.80
1:D:171:LEU:HD13	1:D:232:VAL:HG21	1.64	0.80
1:B:183:PHE:HD1	1:B:268:THR:HG21	1.44	0.79
1:C:21:LEU:HD22	1:C:31:ILE:HD12	1.65	0.79
1:C:61:TYR:HB3	1:C:81:ASN:ND2	1.98	0.79
1:A:94:PRO:HA	1:A:97:LEU:HD23	1.65	0.78
1:B:285:MET:HB3	1:B:289:LEU:HB3	1.66	0.78
1:B:3:ILE:HD11	1:B:31:ILE:HG12	1.66	0.78
1:B:12:ARG:O	1:B:16:MET:HG2	1.84	0.77
1:B:8:ASN:HA	1:B:37:VAL:HG13	1.65	0.77
1:C:93:ASN:HB3	1:C:96:ASP:OD2	1.86	0.76
1:A:38:ASP:O	1:A:90:ALA:HB2	1.86	0.76
3:A:804:S70:H42	4:A:973:HOH:O	1.86	0.75
1:D:19:GLN:HG3	1:D:59:PHE:CE1	2.21	0.75
1:C:26:LEU:HD13	1:C:348:ARG:HD3	1.67	0.75
1:B:183:PHE:CD1	1:B:268:THR:HG21	2.21	0.75
1:C:195:SER:HB2	1:C:252:THR:O	1.86	0.75
1:A:194:HIS:HB3	1:A:249:ARG:HD3	1.68	0.75
1:B:8:ASN:HA	1:B:37:VAL:CG1	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:LYS:CD	1:B:159:HIS:HA	2.18	0.74
1:C:38:ASP:O	1:C:90:ALA:HB2	1.86	0.74
1:B:11:GLY:HA3	2:B:863:NAD:H4B	1.69	0.74
1:C:8:ASN:HA	1:C:37:VAL:HG13	1.68	0.73
1:B:38:ASP:C	1:B:90:ALA:HB2	2.13	0.73
1:B:195:SER:HB2	1:B:252:THR:O	1.87	0.73
1:A:18:PHE:HA	1:A:21:LEU:HD12	1.71	0.73
1:B:91:GLN:HG3	1:B:96:ASP:HB2	1.71	0.73
1:B:129:ARG:HG2	1:B:130:LYS:HG2	1.70	0.72
1:D:117:LYS:HB3	1:D:139:GLY:HA3	1.70	0.72
1:D:39:MET:HE3	1:D:39:MET:HA	1.72	0.72
1:A:183:PHE:CD1	1:A:268:THR:HG21	2.25	0.71
1:C:109:GLU:HG2	1:C:123:HIS:NE2	2.06	0.71
1:A:11:GLY:HA3	2:A:862:NAD:H4B	1.72	0.71
1:B:16:MET:HE2	1:B:16:MET:HA	1.73	0.71
1:D:242:LYS:HE2	4:D:957:HOH:O	1.89	0.71
1:D:263:THR:HA	1:D:325:ARG:O	1.91	0.71
1:B:38:ASP:HA	2:B:863:NAD:H2A	1.73	0.70
1:B:45:TYR:O	1:B:49:GLN:HG3	1.90	0.70
1:D:147:GLY:H	1:D:150:HIS:HD2	1.38	0.70
1:A:61:TYR:HD2	1:A:81:ASN:ND2	1.90	0.70
1:A:41:THR:O	1:A:41:THR:HG22	1.90	0.69
1:A:200:GLN:OE1	1:A:249:ARG:NH1	2.25	0.69
1:C:83:HIS:HB2	4:C:881:HOH:O	1.92	0.69
1:C:194:HIS:HB3	1:C:249:ARG:HD3	1.75	0.68
1:C:107:VAL:CG1	1:C:131:VAL:HG22	2.24	0.68
1:A:183:PHE:HD1	1:A:268:THR:HG21	1.59	0.68
1:A:298:SER:HB2	4:A:955:HOH:O	1.94	0.68
1:A:147:GLY:H	1:A:150:HIS:CD2	2.12	0.67
1:D:147:GLY:H	1:D:150:HIS:CD2	2.12	0.67
1:A:125:ARG:HG2	1:A:125:ARG:NH1	2.07	0.67
1:D:172:ALA:HB3	1:D:173:PRO:HD3	1.77	0.67
1:B:98:PRO:O	1:B:102:LEU:HD23	1.96	0.66
1:D:239:THR:HA	1:D:242:LYS:HD2	1.75	0.66
1:B:185:VAL:HG11	1:B:243:LEU:HD21	1.75	0.66
1:A:39:MET:HA	4:A:900:HOH:O	1.95	0.66
1:C:109:GLU:HG2	1:C:123:HIS:HE2	1.61	0.66
1:D:351:ALA:O	1:D:355:ARG:HD2	1.96	0.65
1:B:67:LYS:HD2	1:B:72:VAL:HB	1.77	0.65
1:B:124:LEU:HA	1:B:128:ALA:O	1.96	0.65
1:A:10:PHE:CE2	1:A:15:ARG:HG2	2.31	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:ASN:ND2	1:C:95:ALA:H	1.94	0.65
1:C:35:ALA:HB2	1:C:86:LEU:HB3	1.79	0.65
1:D:93:ASN:HD22	1:D:94:PRO:HD2	1.62	0.64
1:A:172:ALA:HB3	1:A:173:PRO:HD3	1.77	0.64
1:D:359:LEU:HA	4:D:969:HOH:O	1.97	0.64
1:B:93:ASN:HB3	1:B:96:ASP:OD2	1.98	0.64
1:C:130:LYS:CD	1:C:159:HIS:HA	2.28	0.64
1:A:105:GLU:HG3	1:A:106:TYR:CD1	2.33	0.64
1:A:257:VAL:HG23	1:A:331:SER:O	1.98	0.64
1:B:147:GLY:H	1:B:150:HIS:CD2	2.16	0.63
1:A:8:ASN:HA	1:A:37:VAL:HG13	1.81	0.63
1:B:130:LYS:HD2	1:B:159:HIS:HA	1.80	0.63
1:A:145:VAL:HG13	1:A:169:ASN:OD1	1.97	0.63
1:B:226:THR:HG22	1:B:246:MET:HA	1.80	0.63
1:D:47:ALA:HB1	1:D:51:ARG:NH1	2.13	0.63
1:A:212:ARG:HG2	1:B:295:GLU:HB3	1.81	0.63
1:B:47:ALA:HB1	1:B:51:ARG:NH1	2.14	0.62
1:A:142:LYS:HB3	1:A:161:VAL:HG12	1.81	0.62
1:C:171:LEU:HD22	1:C:175:VAL:HG23	1.82	0.62
1:D:254:ASP:O	1:D:255:VAL:HB	2.00	0.62
1:C:24:ASP:HB2	1:C:26:LEU:HG	1.81	0.62
1:C:215:ARG:NH1	1:D:297:VAL:HG22	2.15	0.62
1:B:37:VAL:HG23	1:B:90:ALA:HA	1.80	0.62
1:C:8:ASN:HA	1:C:37:VAL:CG1	2.29	0.62
1:C:319:ASN:HD21	1:C:328:LYS:N	1.88	0.62
1:D:194:HIS:HB3	1:D:249:ARG:HD3	1.80	0.62
1:B:172:ALA:HB3	1:B:173:PRO:HD3	1.82	0.62
1:C:21:LEU:CD2	1:C:31:ILE:HD12	2.30	0.61
1:C:190:MET:O	1:C:245:GLY:HA3	1.99	0.61
1:D:11:GLY:HA2	1:D:15:ARG:NH1	2.15	0.61
1:B:69:SER:O	1:B:72:VAL:HG23	1.99	0.61
1:B:95:ALA:HA	4:B:939:HOH:O	2.00	0.61
1:B:37:VAL:CG2	1:B:90:ALA:HA	2.30	0.61
1:C:251:PRO:HB2	1:D:251:PRO:HB2	1.82	0.61
1:D:285:MET:HB3	1:D:289:LEU:HB3	1.83	0.61
1:D:341:HIS:HB3	4:D:939:HOH:O	2.02	0.60
1:A:25:GLY:HA2	4:A:889:HOH:O	2.00	0.60
1:A:251:PRO:HB2	1:B:251:PRO:HB2	1.84	0.60
1:A:37:VAL:HB	1:A:88:VAL:HG22	1.82	0.60
1:B:50:MET:O	1:B:59:PHE:HB2	2.01	0.60
1:C:37:VAL:HB	1:C:88:VAL:CG2	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:93:ASN:HB3	1:D:96:ASP:OD2	2.01	0.60
1:B:263:THR:HA	1:B:325:ARG:O	2.02	0.60
1:D:94:PRO:HA	1:D:97:LEU:HD23	1.84	0.60
1:D:107:VAL:HG22	1:D:131:VAL:HG22	1.82	0.59
1:B:115:THR:HG23	4:B:913:HOH:O	2.02	0.59
1:A:93:ASN:ND2	1:A:95:ALA:H	2.01	0.59
1:B:3:ILE:CD1	1:B:31:ILE:HG12	2.32	0.59
1:C:180:LYS:C	1:C:182:GLY:H	2.10	0.59
1:C:239:THR:HA	1:C:242:LYS:HD3	1.85	0.59
1:A:98:PRO:O	1:A:102:LEU:HD23	2.02	0.59
1:A:351:ALA:O	1:A:355:ARG:HG2	2.03	0.59
1:A:4:LYS:HA	1:A:4:LYS:HE2	1.84	0.59
1:B:13:ILE:HD11	2:B:863:NAD:H71N	1.67	0.59
1:C:254:ASP:O	1:C:255:VAL:HB	2.02	0.59
1:A:29:THR:HG22	1:A:30:GLU:HG2	1.85	0.59
1:A:260:LEU:O	1:A:328:LYS:HA	2.02	0.59
1:D:136:PRO:CG	1:D:165:SER:HB3	2.33	0.58
1:B:30:GLU:HG3	1:B:348:ARG:NH1	2.18	0.58
1:B:302:ILE:O	1:B:303:ASN:HB2	2.03	0.58
1:B:324:ARG:HD3	4:B:927:HOH:O	2.03	0.58
1:D:8:ASN:HD22	1:D:37:VAL:HG13	1.67	0.58
1:D:171:LEU:HD22	1:D:175:VAL:HG23	1.84	0.58
1:B:204:ASP:OD1	1:B:215:ARG:HA	2.03	0.58
1:A:107:VAL:HG22	1:A:131:VAL:HG22	1.85	0.58
1:A:105:GLU:HG3	1:A:106:TYR:HD1	1.68	0.58
1:C:267:ASP:OD1	1:C:325:ARG:NE	2.36	0.58
1:C:200:GLN:HB3	4:C:864:HOH:O	2.02	0.58
1:D:319:ASN:HD21	1:D:328:LYS:N	1.98	0.58
1:C:190:MET:HG2	1:C:229:ALA:HB2	1.85	0.58
1:C:268:THR:HG22	1:C:272:GLU:OE2	2.04	0.58
1:B:37:VAL:HB	1:B:88:VAL:CG2	2.33	0.58
1:D:226:THR:HG22	1:D:246:MET:HA	1.86	0.57
1:D:12:ARG:HH22	1:D:53:ASP:CG	2.12	0.57
1:C:130:LYS:HD3	1:C:159:HIS:HA	1.85	0.57
1:B:319:ASN:HD21	1:B:328:LYS:H	1.51	0.57
1:C:270:ILE:HG22	1:C:271:GLN:NE2	2.19	0.57
1:B:268:THR:HG22	1:B:272:GLU:OE1	2.03	0.57
1:A:226:THR:HG22	1:A:245:GLY:O	2.05	0.57
1:A:254:ASP:O	1:A:255:VAL:HB	2.04	0.57
1:B:8:ASN:HD22	1:B:37:VAL:HG13	1.69	0.57
1:D:98:PRO:O	1:D:102:LEU:HD23	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:ILE:O	1:C:17:VAL:HG23	2.04	0.57
1:A:223:PRO:HG2	1:B:332:TRP:HZ2	1.70	0.57
1:B:229:ALA:HB3	4:B:868:HOH:O	2.04	0.57
1:C:172:ALA:HB3	1:C:173:PRO:HD3	1.85	0.56
1:C:201:LYS:HE2	4:C:904:HOH:O	2.04	0.56
1:B:21:LEU:O	1:B:21:LEU:HD23	2.05	0.56
1:B:31:ILE:HG22	1:B:32:ASP:N	2.20	0.56
1:D:256:SER:HB2	1:D:333:TYR:CZ	2.41	0.56
1:D:305:ASN:C	1:D:305:ASN:HD22	2.13	0.56
1:C:95:ALA:HA	1:C:122:GLY:O	2.06	0.56
1:C:124:LEU:HD21	1:C:131:VAL:HG23	1.88	0.55
1:C:144:LEU:HD11	1:C:153:TYR:HB2	1.87	0.55
1:A:319:ASN:HD21	1:A:328:LYS:H	1.53	0.55
1:C:35:ALA:HA	1:C:85:ILE:HG23	1.88	0.55
1:D:3:ILE:O	1:D:3:ILE:HG13	2.06	0.55
1:D:123:HIS:HB2	1:D:131:VAL:HG21	1.86	0.55
1:A:147:GLY:H	1:A:150:HIS:HD2	1.54	0.55
1:B:21:LEU:HD23	1:B:21:LEU:C	2.32	0.55
1:C:302:ILE:HD13	4:C:867:HOH:O	2.06	0.55
1:A:68:SER:HA	4:A:976:HOH:O	2.05	0.55
1:C:6:GLY:HA3	1:C:104:VAL:HG11	1.88	0.55
1:D:45:TYR:O	1:D:49:GLN:HG3	2.06	0.55
1:C:302:ILE:O	1:C:303:ASN:HB2	2.06	0.55
1:A:295:GLU:HB3	1:B:212:ARG:HG3	1.89	0.55
1:C:190:MET:CG	1:C:229:ALA:HB2	2.37	0.55
1:B:154:ASN:HB3	1:B:157:GLU:HB3	1.88	0.55
1:C:215:ARG:HH12	1:D:297:VAL:HG22	1.72	0.54
1:D:46:PHE:O	1:D:50:MET:HG3	2.07	0.54
1:C:32:ASP:OD1	1:C:83:HIS:NE2	2.40	0.54
1:C:93:ASN:OD1	1:C:95:ALA:HB3	2.08	0.54
1:D:94:PRO:HA	1:D:97:LEU:CD2	2.37	0.54
1:A:81:ASN:ND2	4:A:876:HOH:O	2.40	0.54
1:D:339:TYR:O	1:D:343:VAL:HG23	2.07	0.54
1:B:5:VAL:HG22	1:B:106:TYR:HB2	1.89	0.54
1:C:306:ARG:HB2	1:C:309:ILE:CD1	2.37	0.54
1:D:61:TYR:OH	4:D:902:HOH:O	2.18	0.54
1:A:37:VAL:O	1:A:37:VAL:HG22	2.08	0.54
1:C:10:PHE:CE2	1:C:15:ARG:HG2	2.43	0.54
1:D:95:ALA:HA	1:D:122:GLY:O	2.07	0.54
1:C:98:PRO:O	1:C:102:LEU:HD23	2.08	0.53
1:C:318:ASN:OD1	1:D:225:THR:HG22	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:MET:HG2	1:A:229:ALA:HB2	1.90	0.53
1:A:271:GLN:CD	1:A:271:GLN:H	2.16	0.53
1:C:17:VAL:O	1:C:20:ALA:HB3	2.08	0.53
1:C:109:GLU:OE1	1:C:114:PHE:HD2	1.91	0.53
1:B:38:ASP:O	1:B:90:ALA:HB2	2.08	0.53
1:B:67:LYS:HA	1:B:76:ASP:OD1	2.09	0.53
1:C:183:PHE:CD1	1:C:268:THR:HG21	2.44	0.53
1:B:95:ALA:C	1:B:97:LEU:H	2.16	0.53
1:D:65:THR:HG22	1:D:78:LEU:HD12	1.89	0.53
1:A:93:ASN:HB3	1:A:96:ASP:OD2	2.09	0.53
1:B:230:LYS:HB2	1:B:230:LYS:NZ	2.24	0.53
1:C:50:MET:O	1:C:59:PHE:HB2	2.09	0.53
1:A:324:ARG:HG3	1:A:324:ARG:HH11	1.73	0.53
1:C:133:ILE:C	1:C:135:ALA:H	2.16	0.52
1:D:38:ASP:O	1:D:90:ALA:HB2	2.09	0.52
4:D:947:HOH:O	1:A:201:LYS:HG2	2.08	0.52
1:A:125:ARG:NH1	1:A:125:ARG:CG	2.71	0.52
1:C:3:ILE:HG13	1:C:30:GLU:O	2.09	0.52
1:C:263:THR:HA	1:C:325:ARG:O	2.09	0.52
1:A:192:THR:HG23	1:A:192:THR:O	2.08	0.52
1:A:215:ARG:HD2	4:A:896:HOH:O	2.10	0.52
1:B:6:GLY:O	1:B:107:VAL:HA	2.09	0.52
1:D:8:ASN:HA	1:D:37:VAL:CG1	2.35	0.52
1:C:6:GLY:O	1:C:107:VAL:HG23	2.10	0.52
1:D:69:SER:O	1:D:72:VAL:HG23	2.09	0.52
1:C:313:LYS:HB3	1:C:317:GLN:NE2	2.24	0.52
1:A:130:LYS:HD3	1:A:159:HIS:HA	1.92	0.52
1:D:302:ILE:O	1:D:303:ASN:HB2	2.10	0.52
1:C:61:TYR:HB3	1:C:81:ASN:HD22	1.75	0.51
1:D:136:PRO:HG3	1:D:165:SER:HB3	1.92	0.51
1:B:190:MET:HG2	1:B:229:ALA:HB2	1.93	0.51
1:B:195:SER:CB	1:B:252:THR:OG1	2.54	0.51
1:C:36:VAL:HG13	1:C:85:ILE:HG21	1.92	0.51
1:C:146:MET:HE2	1:C:345:ASP:HB3	1.92	0.51
1:C:266:ARG:HD3	4:C:900:HOH:O	2.11	0.51
1:B:194:HIS:HB3	1:B:249:ARG:HD3	1.92	0.51
1:C:93:ASN:HD22	1:C:94:PRO:HD2	1.76	0.51
1:C:270:ILE:HG22	1:C:271:GLN:HE22	1.76	0.51
1:A:221:ILE:HD13	1:B:193:ILE:HG22	1.91	0.51
1:C:123:HIS:HB2	1:C:131:VAL:HG21	1.92	0.51
1:D:163:ASN:O	1:D:164:ALA:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ASN:C	1:A:93:ASN:HD22	2.17	0.51
1:D:208:VAL:HG11	1:A:40:ASN:HD22	1.76	0.51
1:A:295:GLU:HB3	1:B:212:ARG:CG	2.41	0.51
1:B:37:VAL:HB	1:B:88:VAL:HG23	1.93	0.51
1:D:46:PHE:HB3	1:D:78:LEU:HD21	1.94	0.50
1:A:255:VAL:HB	4:A:955:HOH:O	2.12	0.50
1:D:256:SER:HB2	1:D:333:TYR:CE1	2.46	0.50
1:B:130:LYS:HD3	1:B:159:HIS:HA	1.94	0.50
1:B:142:LYS:HB3	1:B:161:VAL:HG12	1.94	0.50
1:B:274:ASP:O	1:B:278:LYS:HG3	2.11	0.50
1:D:93:ASN:HD22	1:D:94:PRO:CD	2.24	0.50
1:B:9:GLY:HA3	1:B:111:THR:HG22	1.93	0.50
1:C:144:LEU:CD1	1:C:153:TYR:HB2	2.41	0.50
1:A:121:GLU:HG2	1:A:160:VAL:HG21	1.93	0.50
1:B:8:ASN:O	1:B:111:THR:HG23	2.11	0.50
1:B:117:LYS:HB2	1:B:141:ALA:HB2	1.93	0.50
1:D:194:HIS:O	1:D:249:ARG:HA	2.12	0.50
1:B:278:LYS:O	1:B:282:LYS:HG3	2.11	0.50
1:C:295:GLU:HB3	1:D:212:ARG:HG2	1.93	0.50
1:B:227:GLY:O	1:B:228:ALA:C	2.54	0.50
1:C:298:SER:CB	1:D:221:ILE:H	2.24	0.49
1:C:354:ASP:O	1:C:357:ALA:HB3	2.11	0.49
4:D:863:HOH:O	1:A:218:ALA:HB3	2.12	0.49
1:A:193:ILE:HG22	1:B:221:ILE:HD13	1.94	0.49
1:B:149:ASN:OD1	1:B:235:VAL:HG13	2.12	0.49
1:C:117:LYS:O	1:C:121:GLU:HG3	2.11	0.49
1:D:180:LYS:HD2	4:D:973:HOH:O	2.12	0.49
1:A:4:LYS:HZ1	1:A:32:ASP:CG	2.20	0.49
1:A:94:PRO:HA	1:A:97:LEU:CD2	2.40	0.49
1:A:266:ARG:NH1	1:A:268:THR:HG22	2.28	0.49
1:A:349:HIS:HA	4:A:890:HOH:O	2.11	0.49
1:A:8:ASN:HD22	1:A:37:VAL:HG13	1.76	0.49
1:A:93:ASN:ND2	1:A:93:ASN:C	2.69	0.49
1:C:298:SER:HB2	1:D:221:ILE:H	1.77	0.49
1:D:149:ASN:O	1:D:152:GLU:HG3	2.12	0.49
1:A:4:LYS:NZ	1:A:32:ASP:CG	2.70	0.49
1:C:30:GLU:OE1	1:C:348:ARG:HD2	2.13	0.49
1:C:99:TRP:NE1	1:C:123:HIS:ND1	2.59	0.49
1:B:3:ILE:HG12	1:B:30:GLU:O	2.13	0.49
1:A:319:ASN:HD21	1:A:327:PHE:HA	1.78	0.49
1:B:106:TYR:CD2	1:B:130:LYS:HB2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:LEU:HD22	1:C:129:ARG:O	2.13	0.48
1:D:105:GLU:HG3	1:D:106:TYR:CD1	2.47	0.48
1:D:154:ASN:HB3	1:D:157:GLU:HB3	1.93	0.48
1:A:307:SER:OG	1:A:342:ARG:HD2	2.13	0.48
1:C:226:THR:HG22	1:C:246:MET:HA	1.96	0.48
1:D:116:ALA:O	1:D:117:LYS:C	2.54	0.48
1:A:249:ARG:HG3	1:A:249:ARG:HH11	1.78	0.48
1:C:254:ASP:O	1:C:255:VAL:CB	2.60	0.48
1:A:95:ALA:CB	1:A:125:ARG:HD2	2.43	0.48
1:A:326:PHE:CE1	1:B:326:PHE:CE1	3.01	0.48
1:C:253:PRO:HD2	4:D:929:HOH:O	2.13	0.48
1:B:124:LEU:HD11	1:B:160:VAL:HG23	1.94	0.48
1:C:171:LEU:HD13	1:C:232:VAL:HG21	1.93	0.48
1:D:47:ALA:HB1	1:D:51:ARG:HH11	1.77	0.48
1:A:85:ILE:HD12	1:A:85:ILE:N	2.29	0.48
1:B:35:ALA:HB2	1:B:86:LEU:HB3	1.96	0.48
1:B:194:HIS:O	1:B:249:ARG:HA	2.14	0.48
1:C:52:TYR:HB3	1:A:300:ASP:OD1	2.14	0.48
1:A:263:THR:HG23	1:B:326:PHE:CE1	2.49	0.48
1:C:12:ARG:NH2	1:C:55:VAL:HG12	2.29	0.48
1:C:154:ASN:HB3	1:C:157:GLU:HB3	1.95	0.48
1:D:200:GLN:HG2	4:D:862:HOH:O	2.14	0.48
1:D:55:VAL:HG13	1:D:56:HIS:CD2	2.49	0.48
1:C:61:TYR:HD2	1:C:81:ASN:HD21	1.61	0.47
1:C:261:THR:HA	1:C:327:PHE:O	2.13	0.47
1:D:254:ASP:O	1:D:255:VAL:CB	2.62	0.47
1:B:107:VAL:HG21	1:B:123:HIS:CG	2.48	0.47
1:B:147:GLY:H	1:B:150:HIS:HD2	1.60	0.47
1:C:270:ILE:CG2	1:C:271:GLN:NE2	2.77	0.47
1:A:81:ASN:CB	4:A:876:HOH:O	2.63	0.47
1:C:319:ASN:ND2	1:C:328:LYS:H	1.92	0.47
1:D:43:ALA:HB1	1:D:78:LEU:HD13	1.97	0.47
1:B:91:GLN:CD	1:B:92:ARG:H	2.23	0.47
1:B:305:ASN:H	1:B:305:ASN:ND2	2.11	0.47
1:C:85:ILE:N	1:C:85:ILE:HD12	2.29	0.47
1:C:106:TYR:CD2	1:C:130:LYS:HB2	2.48	0.47
1:B:274:ASP:OD2	1:B:278:LYS:HE3	2.13	0.47
1:C:55:VAL:HG21	1:C:254:ASP:HB3	1.97	0.47
1:B:200:GLN:HG2	4:B:865:HOH:O	2.15	0.47
1:D:8:ASN:HD22	1:D:37:VAL:CG1	2.28	0.47
1:B:186:GLN:NE2	4:B:925:HOH:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:ASP:O	1:B:255:VAL:HB	2.14	0.47
1:C:15:ARG:O	1:C:19:GLN:HG3	2.14	0.47
1:C:188:GLY:O	1:C:243:LEU:HA	2.14	0.47
1:C:98:PRO:HB2	1:C:102:LEU:HD23	1.97	0.47
1:B:15:ARG:O	1:B:19:GLN:HG3	2.15	0.47
1:C:180:LYS:C	1:C:182:GLY:N	2.73	0.47
1:A:8:ASN:HA	1:A:37:VAL:CG1	2.45	0.46
1:A:171:LEU:HD13	1:A:232:VAL:HG21	1.97	0.46
1:B:26:LEU:HD13	1:B:348:ARG:HD3	1.96	0.46
1:B:171:LEU:HD13	1:B:232:VAL:HG21	1.98	0.46
1:D:107:VAL:O	1:D:107:VAL:HG23	2.14	0.46
1:A:41:THR:O	1:A:41:THR:CG2	2.62	0.46
1:A:208:VAL:HB	4:A:929:HOH:O	2.15	0.46
1:B:65:THR:HG22	1:B:78:LEU:HD12	1.97	0.46
1:C:216:ALA:HB3	1:B:54:THR:HG21	1.96	0.46
1:C:280:ALA:O	1:C:285:MET:HB2	2.15	0.46
1:A:130:LYS:CD	1:A:159:HIS:HA	2.46	0.46
1:B:99:TRP:HB2	1:B:126:GLY:O	2.15	0.46
1:B:148:VAL:CG1	1:B:235:VAL:HG12	2.46	0.46
1:C:93:ASN:ND2	1:C:93:ASN:C	2.73	0.46
1:A:319:ASN:HD21	1:A:328:LYS:N	2.13	0.46
1:B:226:THR:HG22	1:B:246:MET:CA	2.44	0.46
1:A:339:TYR:O	1:A:343:VAL:HG23	2.16	0.46
1:C:35:ALA:CB	1:C:86:LEU:HB3	2.45	0.46
1:D:283:THR:OG1	1:D:284:TYR:N	2.48	0.46
1:B:212:ARG:NH1	1:B:223:PRO:HG2	2.31	0.46
1:C:353:LYS:O	1:C:357:ALA:HB2	2.16	0.46
1:B:3:ILE:HG13	1:B:31:ILE:HA	1.98	0.46
1:C:307:SER:OG	1:C:342:ARG:HD2	2.15	0.46
1:B:123:HIS:O	1:B:128:ALA:HB3	2.16	0.46
1:A:81:ASN:HB2	4:A:876:HOH:O	2.16	0.46
1:B:94:PRO:HD3	1:B:114:PHE:CE1	2.51	0.46
1:B:193:ILE:HD12	1:B:193:ILE:N	2.31	0.46
1:C:94:PRO:O	1:C:97:LEU:HD23	2.15	0.45
1:C:194:HIS:O	1:C:249:ARG:HA	2.15	0.45
1:C:215:ARG:O	1:C:216:ALA:C	2.58	0.45
1:D:194:HIS:HB3	1:D:249:ARG:CD	2.44	0.45
1:B:319:ASN:HD22	1:B:327:PHE:HD1	1.64	0.45
1:C:37:VAL:HB	1:C:88:VAL:HG23	1.98	0.45
1:C:268:THR:O	1:C:324:ARG:HA	2.16	0.45
1:A:7:ILE:O	1:A:37:VAL:HG12	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:GLY:HA2	1:B:125:ARG:HH21	1.81	0.45
1:B:168:THR:OG1	1:B:228:ALA:HA	2.16	0.45
1:C:153:TYR:CE1	1:C:158:HIS:HB2	2.51	0.45
1:B:342:ARG:HA	1:B:342:ARG:NE	2.30	0.45
1:C:114:PHE:C	1:C:116:ALA:H	2.23	0.45
1:D:11:GLY:HA2	1:D:15:ARG:HH12	1.79	0.45
1:D:34:VAL:CG1	1:D:102:LEU:HD12	2.47	0.45
1:B:166:CYS:HB3	2:B:863:NAD:H4N	1.99	0.45
1:D:92:ARG:HG3	1:D:92:ARG:HH11	1.82	0.45
1:C:109:GLU:CG	1:C:123:HIS:HE2	2.29	0.45
1:C:203:VAL:O	1:C:204:ASP:C	2.60	0.45
1:D:12:ARG:HA	1:D:12:ARG:HD2	1.73	0.45
1:D:268:THR:O	1:D:324:ARG:HA	2.16	0.45
1:A:163:ASN:O	1:A:164:ALA:HB3	2.16	0.45
1:B:99:TRP:NE1	1:B:123:HIS:ND1	2.61	0.45
1:B:313:LYS:HE3	4:B:886:HOH:O	2.17	0.45
1:C:93:ASN:HD22	1:C:94:PRO:CD	2.30	0.45
1:C:137:ALA:HB2	1:C:162:SER:HB2	1.98	0.45
1:D:190:MET:CG	1:D:229:ALA:HB2	2.39	0.45
1:A:194:HIS:O	1:A:249:ARG:HA	2.17	0.45
1:C:133:ILE:O	1:C:135:ALA:N	2.41	0.45
1:B:212:ARG:HH11	1:B:223:PRO:HG2	1.81	0.45
1:B:336:GLU:HG2	2:B:863:NAD:H72N	1.82	0.45
1:A:306:ARG:HB2	1:A:309:ILE:CD1	2.46	0.45
1:A:324:ARG:HG3	1:A:324:ARG:NH1	2.32	0.45
1:A:158:HIS:HB3	1:A:161:VAL:CG1	2.47	0.44
1:A:319:ASN:ND2	1:A:327:PHE:HA	2.31	0.44
1:C:183:PHE:CE1	1:C:268:THR:HG21	2.52	0.44
1:D:50:MET:O	1:D:59:PHE:HB2	2.16	0.44
1:D:315:THR:HG23	1:D:328:LYS:O	2.17	0.44
1:A:136:PRO:HG3	4:A:885:HOH:O	2.18	0.44
1:B:47:ALA:O	1:B:51:ARG:HB2	2.17	0.44
1:C:148:VAL:HG12	1:C:235:VAL:HG12	1.99	0.44
1:C:324:ARG:C	1:C:325:ARG:HG2	2.42	0.44
1:A:12:ARG:HB2	2:A:862:NAD:O2A	2.18	0.44
1:B:171:LEU:HD22	1:B:175:VAL:HG23	2.00	0.44
1:C:201:LYS:HD2	1:C:201:LYS:N	2.32	0.44
1:A:106:TYR:OH	1:A:351:ALA:HA	2.18	0.44
1:C:65:THR:HG22	1:C:78:LEU:HD12	1.99	0.44
1:B:5:VAL:HA	1:B:106:TYR:O	2.18	0.44
1:C:257:VAL:HB	1:C:332:TRP:CE3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:37:VAL:HB	1:C:88:VAL:HG22	1.98	0.44
1:C:43:ALA:HB1	1:C:65:THR:CG2	2.47	0.44
1:C:118:ALA:O	1:C:121:GLU:HB2	2.18	0.44
1:A:117:LYS:O	1:A:121:GLU:HG3	2.18	0.44
1:C:31:ILE:HG22	1:C:32:ASP:N	2.33	0.43
1:A:23:GLU:OE1	1:A:337:TRP:HZ2	2.01	0.43
1:A:45:TYR:O	1:A:49:GLN:HG3	2.18	0.43
1:A:302:ILE:O	1:A:303:ASN:HB2	2.17	0.43
1:D:180:LYS:NZ	4:D:979:HOH:O	2.49	0.43
1:A:41:THR:HG23	1:A:87:CYS:O	2.17	0.43
1:B:16:MET:HB3	1:B:340:SER:OG	2.19	0.43
1:C:169:ASN:O	1:C:307:SER:HB3	2.18	0.43
1:B:89:LYS:HD3	1:B:90:ALA:N	2.33	0.43
1:B:274:ASP:CG	1:B:278:LYS:HE3	2.43	0.43
1:C:107:VAL:HG13	1:C:131:VAL:HG22	2.00	0.43
1:D:2:PRO:HG3	1:D:29:THR:O	2.19	0.43
1:A:136:PRO:CG	1:A:165:SER:HB3	2.48	0.43
1:B:51:ARG:HG3	4:B:882:HOH:O	2.19	0.43
1:C:215:ARG:HG3	1:C:215:ARG:HH11	1.82	0.43
1:B:9:GLY:CA	1:B:111:THR:HG22	2.48	0.43
1:C:133:ILE:C	1:C:135:ALA:N	2.76	0.43
1:D:89:LYS:HE2	4:D:875:HOH:O	2.19	0.43
1:A:5:VAL:HG12	1:A:6:GLY:N	2.34	0.43
1:B:280:ALA:HB1	1:B:285:MET:HG2	2.01	0.43
1:A:271:GLN:CD	1:A:271:GLN:N	2.76	0.43
1:A:328:LYS:HB2	1:B:189:LEU:HD13	2.00	0.43
1:B:268:THR:HB	1:B:269:SER:H	1.54	0.43
1:C:189:LEU:HA	1:C:244:THR:O	2.18	0.43
1:B:105:GLU:HB2	1:B:129:ARG:HB3	2.01	0.43
1:C:84:ARG:C	1:C:85:ILE:HD12	2.44	0.42
1:C:175:VAL:O	1:C:179:VAL:HG23	2.19	0.42
1:C:281:SER:O	1:C:286:LYS:HA	2.18	0.42
1:C:324:ARG:HG3	1:C:324:ARG:HH11	1.84	0.42
1:D:39:MET:HA	1:D:39:MET:CE	2.42	0.42
1:A:23:GLU:OE1	1:A:337:TRP:CZ2	2.72	0.42
1:A:268:THR:HB	1:A:269:SER:H	1.69	0.42
1:C:297:VAL:O	1:C:298:SER:C	2.62	0.42
1:A:189:LEU:HD23	1:A:244:THR:HG22	2.02	0.42
1:B:13:ILE:HG22	1:B:110:SER:HB2	2.01	0.42
1:B:31:ILE:CG2	1:B:32:ASP:N	2.82	0.42
1:B:38:ASP:HA	2:B:863:NAD:C2A	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:ASP:HB3	1:B:46:PHE:HZ	1.83	0.42
1:C:124:LEU:CD2	1:C:131:VAL:HG23	2.49	0.42
1:D:107:VAL:HG23	1:D:131:VAL:HG13	2.01	0.42
1:B:114:PHE:O	1:B:120:ALA:HB2	2.19	0.42
1:B:285:MET:CB	1:B:289:LEU:HB3	2.44	0.42
1:C:154:ASN:CG	1:C:157:GLU:HB2	2.44	0.42
1:A:5:VAL:CG1	1:A:6:GLY:N	2.81	0.42
1:B:12:ARG:O	1:B:16:MET:CG	2.62	0.42
1:B:93:ASN:HD22	1:B:94:PRO:CD	2.32	0.42
1:C:55:VAL:HG13	1:C:56:HIS:CD2	2.54	0.42
1:A:116:ALA:O	1:A:117:LYS:C	2.63	0.42
1:A:180:LYS:HD3	1:A:284:TYR:CE2	2.54	0.42
1:D:344:VAL:O	1:D:348:ARG:HG3	2.20	0.42
1:A:95:ALA:HB2	1:A:125:ARG:HD2	2.01	0.42
1:A:320:LEU:HB2	1:A:323:GLU:HB2	2.02	0.42
1:C:151:HIS:HB2	4:C:884:HOH:O	2.19	0.42
1:C:186:GLN:O	1:C:242:LYS:HE2	2.20	0.42
1:C:260:LEU:O	1:C:328:LYS:HA	2.20	0.42
1:C:17:VAL:HG11	1:C:108:ILE:HD13	2.02	0.42
1:C:178:LEU:O	1:C:183:PHE:HB2	2.20	0.42
1:B:124:LEU:HD22	1:B:129:ARG:O	2.19	0.42
1:C:107:VAL:HG21	1:C:123:HIS:ND1	2.35	0.42
1:D:321:PRO:O	1:D:322:LYS:HB2	2.20	0.42
1:B:129:ARG:NH1	1:B:355:ARG:NH2	2.68	0.42
1:B:153:TYR:CE2	1:B:350:MET:HA	2.55	0.42
1:C:144:LEU:HD12	1:C:346:LEU:HD11	2.02	0.41
1:C:342:ARG:HA	1:C:342:ARG:NE	2.35	0.41
1:D:34:VAL:HB	1:D:104:VAL:HG22	2.02	0.41
1:D:45:TYR:CD1	1:A:211:TRP:CE2	3.08	0.41
1:A:107:VAL:CG2	1:A:131:VAL:HG22	2.47	0.41
1:C:256:SER:HB2	1:C:333:TYR:CE1	2.55	0.41
1:D:130:LYS:CD	1:D:159:HIS:HA	2.49	0.41
1:B:38:ASP:OD1	1:B:39:MET:N	2.54	0.41
1:B:93:ASN:HD22	1:B:94:PRO:HD2	1.85	0.41
1:C:36:VAL:O	1:C:88:VAL:HG22	2.21	0.41
1:C:133:ILE:HD12	1:C:137:ALA:HB2	2.03	0.41
1:C:200:GLN:C	1:C:201:LYS:HD2	2.46	0.41
1:C:256:SER:HB2	1:C:333:TYR:CZ	2.55	0.41
1:C:332:TRP:O	1:C:333:TYR:HB3	2.20	0.41
1:A:17:VAL:HG11	1:A:108:ILE:HD13	2.02	0.41
1:C:32:ASP:CG	1:C:83:HIS:NE2	2.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:TYR:CD1	1:C:61:TYR:N	2.88	0.41
1:C:153:TYR:O	1:C:155:PRO:HD3	2.21	0.41
1:A:74:LYS:O	1:A:75:ASP:C	2.64	0.41
1:A:296:LEU:O	1:B:212:ARG:HD3	2.20	0.41
1:B:306:ARG:HB2	1:B:309:ILE:CD1	2.50	0.41
1:C:2:PRO:HB3	1:C:32:ASP:HB2	2.02	0.41
1:D:40:ASN:ND2	1:A:208:VAL:HG11	2.35	0.41
1:B:30:GLU:HG3	1:B:348:ARG:HH11	1.85	0.41
1:A:193:ILE:CG2	1:B:221:ILE:HD13	2.50	0.41
1:C:74:LYS:O	1:C:75:ASP:C	2.64	0.41
1:D:137:ALA:CB	1:D:162:SER:HB2	2.50	0.41
1:D:208:VAL:HG11	1:A:40:ASN:ND2	2.35	0.41
1:A:29:THR:HG22	1:A:30:GLU:N	2.35	0.41
1:A:297:VAL:O	1:A:298:SER:C	2.63	0.41
1:B:132:VAL:HG21	1:B:347:VAL:HG22	2.03	0.41
1:C:28:GLY:N	1:C:31:ILE:O	2.54	0.41
1:D:81:ASN:C	1:D:81:ASN:HD22	2.28	0.41
1:A:209:LYS:HA	1:A:209:LYS:HD3	1.86	0.41
1:C:137:ALA:CB	1:C:162:SER:HB2	2.51	0.41
1:C:149:ASN:CG	1:C:235:VAL:HG13	2.45	0.41
1:C:223:PRO:CG	1:D:332:TRP:HZ2	2.33	0.41
1:D:200:GLN:HB3	1:D:217:ALA:HB2	2.02	0.41
1:D:313:LYS:O	1:D:317:GLN:HG3	2.20	0.41
1:A:149:ASN:O	1:A:152:GLU:HG3	2.21	0.41
1:B:30:GLU:HG2	4:B:872:HOH:O	2.20	0.41
1:B:137:ALA:HB3	1:B:162:SER:HB2	2.03	0.41
1:C:186:GLN:HG2	4:C:903:HOH:O	2.20	0.41
1:D:58:LYS:HB3	4:D:887:HOH:O	2.20	0.41
1:D:99:TRP:CE3	1:D:99:TRP:HA	2.56	0.41
1:A:125:ARG:C	1:A:127:GLY:N	2.79	0.40
1:B:122:GLY:HA2	1:B:125:ARG:HE	1.86	0.40
1:B:169:ASN:O	1:B:307:SER:HB3	2.22	0.40
1:C:107:VAL:HG11	1:C:123:HIS:HB3	2.02	0.40
1:A:27:LEU:HD11	1:A:83:HIS:ND1	2.37	0.40
1:B:13:ILE:HD13	1:B:13:ILE:HA	1.93	0.40
1:A:36:VAL:HG13	1:A:85:ILE:HG21	2.02	0.40
1:A:305:ASN:OD1	1:A:337:TRP:NE1	2.54	0.40
1:A:16:MET:SD	1:A:337:TRP:CZ3	3.15	0.40
1:A:290:GLY:HA3	1:A:306:ARG:NH1	2.37	0.40
1:B:323:GLU:HG2	1:B:326:PHE:O	2.21	0.40
1:C:188:GLY:CA	1:C:243:LEU:HD23	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:326:PHE:CZ	1:D:262:PHE:HA	2.56	0.40
1:D:12:ARG:O	1:D:16:MET:HG2	2.21	0.40
1:A:4:LYS:CE	1:A:32:ASP:HB3	2.52	0.40
1:A:221:ILE:HD13	1:B:193:ILE:CG2	2.52	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	357/359 (99%)	317 (89%)	33 (9%)	7 (2%)	6	11
1	B	357/359 (99%)	320 (90%)	31 (9%)	6 (2%)	7	14
1	C	357/359 (99%)	310 (87%)	40 (11%)	7 (2%)	6	11
1	D	357/359 (99%)	317 (89%)	39 (11%)	1 (0%)	36	56
All	All	1428/1436 (99%)	1264 (88%)	143 (10%)	21 (2%)	8	15

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	90	ALA
1	C	29	THR
1	C	157	GLU
1	C	255	VAL
1	D	255	VAL
1	A	37	VAL
1	A	255	VAL
1	A	298	SER
1	B	96	ASP
1	B	255	VAL
1	C	134	SER

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Mol	Chain	Res	Type
1	A	90	ALA
1	C	335	ASN
1	A	164	ALA
1	B	42	ASP
1	B	228	ALA
1	C	216	ALA
1	A	73	ALA
1	B	351	ALA
1	A	237	PRO
1	C	37	VAL

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	293/296 (99%)	280 (96%)	13 (4%)	25	47
1	B	293/296 (99%)	281 (96%)	12 (4%)	27	49
1	C	293/296 (99%)	282 (96%)	11 (4%)	29	52
1	D	293/296 (99%)	277 (94%)	16 (6%)	19	37
All	All	1172/1184 (99%)	1120 (96%)	52 (4%)	25	47

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	4	LYS
1	C	30	GLU
1	C	37	VAL
1	C	157	GLU
1	C	171	LEU
1	C	187	THR
1	C	190	MET
1	C	238	SER
1	C	305	ASN
1	C	309	ILE
1	C	355	ARG

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Mol	Chain	Res	Type
1	D	12	ARG
1	D	30	GLU
1	D	37	VAL
1	D	39	MET
1	D	60	LYS
1	D	78	LEU
1	D	89	LYS
1	D	97	LEU
1	D	171	LEU
1	D	187	THR
1	D	190	MET
1	D	192	THR
1	D	267	ASP
1	D	305	ASN
1	D	309	ILE
1	D	355	ARG
1	A	37	VAL
1	A	55	VAL
1	A	60	LYS
1	A	89	LYS
1	A	109	GLU
1	A	157	GLU
1	A	171	LEU
1	A	190	MET
1	A	230	LYS
1	A	237	PRO
1	A	309	ILE
1	A	319	ASN
1	A	335	ASN
1	B	13	ILE
1	B	16	MET
1	B	30	GLU
1	B	37	VAL
1	B	55	VAL
1	B	91	GLN
1	B	171	LEU
1	B	190	MET
1	B	230	LYS
1	B	305	ASN
1	B	309	ILE
1	B	355	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32)

such sidechains are listed below:

Mol	Chain	Res	Type
1	C	19	GLN
1	C	81	ASN
1	C	93	ASN
1	C	150	HIS
1	C	271	GLN
1	C	317	GLN
1	C	319	ASN
1	C	349	HIS
1	D	8	ASN
1	D	81	ASN
1	D	93	ASN
1	D	150	HIS
1	D	151	HIS
1	D	176	HIS
1	D	303	ASN
1	D	305	ASN
1	D	319	ASN
1	A	19	GLN
1	A	81	ASN
1	A	93	ASN
1	A	150	HIS
1	A	151	HIS
1	A	154	ASN
1	A	317	GLN
1	A	319	ASN
1	B	19	GLN
1	B	81	ASN
1	B	93	ASN
1	B	150	HIS
1	B	159	HIS
1	B	305	ASN
1	B	319	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	S70	A	804	-	13,14,14	3.35	10 (76%)	15,21,21	1.28	3 (20%)
2	NAD	D	861	-	46,48,48	1.65	6 (13%)	64,73,73	1.61	8 (12%)
2	NAD	B	863	-	46,48,48	1.64	6 (13%)	64,73,73	1.60	8 (12%)
2	NAD	A	862	-	46,48,48	1.61	7 (15%)	64,73,73	1.62	9 (14%)
2	NAD	C	860	-	46,48,48	1.68	6 (13%)	64,73,73	1.62	9 (14%)

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
3	S70	A	804	-	-	3/14/15/15	-
2	NAD	D	861	-	-	13/30/62/62	0/5/5/5
2	NAD	B	863	-	-	10/30/62/62	0/5/5/5
2	NAD	A	862	-	-	8/30/62/62	0/5/5/5
2	NAD	C	860	-	-	9/30/62/62	0/5/5/5

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	804	S70	C2-C3	6.16	1.63	1.53
3	A	804	S70	P8-C4	-5.51	1.71	1.79
3	A	804	S70	P8-O10	5.47	1.61	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	860	NAD	C3N-C7N	5.16	1.58	1.50
2	B	863	NAD	C3N-C7N	5.10	1.58	1.50
2	D	861	NAD	C3N-C7N	5.06	1.58	1.50
2	A	862	NAD	C3N-C7N	5.05	1.58	1.50
2	C	860	NAD	C2N-N1N	5.00	1.40	1.35
2	D	861	NAD	C2N-N1N	4.99	1.40	1.35
2	B	863	NAD	C2N-N1N	4.95	1.40	1.35
2	A	862	NAD	C2N-N1N	4.52	1.40	1.35
2	D	861	NAD	C6N-N1N	4.13	1.44	1.35
2	B	863	NAD	C6N-N1N	3.93	1.44	1.35
2	C	860	NAD	C6N-N1N	3.92	1.44	1.35
2	C	860	NAD	PA-O3	3.82	1.63	1.59
2	A	862	NAD	C6N-N1N	3.78	1.43	1.35
2	D	861	NAD	C4N-C3N	3.70	1.45	1.39
2	C	860	NAD	C4N-C3N	3.36	1.44	1.39
2	A	862	NAD	C4N-C3N	3.29	1.44	1.39
3	A	804	S70	P12-O15	3.22	1.66	1.54
2	B	863	NAD	C4N-C3N	3.14	1.44	1.39
3	A	804	S70	P12-O13	3.04	1.59	1.50
2	A	862	NAD	PA-O3	2.82	1.62	1.59
2	A	862	NAD	C2A-N1A	2.68	1.38	1.33
2	B	863	NAD	PA-O3	2.67	1.62	1.59
2	B	863	NAD	C2A-N1A	2.56	1.38	1.33
3	A	804	S70	P8-O9	2.56	1.60	1.55
2	C	860	NAD	C2A-N1A	2.50	1.38	1.33
2	D	861	NAD	PA-O3	2.39	1.62	1.59
2	D	861	NAD	C2A-N1A	2.39	1.38	1.33
3	A	804	S70	P12-O14	2.18	1.62	1.54
3	A	804	S70	C4-C3	2.10	1.53	1.51
3	A	804	S70	P8-O11	2.09	1.59	1.55
2	A	862	NAD	PA-O2A	-2.05	1.45	1.55
3	A	804	S70	O6-C3	2.01	1.25	1.21

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	862	NAD	C5N-C4N-C3N	-6.25	114.22	120.36
2	D	861	NAD	C5N-C4N-C3N	-6.14	114.33	120.36
2	C	860	NAD	C5N-C4N-C3N	-6.00	114.47	120.36
2	B	863	NAD	C5N-C4N-C3N	-5.90	114.56	120.36
2	A	862	NAD	C6N-C5N-C4N	5.49	127.35	119.45
2	C	860	NAD	C6N-C5N-C4N	5.45	127.30	119.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	863	NAD	C6N-C5N-C4N	5.36	127.17	119.45
2	D	861	NAD	C6N-C5N-C4N	5.34	127.14	119.45
2	A	862	NAD	C5N-C6N-N1N	-4.73	113.93	120.38
2	B	863	NAD	C5N-C6N-N1N	-4.65	114.04	120.38
2	C	860	NAD	C5N-C6N-N1N	-4.64	114.05	120.38
2	D	861	NAD	C5N-C6N-N1N	-4.28	114.55	120.38
2	B	863	NAD	C4N-C3N-C7N	-3.93	110.36	121.06
2	D	861	NAD	C4N-C3N-C7N	-3.91	110.42	121.06
2	C	860	NAD	C4N-C3N-C7N	-3.75	110.83	121.06
2	A	862	NAD	C4N-C3N-C7N	-3.68	111.03	121.06
2	D	861	NAD	C2N-C3N-C4N	3.49	122.32	118.26
2	A	862	NAD	C2N-C3N-C4N	3.22	122.00	118.26
2	B	863	NAD	C2N-C3N-C4N	3.20	121.98	118.26
2	C	860	NAD	C2N-C3N-C4N	3.14	121.91	118.26
2	B	863	NAD	C2N-C3N-C7N	2.84	127.66	119.46
2	D	861	NAD	C2N-C3N-C7N	2.70	127.26	119.46
2	C	860	NAD	C2N-C3N-C7N	2.70	127.26	119.46
2	A	862	NAD	C2N-C3N-C7N	2.60	126.97	119.46
2	B	863	NAD	C3B-C2B-C1B	2.60	106.38	101.46
2	C	860	NAD	O4B-C1B-N9A	-2.41	103.45	108.09
2	C	860	NAD	C3B-C2B-C1B	2.39	105.99	101.46
2	A	862	NAD	O4B-C1B-C2B	-2.28	101.74	106.62
3	A	804	S70	O14-P12-O7	2.27	112.58	106.67
2	A	862	NAD	C3B-C2B-C1B	2.25	105.73	101.46
2	D	861	NAD	O4B-C1B-C2B	-2.20	101.90	106.62
2	D	861	NAD	O4B-C1B-N9A	-2.17	103.92	108.09
2	B	863	NAD	O4B-C1B-N9A	-2.14	103.97	108.09
3	A	804	S70	O11-P8-C4	2.13	111.31	106.84
2	C	860	NAD	O4B-C1B-C2B	-2.10	102.12	106.62
3	A	804	S70	O6-C3-C2	2.04	122.17	120.07
2	A	862	NAD	O4B-C1B-N9A	-2.04	104.17	108.09

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	860	NAD	O4D-C1D-N1N-C2N
2	C	860	NAD	O4D-C1D-N1N-C6N
2	C	860	NAD	C2D-C1D-N1N-C2N
2	C	860	NAD	C2D-C1D-N1N-C6N
2	D	861	NAD	C5B-O5B-PA-O1A
2	D	861	NAD	O4D-C1D-N1N-C2N

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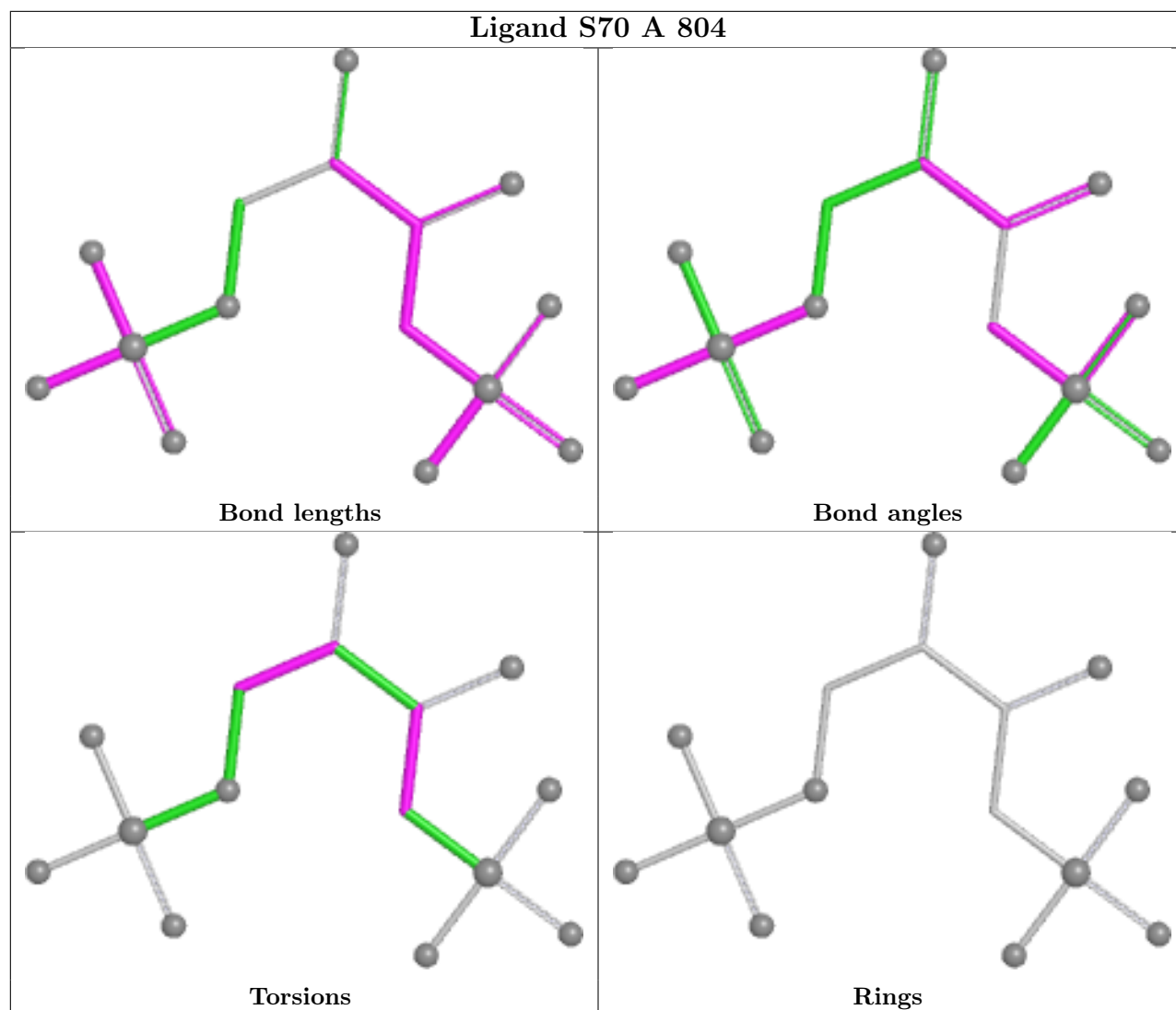
Mol	Chain	Res	Type	Atoms
2	D	861	NAD	O4D-C1D-N1N-C6N
2	D	861	NAD	C2D-C1D-N1N-C2N
2	D	861	NAD	C2D-C1D-N1N-C6N
2	A	862	NAD	O4D-C1D-N1N-C2N
2	A	862	NAD	O4D-C1D-N1N-C6N
2	A	862	NAD	C2D-C1D-N1N-C2N
2	A	862	NAD	C2D-C1D-N1N-C6N
2	A	862	NAD	C2N-C3N-C7N-O7N
2	A	862	NAD	C2N-C3N-C7N-N7N
2	B	863	NAD	O4D-C1D-N1N-C2N
2	B	863	NAD	O4D-C1D-N1N-C6N
2	B	863	NAD	C2D-C1D-N1N-C2N
2	B	863	NAD	C2D-C1D-N1N-C6N
2	B	863	NAD	C2N-C3N-C7N-O7N
2	B	863	NAD	C2N-C3N-C7N-N7N
3	A	804	S70	O7-C1-C2-C3
3	A	804	S70	O7-C1-C2-O5
2	B	863	NAD	C4N-C3N-C7N-O7N
2	B	863	NAD	C4N-C3N-C7N-N7N
2	C	860	NAD	C2N-C3N-C7N-O7N
2	A	862	NAD	C4N-C3N-C7N-N7N
2	A	862	NAD	C4N-C3N-C7N-O7N
2	C	860	NAD	C2N-C3N-C7N-N7N
2	C	860	NAD	C4N-C3N-C7N-N7N
2	C	860	NAD	C4N-C3N-C7N-O7N
2	B	863	NAD	O4B-C4B-C5B-O5B
2	B	863	NAD	C3B-C4B-C5B-O5B
3	A	804	S70	O6-C3-C4-P8
2	C	860	NAD	C2B-C1B-N9A-C8A
2	D	861	NAD	C2B-C1B-N9A-C8A
2	D	861	NAD	C5B-O5B-PA-O2A
2	D	861	NAD	C5B-O5B-PA-O3
2	D	861	NAD	O4B-C4B-C5B-O5B
2	D	861	NAD	C4N-C3N-C7N-N7N
2	D	861	NAD	C2B-C1B-N9A-C4A
2	D	861	NAD	C4N-C3N-C7N-O7N
2	D	861	NAD	C3B-C4B-C5B-O5B

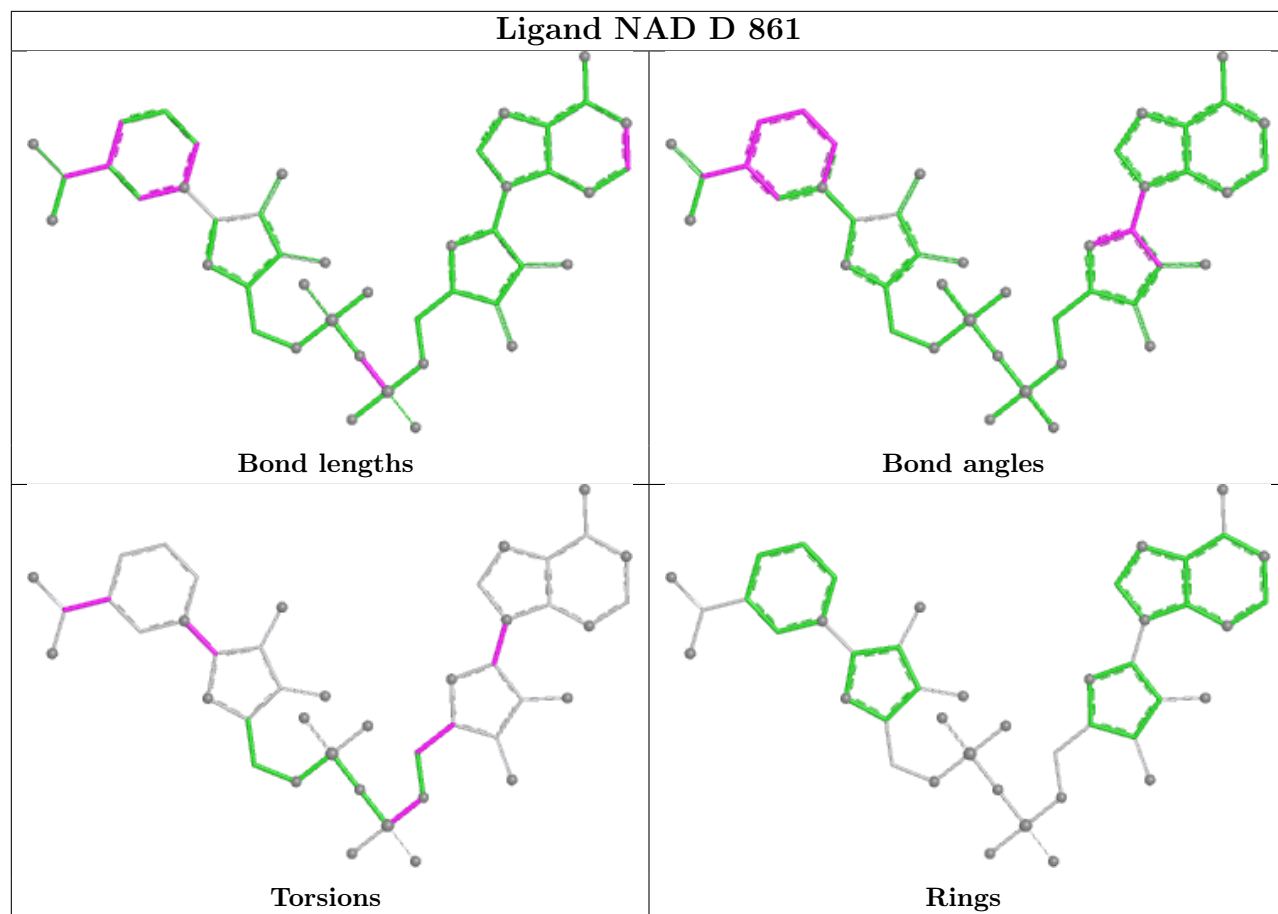
There are no ring outliers.

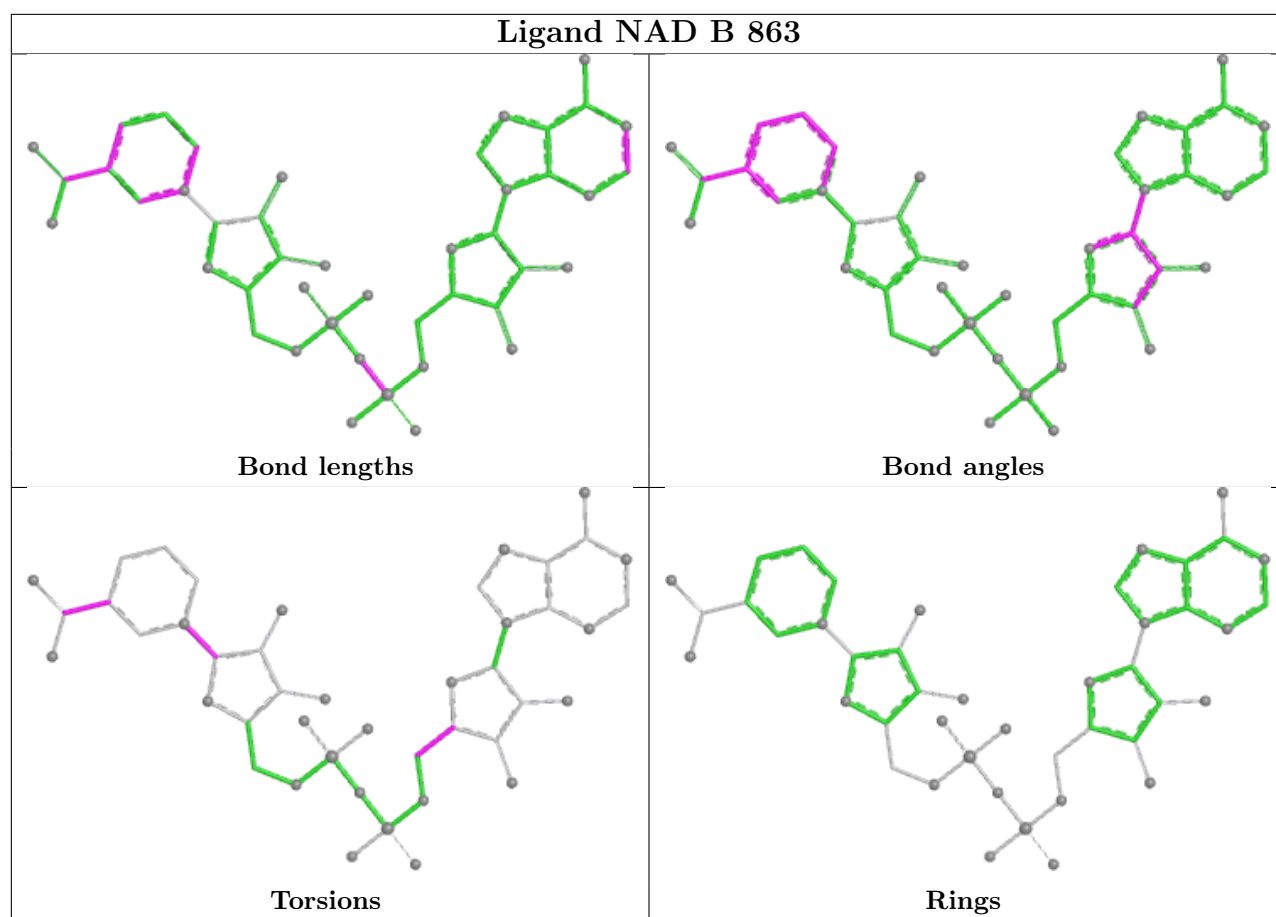
3 monomers are involved in 9 short contacts:

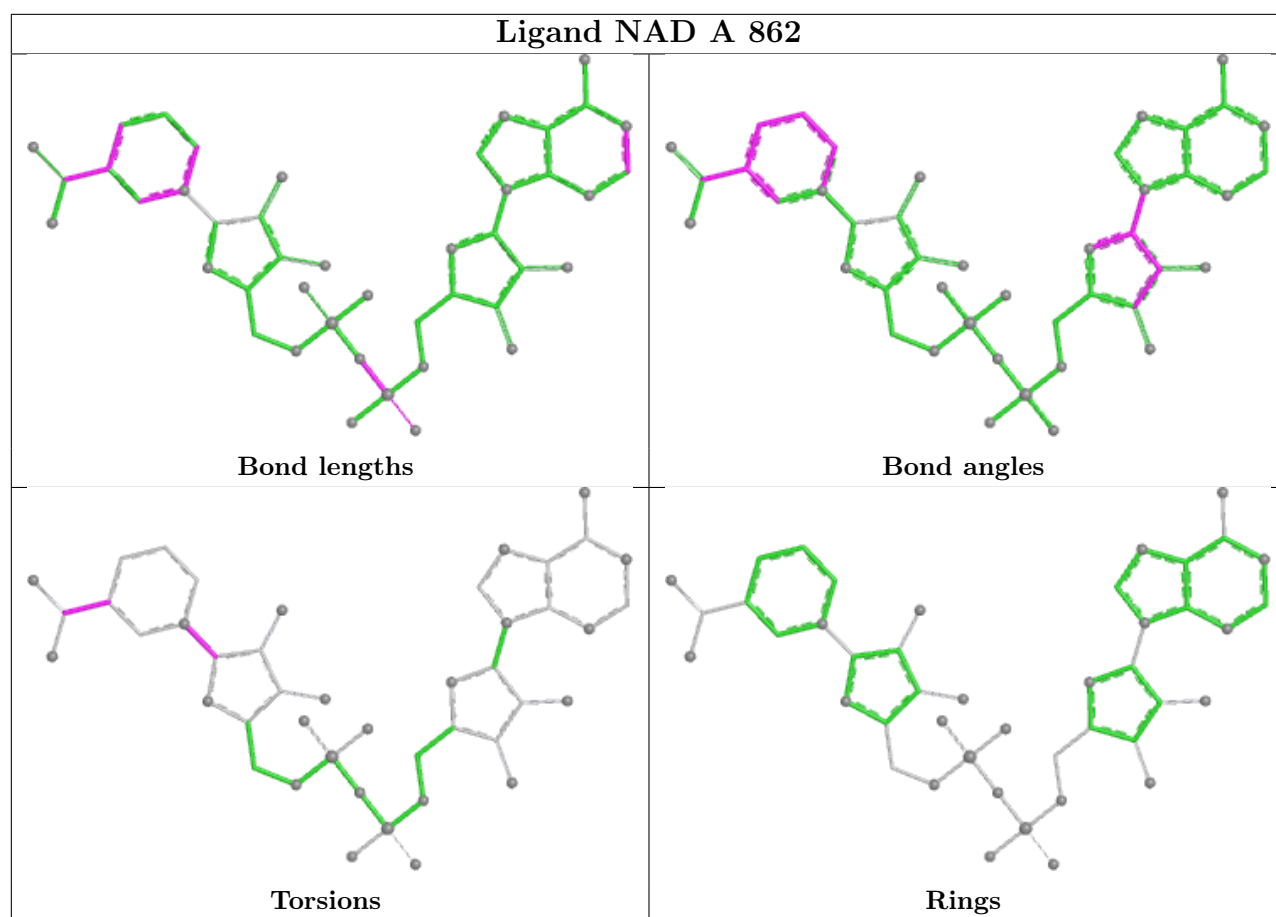
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	804	S70	1	0
2	B	863	NAD	6	0
2	A	862	NAD	2	0

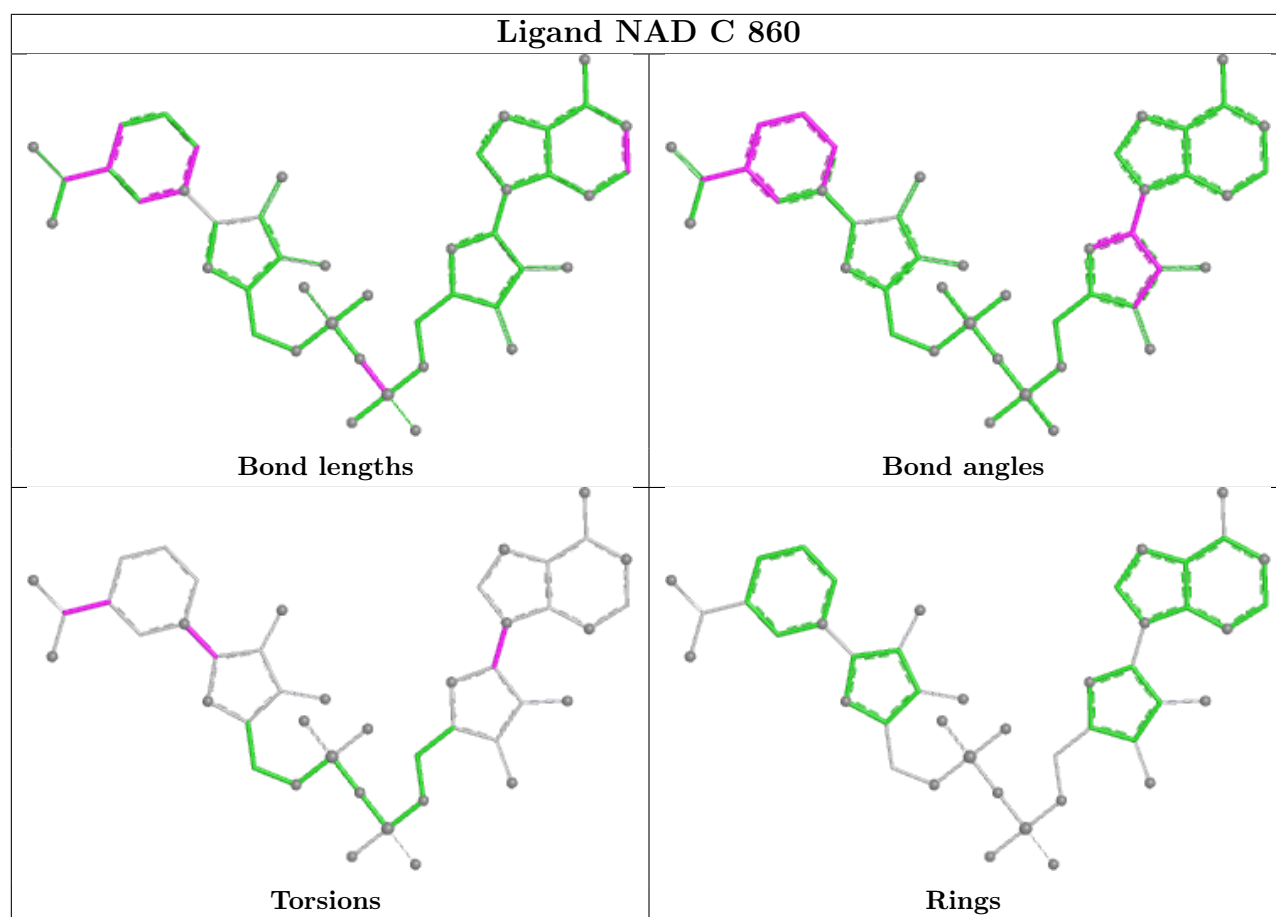
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	359/359 (100%)	-0.62	0 100 100	30, 51, 86, 102	0
1	B	359/359 (100%)	-0.52	0 100 100	29, 52, 95, 106	0
1	C	359/359 (100%)	-0.46	0 100 100	33, 59, 97, 108	0
1	D	359/359 (100%)	-0.69	0 100 100	32, 51, 79, 95	0
All	All	1436/1436 (100%)	-0.57	0 100 100	29, 53, 90, 108	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

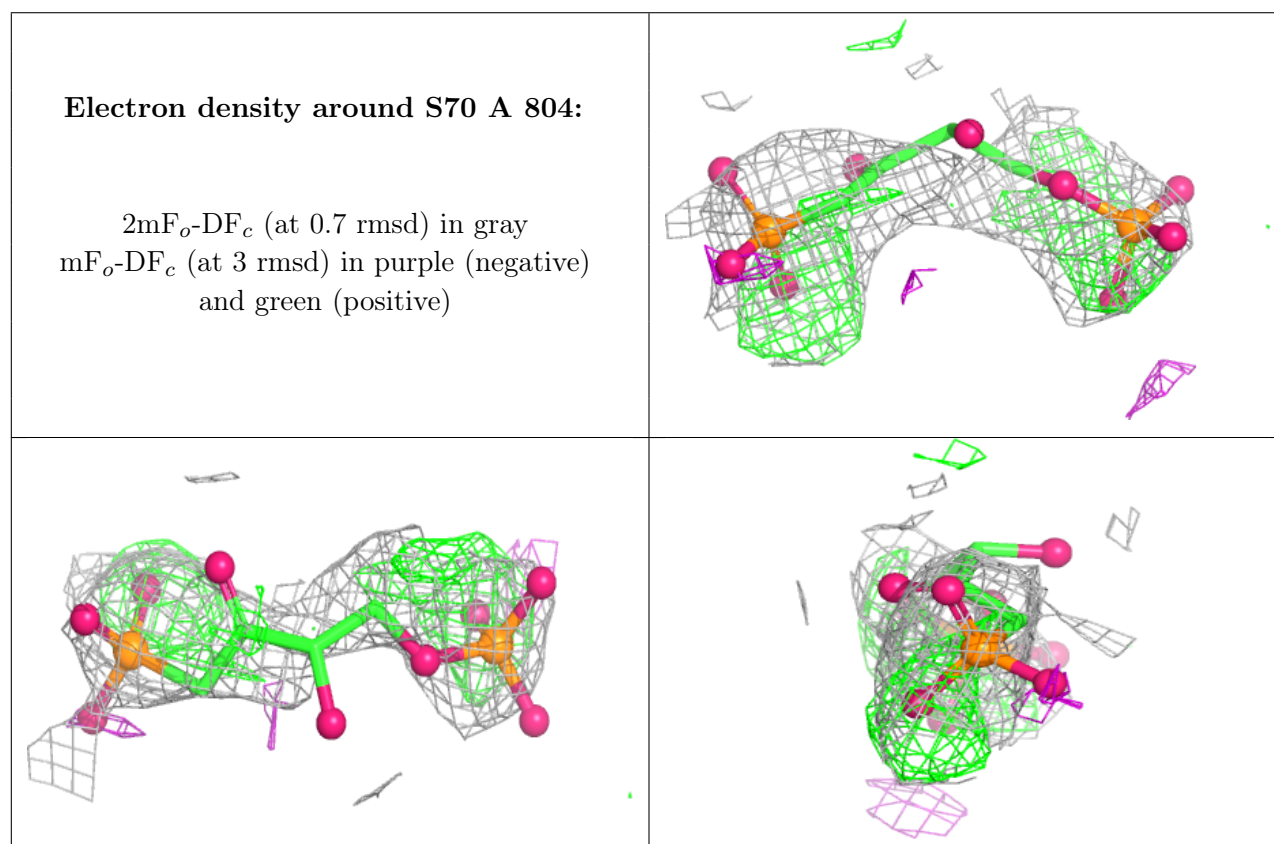
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	S70	A	804	15/15	0.78	0.24	84,86,86,86	15
2	NAD	B	863	44/44	0.83	0.12	82,88,91,92	44
2	NAD	A	862	44/44	0.83	0.11	66,71,74,74	0
2	NAD	C	860	44/44	0.90	0.08	66,77,81,82	0

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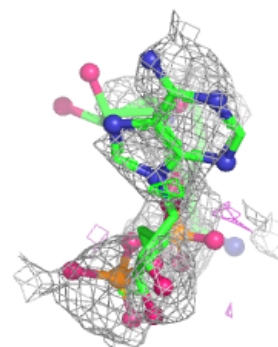
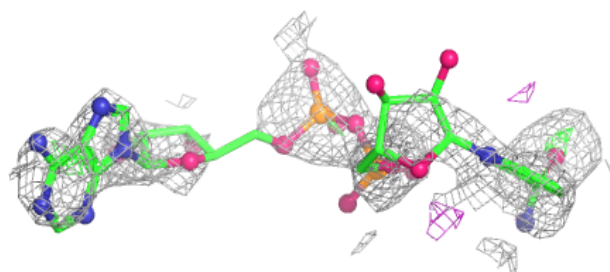
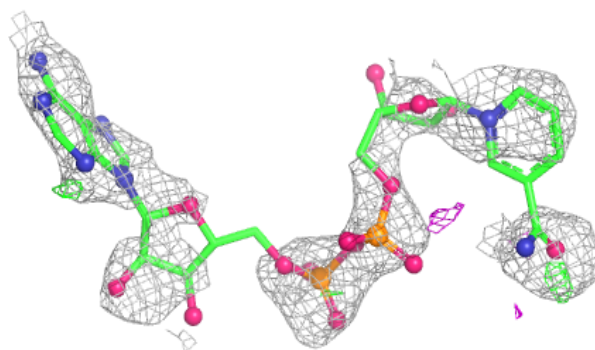
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAD	D	861	44/44	0.94	0.07	51,53,55,55	0

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.

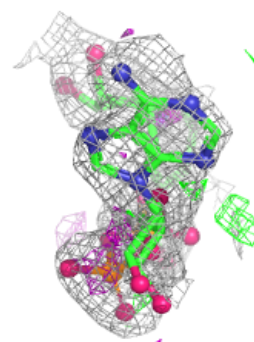
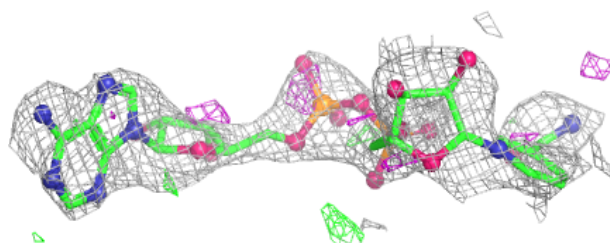
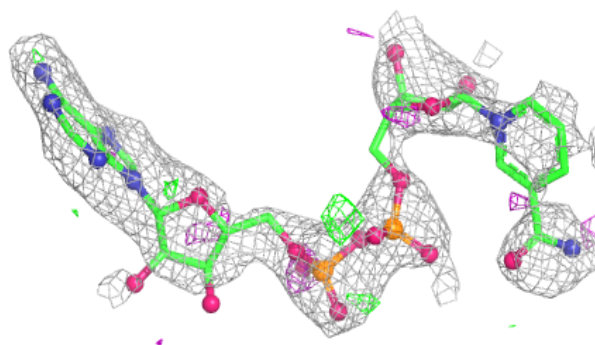


Electron density around NAD B 863:

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

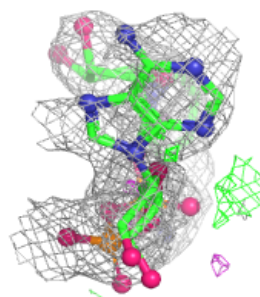
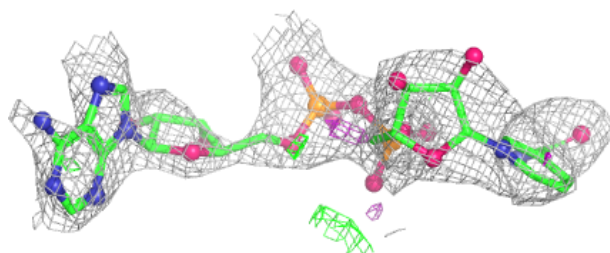
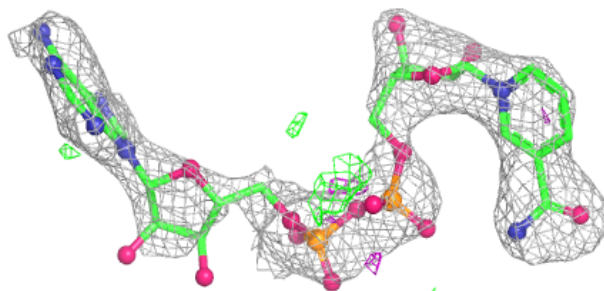
**Electron density around NAD A 862:**

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

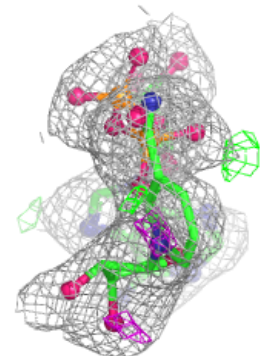
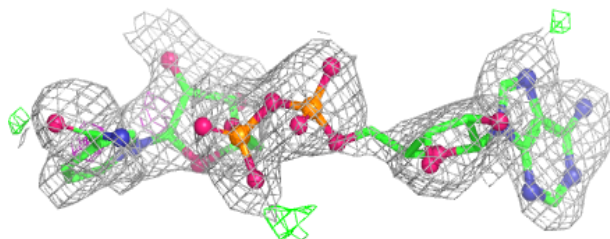
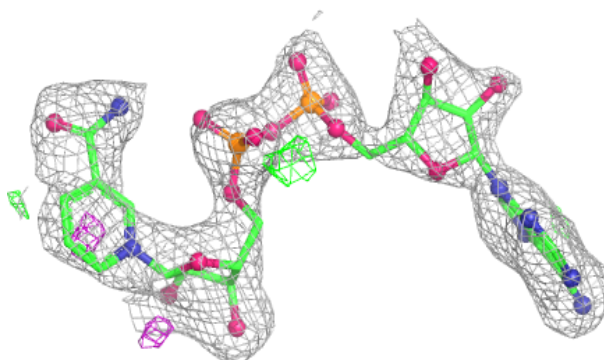


Electron density around NAD C 860:

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

**Electron density around NAD D 861:**

$2mF_o-DF_c$ (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.