



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:03 AM UTC

PDB ID : 1PIW / pdb_00001piw
Title : APO AND HOLO STRUCTURES OF AN NADP(H)-DEPENDENT CIN-
NAMYL ALCOHOL DEHYDROGENASE FROM SACCHAROMYCES
CEREVISIAE
Authors : Valencia, E.; Larroy, C.; Ochoa, W.F.; Pares, X.; Fita, I.; Biosca, J.A.
Deposited on : 2003-05-30
Resolution : 3.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

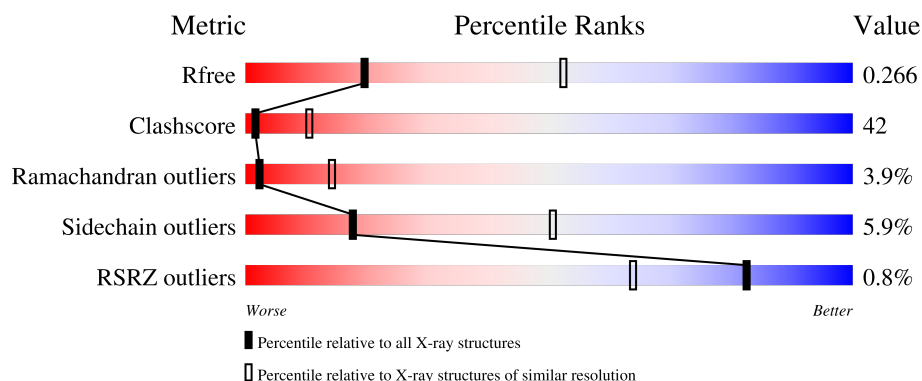
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	360	46% 47% 6%
1	B	360	34% 55% 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAP	B	4293	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5660 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.

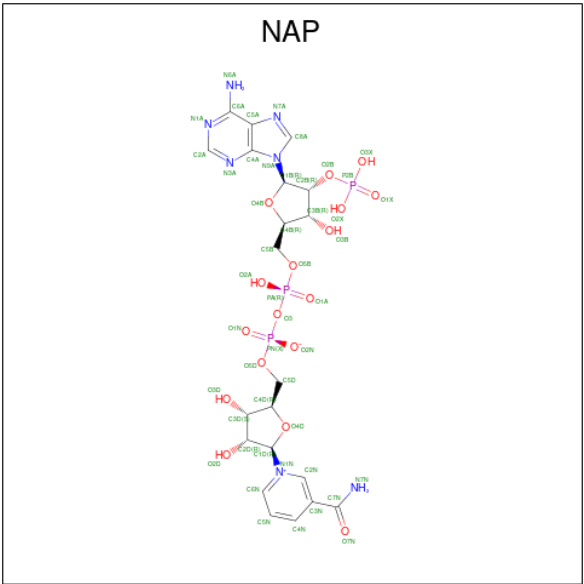
- Molecule 1 is a protein called Hypothetical zinc-type alcohol dehydrogenase-like protein in PRE5-FET4 intergenic region.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	0	0
			2780	1770	466	523	21			
1	B	360	Total	C	N	O	S	0	0	0
			2780	1770	466	523	21			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).

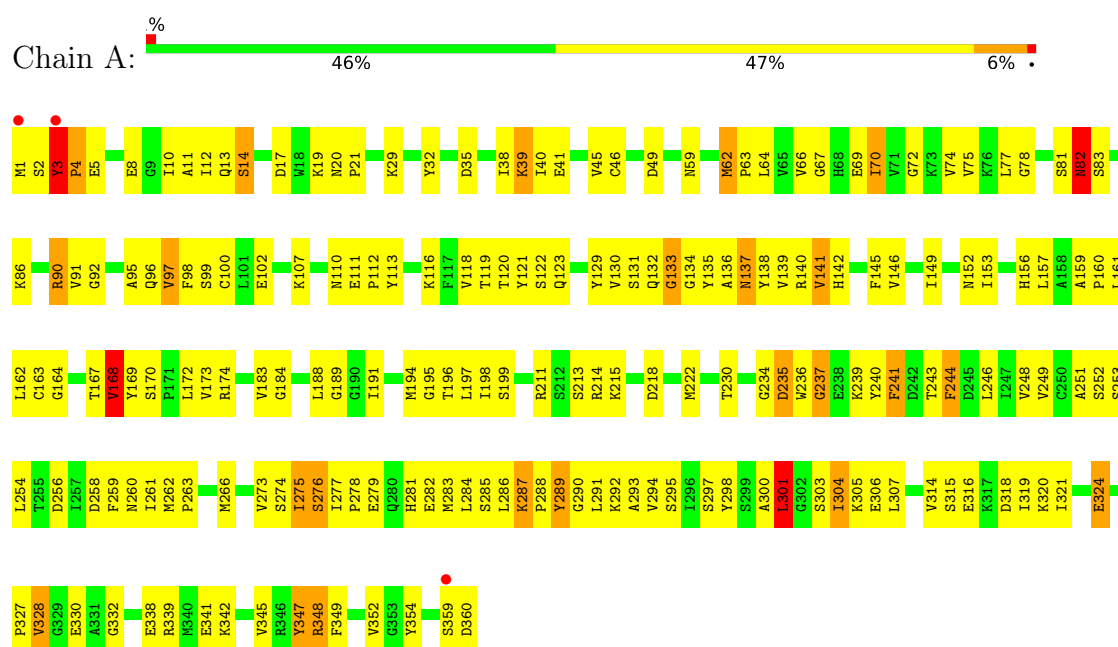


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
3	B	1	Total 48	C 21	N 7	O 17	P 3	0	0

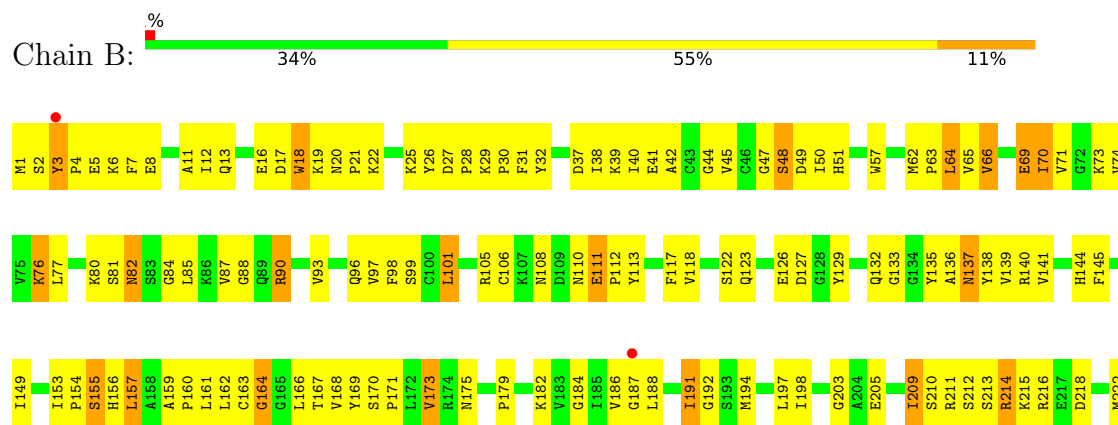
3 Residue-property plots

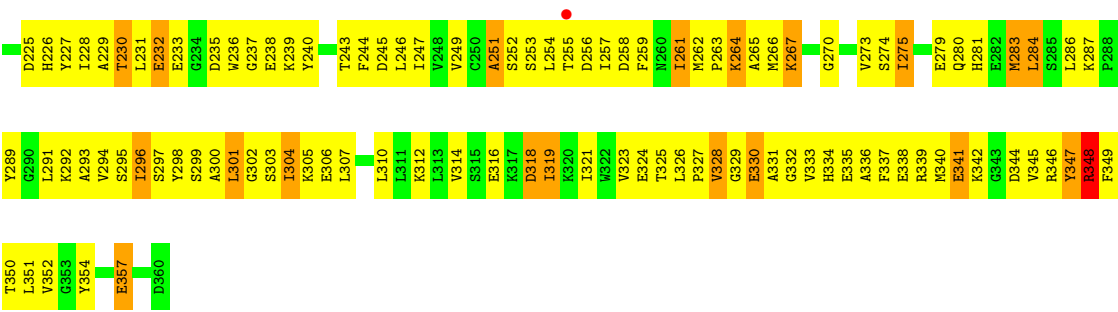
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Hypothetical zinc-type alcohol dehydrogenase-like protein in PRE5-FET4 intergenic region



- Molecule 1: Hypothetical zinc-type alcohol dehydrogenase-like protein in PRE5-FET4 intergenic region





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	234.98Å 100.35Å 52.49Å 90.00° 92.18° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 30.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-3.00) 98.7 (30.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 3.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.218 , 0.271 0.219 , 0.266	Depositor DCC
R_{free} test set	2393 reflections (9.87%)	wwPDB-VP
Wilson B-factor (Å ²)	60.7	Xtriage
Anisotropy	0.762	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 52.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5660	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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: ZN, NAP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.51	0/2844	1.10	18/3839 (0.5%)
1	B	0.43	0/2844	0.99	9/3839 (0.2%)
All	All	0.47	0/5688	1.05	27/7678 (0.4%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	TYR	CA-C-N	-8.74	108.91	119.84
1	A	3	TYR	C-N-CA	-8.74	108.91	119.84
1	B	70	ILE	N-CA-C	8.01	119.31	109.30
1	A	70	ILE	N-CA-C	7.30	118.53	108.89
1	A	276	SER	N-CA-C	7.21	120.29	110.55
1	B	348	ARG	N-CA-C	6.81	119.84	108.13
1	B	245	ASP	N-CA-C	-6.56	103.33	112.45
1	A	168	VAL	CB-CA-C	-6.56	102.41	111.65
1	A	152	ASN	N-CA-C	6.52	120.23	112.93
1	A	287	LYS	CA-C-N	6.40	127.83	119.84
1	A	287	LYS	C-N-CA	6.40	127.83	119.84
1	A	17	ASP	N-CA-C	-6.14	105.63	112.57
1	A	11	ALA	N-CA-C	5.88	119.05	109.59
1	B	84	GLY	N-CA-C	-5.58	107.32	115.63
1	A	244	PHE	N-CA-C	5.55	119.28	109.96
1	B	69	GLU	N-CA-C	-5.35	104.02	110.44
1	A	78	GLY	CA-C-N	5.33	126.50	119.84
1	A	78	GLY	C-N-CA	5.33	126.50	119.84
1	B	49	ASP	N-CA-C	-5.29	104.92	111.33
1	B	64	LEU	N-CA-C	5.29	116.65	107.93
1	A	39	LYS	N-CA-C	-5.25	100.62	108.96
1	A	239	LYS	N-CA-C	5.17	116.99	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	232	GLU	N-CA-C	5.16	116.98	111.36
1	A	237	GLY	N-CA-C	-5.16	106.52	112.50
1	B	209	ILE	N-CA-C	5.11	114.84	106.72
1	A	141	VAL	N-CA-C	5.05	115.92	108.45
1	A	324	GLU	N-CA-C	-5.01	101.71	109.72

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	2780	0	2758	180	0
1	B	2780	0	2758	299	1
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	48	0	25	11	0
3	B	48	0	25	27	0
All	All	5660	0	5566	467	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 42.

All (467) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:SER:HB2	3:B:4293:NAP:C6A	1.26	1.56
1:B:252:SER:CB	3:B:4293:NAP:C6A	2.23	1.17
1:B:300:ALA:HA	3:B:4293:NAP:H72N	1.15	1.08
1:B:252:SER:HB2	3:B:4293:NAP:N6A	1.69	1.06
1:B:301:LEU:HD23	3:B:4293:NAP:O7N	1.53	1.06
1:B:8:GLU:HB2	1:B:25:LYS:HG2	1.39	1.02
1:B:97:VAL:HG23	1:B:141:VAL:HG13	1.39	0.99
1:B:251:ALA:HB1	3:B:4293:NAP:N3A	1.75	0.98
1:B:300:ALA:HA	3:B:4293:NAP:N7N	1.81	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:LEU:N	3:B:4293:NAP:O7N	1.98	0.94
1:B:252:SER:HB2	3:B:4293:NAP:N1A	1.82	0.93
1:B:6:LYS:HA	1:B:28:PRO:HD3	1.50	0.92
1:B:62:MET:HG3	1:B:63:PRO:HA	1.52	0.92
1:B:249:VAL:HG21	1:B:257:ILE:HD11	1.53	0.90
1:B:251:ALA:HB1	3:B:4293:NAP:C2A	2.01	0.90
1:B:304:ILE:HD13	1:B:305:LYS:H	1.35	0.89
1:B:1:MET:SD	1:B:5:GLU:HG3	2.13	0.89
1:B:328:VAL:O	1:B:357:GLU:HG2	1.80	0.80
1:B:191:ILE:HD12	1:B:191:ILE:H	1.47	0.80
1:A:314:VAL:HA	1:A:319:ILE:HG12	1.62	0.79
1:A:168:VAL:HG22	1:A:191:ILE:HG23	1.65	0.79
1:A:97:VAL:HG22	1:A:141:VAL:HG13	1.65	0.79
1:A:279:GLU:OE2	1:A:281:HIS:HB2	1.83	0.78
1:B:252:SER:HB2	3:B:4293:NAP:C5A	2.10	0.78
1:B:314:VAL:HA	1:B:319:ILE:HD11	1.64	0.77
1:B:210:SER:O	1:B:229:ALA:HA	1.84	0.76
1:B:252:SER:H	3:B:4293:NAP:C2A	1.96	0.76
1:B:341:GLU:O	1:B:341:GLU:HG2	1.84	0.76
1:B:168:VAL:CG2	1:B:191:ILE:HG22	2.16	0.76
1:B:264:LYS:HB2	1:B:264:LYS:NZ	2.00	0.76
1:B:90:ARG:NH2	1:B:153:ILE:O	2.19	0.76
1:A:159:ALA:HB3	1:A:160:PRO:HD3	1.67	0.76
1:B:188:LEU:HD23	1:B:215:LYS:HB3	1.68	0.76
1:B:211:ARG:HB3	1:B:230:THR:HB	1.68	0.76
1:B:240:TYR:O	1:B:243:THR:HG22	1.84	0.76
1:A:314:VAL:HG13	1:A:319:ILE:HB	1.68	0.75
1:B:186:VAL:HG22	1:B:209:ILE:HB	1.67	0.75
1:A:304:ILE:HD12	1:A:305:LYS:H	1.51	0.74
1:B:166:LEU:HD22	1:B:307:LEU:HD13	1.69	0.74
1:B:167:THR:O	1:B:275:ILE:HD11	1.87	0.74
1:B:255:THR:HG22	1:B:284:LEU:HG	1.70	0.74
1:B:252:SER:CB	3:B:4293:NAP:N6A	2.44	0.74
1:A:240:TYR:O	1:A:243:THR:HG22	1.88	0.73
1:B:303:SER:OG	1:B:306:GLU:HG3	1.88	0.73
1:B:170:SER:HB3	1:B:171:PRO:HD3	1.69	0.73
1:A:66:VAL:HG12	1:A:67:GLY:H	1.53	0.73
1:A:81:SER:C	1:A:82:ASN:HD22	1.97	0.72
1:A:287:LYS:HD2	1:B:281:HIS:HA	1.71	0.72
1:B:62:MET:CG	1:B:63:PRO:HA	2.19	0.72
1:A:252:SER:OG	3:A:4292:NAP:H51A	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:VAL:HG12	1:A:67:GLY:N	2.05	0.71
1:B:247:ILE:HG21	1:B:262:MET:HE2	1.73	0.70
1:A:95:ALA:HB3	1:A:301:LEU:HD22	1.74	0.70
1:B:252:SER:CB	3:B:4293:NAP:N1A	2.46	0.70
1:B:156:HIS:HA	1:B:350:THR:HG21	1.74	0.69
1:B:17:ASP:OD2	1:B:20:ASN:HB2	1.93	0.68
1:A:345:VAL:C	1:A:347:TYR:H	2.01	0.68
1:B:66:VAL:HG23	1:B:135:TYR:CE2	2.29	0.68
1:B:244:PHE:O	1:B:267:LYS:HG2	1.94	0.67
1:B:64:LEU:CD2	1:B:122:SER:HB3	2.24	0.67
1:B:118:VAL:HA	1:B:123:GLN:NE2	2.10	0.67
1:B:137:ASN:HD22	1:B:138:TYR:N	1.92	0.67
1:B:26:TYR:CE2	1:B:65:VAL:HG21	2.30	0.67
1:A:290:GLY:C	1:A:292:LYS:H	2.03	0.67
1:A:211:ARG:NH1	3:A:4292:NAP:O1X	2.28	0.67
1:B:314:VAL:HG13	1:B:319:ILE:HG12	1.76	0.66
1:B:211:ARG:HB2	1:B:231:LEU:HG	1.78	0.66
1:B:133:GLY:HA3	1:B:139:VAL:HG23	1.77	0.66
1:B:251:ALA:CB	3:B:4293:NAP:N3A	2.55	0.66
1:B:137:ASN:HD22	1:B:138:TYR:H	1.43	0.66
1:B:251:ALA:HA	3:B:4293:NAP:H1B	1.77	0.66
1:B:69:GLU:OE2	1:B:163:CYS:HB3	1.96	0.66
1:A:314:VAL:HA	1:A:319:ILE:CG1	2.25	0.66
1:B:118:VAL:HA	1:B:123:GLN:HE22	1.61	0.66
1:B:188:LEU:HD11	1:B:222:MET:HE1	1.78	0.65
1:B:105:ARG:O	1:B:110:ASN:HB2	1.96	0.65
1:B:314:VAL:HA	1:B:319:ILE:CD1	2.26	0.65
1:B:213:SER:O	1:B:216:ARG:HB3	1.97	0.65
1:B:64:LEU:HD21	1:B:122:SER:HB3	1.79	0.65
1:B:249:VAL:HG21	1:B:257:ILE:CD1	2.26	0.64
1:A:324:GLU:OE1	1:A:339:ARG:NH2	2.29	0.64
1:B:194:MET:HG2	1:B:321:ILE:HG21	1.78	0.64
1:A:237:GLY:HA3	1:A:261:ILE:HG22	1.78	0.64
1:B:229:ALA:CB	1:B:232:GLU:HB2	2.27	0.64
1:B:325:THR:O	1:B:326:LEU:HD23	1.98	0.64
1:A:328:VAL:HG23	1:A:352:VAL:O	1.98	0.63
1:A:19:LYS:HD3	1:A:338:GLU:OE1	1.98	0.63
1:B:77:LEU:HD11	1:B:85:LEU:O	1.99	0.63
1:B:21:PRO:HG2	1:B:333:VAL:HG11	1.79	0.63
1:B:328:VAL:HG23	1:B:352:VAL:O	1.98	0.63
1:B:97:VAL:CG2	1:B:141:VAL:HG13	2.23	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ILE:HG22	1:A:21:PRO:HB2	1.80	0.62
1:A:189:GLY:HA3	3:A:4292:NAP:H4B	1.80	0.62
1:B:50:ILE:HD11	1:B:337:PHE:HA	1.81	0.62
1:B:31:PHE:HD1	1:B:140:ARG:HD3	1.65	0.62
1:A:1:MET:HG3	1:A:2:SER:H	1.63	0.62
1:B:300:ALA:HA	3:B:4293:NAP:C7N	2.29	0.62
1:B:31:PHE:HB2	1:B:140:ARG:NH1	2.14	0.62
1:A:113:TYR:CE2	1:B:293:ALA:HB2	2.35	0.61
1:B:41:GLU:OE1	1:B:73:LYS:HE2	2.00	0.61
1:B:163:CYS:O	1:B:164:GLY:C	2.44	0.61
1:B:161:LEU:HD13	1:B:314:VAL:HG21	1.82	0.61
1:B:97:VAL:O	1:B:97:VAL:HG22	2.00	0.61
1:B:296:ILE:HD13	1:B:297:SER:N	2.16	0.61
1:A:100:CYS:HB3	1:A:116:LYS:HG3	1.81	0.60
1:A:211:ARG:HH12	3:A:4292:NAP:P2B	2.24	0.60
1:A:324:GLU:CD	1:A:339:ARG:HH22	2.09	0.60
1:B:6:LYS:HE2	1:B:27:ASP:OD2	2.01	0.60
1:B:340:MET:C	1:B:342:LYS:H	2.09	0.60
1:A:99:SER:HB2	1:A:145:PHE:HZ	1.66	0.60
1:B:13:GLN:O	1:B:62:MET:HG2	2.01	0.60
1:B:304:ILE:HD13	1:B:305:LYS:N	2.12	0.60
1:A:14:SER:HA	1:A:62:MET:HE3	1.83	0.60
1:A:69:GLU:OE1	1:A:348:ARG:NH1	2.35	0.60
1:B:300:ALA:CA	3:B:4293:NAP:H72N	2.04	0.59
1:B:110:ASN:C	1:B:112:PRO:HD2	2.27	0.59
1:B:211:ARG:HA	1:B:230:THR:N	2.17	0.59
1:B:314:VAL:HA	1:B:319:ILE:CG1	2.32	0.59
1:A:287:LYS:HG2	1:B:283:MET:HG2	1.84	0.59
1:B:247:ILE:HD13	1:B:262:MET:HE2	1.85	0.59
1:A:111:GLU:HG2	1:A:112:PRO:HD3	1.85	0.59
1:B:160:PRO:C	1:B:194:MET:HE1	2.28	0.59
1:B:3:TYR:HB2	1:B:138:TYR:CZ	2.38	0.58
1:A:320:LYS:O	1:A:320:LYS:HD3	2.03	0.58
1:B:156:HIS:CE1	1:B:157:LEU:HD13	2.38	0.58
1:A:304:ILE:HD12	1:A:305:LYS:N	2.17	0.58
1:A:35:ASP:HA	1:A:142:HIS:HA	1.85	0.58
1:A:289:TYR:CG	1:B:280:GLN:HA	2.38	0.58
1:B:12:ILE:HD13	1:B:21:PRO:HG3	1.86	0.58
1:B:99:SER:O	1:B:101:LEU:HG	2.04	0.58
1:B:161:LEU:HD23	1:B:194:MET:CE	2.33	0.58
1:B:237:GLY:HA2	1:B:265:ALA:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:ILE:HG21	1:B:262:MET:CE	2.33	0.58
1:A:66:VAL:HG13	1:A:121:TYR:CZ	2.39	0.57
1:A:120:THR:O	1:A:131:SER:HB2	2.03	0.57
1:A:253:SER:HB3	3:A:4292:NAP:N6A	2.18	0.57
1:A:286:LEU:HD12	1:A:286:LEU:O	2.04	0.57
1:B:211:ARG:CB	1:B:230:THR:HB	2.34	0.57
1:A:258:ASP:OD1	1:A:261:ILE:HG13	2.04	0.57
1:B:246:LEU:HD12	1:B:247:ILE:H	1.69	0.57
1:A:19:LYS:NZ	1:A:341:GLU:OE1	2.35	0.57
1:A:170:SER:O	1:A:174:ARG:HG3	2.04	0.57
1:B:252:SER:N	3:B:4293:NAP:C4A	2.68	0.57
1:A:197:LEU:HD12	1:A:321:ILE:HG22	1.86	0.57
1:A:254:LEU:HD13	1:A:284:LEU:HD13	1.86	0.57
1:A:14:SER:CA	1:A:62:MET:HE3	2.35	0.57
1:A:167:THR:HG23	1:A:275:ILE:HD12	1.87	0.57
1:B:316:GLU:C	1:B:318:ASP:H	2.12	0.57
1:A:90:ARG:NH2	1:A:153:ILE:O	2.38	0.56
1:A:289:TYR:CZ	1:B:280:GLN:HG2	2.39	0.56
1:B:229:ALA:HB3	1:B:232:GLU:HB2	1.86	0.56
1:A:156:HIS:CE1	1:A:157:LEU:HD13	2.40	0.56
1:B:188:LEU:HD23	1:B:215:LYS:CB	2.34	0.56
1:A:256:ASP:HB2	3:A:4292:NAP:N1A	2.20	0.56
1:A:29:LYS:HE3	1:A:140:ARG:HG2	1.87	0.56
1:B:167:THR:HG22	1:B:275:ILE:CD1	2.36	0.56
1:B:111:GLU:OE1	1:B:303:SER:HB2	2.05	0.56
1:B:168:VAL:HG23	1:B:191:ILE:HG22	1.87	0.56
1:B:19:LYS:NZ	1:B:19:LYS:HB3	2.21	0.56
1:B:254:LEU:C	1:B:254:LEU:HD23	2.31	0.56
1:B:264:LYS:HB2	1:B:264:LYS:HZ2	1.66	0.56
1:B:211:ARG:HG2	1:B:211:ARG:HH11	1.70	0.55
1:A:66:VAL:CG1	1:A:67:GLY:H	2.17	0.55
1:A:119:THR:HG22	1:A:123:GLN:OE1	2.07	0.55
1:A:259:PHE:HB3	1:A:286:LEU:HD23	1.89	0.55
1:B:161:LEU:HD23	1:B:194:MET:HE2	1.89	0.55
1:A:41:GLU:HG2	1:A:72:GLY:HA2	1.88	0.55
1:A:316:GLU:C	1:A:318:ASP:H	2.14	0.55
1:B:98:PHE:HB3	1:B:118:VAL:HB	1.88	0.55
1:B:111:GLU:N	1:B:112:PRO:HD2	2.21	0.55
1:A:98:PHE:HB3	1:A:118:VAL:HB	1.88	0.55
1:A:137:ASN:N	1:A:137:ASN:HD22	2.04	0.55
1:A:137:ASN:HD22	1:A:138:TYR:H	1.53	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:TYR:O	1:A:173:VAL:HG23	2.07	0.55
1:B:111:GLU:HG2	1:B:112:PRO:CD	2.37	0.54
1:B:197:LEU:CD2	1:B:222:MET:HB3	2.37	0.54
1:B:21:PRO:O	1:B:330:GLU:HG2	2.07	0.54
1:B:236:TRP:H	1:B:236:TRP:CD1	2.23	0.54
1:A:82:ASN:HD22	1:A:82:ASN:N	2.05	0.54
1:B:166:LEU:HD13	1:B:307:LEU:HA	1.88	0.54
1:A:338:GLU:O	1:A:342:LYS:HD3	2.07	0.54
1:A:290:GLY:O	1:A:292:LYS:N	2.35	0.54
1:B:211:ARG:HA	1:B:230:THR:H	1.72	0.54
1:B:312:LYS:HE3	1:B:316:GLU:OE2	2.07	0.54
1:A:234:GLY:O	1:A:235:ASP:HB2	2.07	0.54
1:A:286:LEU:O	1:A:288:PRO:HD3	2.08	0.54
1:B:71:VAL:HG13	1:B:149:ILE:CD1	2.37	0.54
1:A:12:ILE:HD13	1:A:21:PRO:HG3	1.90	0.54
1:A:2:SER:C	1:A:4:PRO:HD2	2.32	0.54
1:A:21:PRO:O	1:A:330:GLU:HG2	2.08	0.54
1:B:30:PRO:HB2	1:B:32:TYR:CE1	2.43	0.54
1:B:194:MET:O	1:B:198:ILE:HG13	2.08	0.54
1:B:329:GLY:O	1:B:332:GLY:N	2.41	0.54
1:B:304:ILE:HD13	1:B:304:ILE:N	2.23	0.53
1:B:345:VAL:C	1:B:347:TYR:H	2.15	0.53
1:A:102:GLU:HA	1:A:107:LYS:HE3	1.88	0.53
1:A:137:ASN:HD22	1:A:138:TYR:N	2.06	0.53
1:A:194:MET:O	1:A:198:ILE:HG13	2.08	0.53
1:A:277:ILE:HB	3:A:4292:NAP:O2D	2.08	0.53
1:B:38:ILE:O	1:B:138:TYR:HA	2.09	0.53
1:B:346:ARG:O	1:B:347:TYR:HB2	2.09	0.53
1:B:6:LYS:HA	1:B:28:PRO:CD	2.31	0.53
1:A:74:VAL:HG11	1:A:77:LEU:HD21	1.91	0.53
1:A:111:GLU:N	1:A:112:PRO:CD	2.72	0.53
1:B:71:VAL:HG23	1:B:159:ALA:HB2	1.90	0.53
1:A:303:SER:OG	1:A:306:GLU:HG3	2.09	0.53
1:A:160:PRO:C	1:A:194:MET:HE1	2.34	0.52
1:B:329:GLY:O	1:B:331:ALA:N	2.43	0.52
1:B:225:ASP:O	1:B:226:HIS:HB2	2.09	0.52
1:A:1:MET:CG	1:A:2:SER:H	2.22	0.52
1:A:40:ILE:HD12	1:A:136:ALA:O	2.10	0.52
1:A:279:GLU:HG2	1:A:282:GLU:HG3	1.91	0.52
1:A:345:VAL:C	1:A:347:TYR:N	2.68	0.52
1:B:31:PHE:CE1	1:B:37:ASP:OD1	2.62	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:GLU:HG2	1:B:112:PRO:HD3	1.91	0.52
1:B:22:LYS:HE2	1:B:330:GLU:OE2	2.10	0.52
1:A:66:VAL:CG1	1:A:67:GLY:N	2.71	0.52
1:B:144:HIS:HD2	1:B:145:PHE:CE1	2.28	0.52
1:B:257:ILE:HD12	1:B:262:MET:SD	2.50	0.52
1:B:71:VAL:CG2	1:B:159:ALA:HB2	2.40	0.52
1:B:218:ASP:CG	1:B:346:ARG:NH2	2.68	0.52
1:B:12:ILE:HD11	1:B:337:PHE:HE1	1.75	0.52
1:B:226:HIS:ND1	1:B:227:TYR:N	2.57	0.51
1:B:229:ALA:HB1	1:B:232:GLU:HB2	1.91	0.51
1:B:169:TYR:O	1:B:173:VAL:HG22	2.10	0.51
1:B:179:PRO:HA	1:B:203:GLY:O	2.11	0.51
1:B:251:ALA:HA	3:B:4293:NAP:C1B	2.39	0.51
1:B:233:GLU:OE2	1:B:239:LYS:HE3	2.11	0.51
1:A:240:TYR:O	1:A:243:THR:CG2	2.58	0.51
1:B:99:SER:OG	1:B:106:CYS:SG	2.69	0.51
1:B:259:PHE:CD1	1:B:286:LEU:HD21	2.46	0.51
1:B:188:LEU:HB3	1:B:215:LYS:HD2	1.93	0.51
1:A:137:ASN:H	1:A:137:ASN:ND2	2.09	0.50
1:B:64:LEU:HD22	1:B:122:SER:HB3	1.93	0.50
1:B:187:GLY:O	1:B:192:GLY:HA3	2.11	0.50
1:B:252:SER:N	3:B:4293:NAP:N3A	2.59	0.50
1:B:182:LYS:HE2	1:B:205:GLU:OE2	2.11	0.50
1:A:92:GLY:HA3	1:A:149:ILE:HD11	1.93	0.50
1:B:31:PHE:HD1	1:B:140:ARG:CD	2.24	0.50
1:B:40:ILE:HD12	1:B:136:ALA:O	2.11	0.50
1:B:213:SER:O	1:B:216:ARG:CB	2.59	0.50
1:A:161:LEU:HA	1:A:194:MET:HE1	1.93	0.50
1:B:18:TRP:CZ3	1:B:337:PHE:HB3	2.46	0.50
1:B:328:VAL:HG11	1:B:354:TYR:CE1	2.46	0.50
1:B:328:VAL:HG23	1:B:352:VAL:C	2.36	0.50
1:B:74:VAL:CG1	1:B:87:VAL:HA	2.42	0.50
1:B:211:ARG:CB	1:B:231:LEU:HG	2.42	0.50
1:B:19:LYS:HB3	1:B:19:LYS:HZ2	1.76	0.49
1:A:132:GLN:O	1:A:133:GLY:O	2.29	0.49
1:B:264:LYS:HB2	1:B:264:LYS:HZ3	1.75	0.49
1:B:336:ALA:HA	1:B:349:PHE:HE1	1.77	0.49
1:A:13:GLN:C	1:A:62:MET:HG2	2.38	0.49
1:A:188:LEU:HG	1:A:188:LEU:O	2.11	0.49
1:A:195:GLY:O	1:A:199:SER:HB2	2.12	0.49
1:A:39:LYS:HG3	1:A:138:TYR:CE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:235:ASP:OD1	1:B:238:GLU:HB2	2.13	0.49
1:B:291:LEU:HB3	1:B:294:VAL:O	2.13	0.49
1:B:304:ILE:N	1:B:304:ILE:CD1	2.75	0.49
1:A:169:TYR:CE1	1:A:173:VAL:CG2	2.96	0.49
1:B:321:ILE:HG13	1:B:323:VAL:HG23	1.94	0.49
1:A:211:ARG:HG2	1:A:211:ARG:HH11	1.78	0.49
1:B:161:LEU:CD1	1:B:314:VAL:HG21	2.42	0.49
1:A:289:TYR:CD2	1:B:280:GLN:HA	2.48	0.49
1:A:3:TYR:HB2	1:A:138:TYR:CZ	2.49	0.48
1:B:20:ASN:N	1:B:21:PRO:HD3	2.28	0.48
1:B:40:ILE:HG23	1:B:70:ILE:HG23	1.95	0.48
1:B:211:ARG:HA	1:B:230:THR:HB	1.95	0.48
1:A:156:HIS:NE2	1:A:157:LEU:HD13	2.28	0.48
1:B:96:GLN:CD	1:B:145:PHE:HE2	2.21	0.48
1:B:216:ARG:HA	1:B:227:TYR:CE1	2.48	0.48
1:A:251:ALA:O	1:A:276:SER:HB3	2.14	0.48
1:A:295:SER:HA	1:B:297:SER:HB3	1.96	0.48
1:B:243:THR:O	1:B:267:LYS:HE2	2.13	0.48
1:A:69:GLU:OE2	1:A:163:CYS:HB3	2.14	0.48
1:A:97:VAL:CG2	1:A:141:VAL:HG13	2.40	0.48
1:B:31:PHE:CD1	1:B:140:ARG:HD3	2.47	0.48
1:B:87:VAL:HG12	1:B:88:GLY:N	2.29	0.48
1:B:262:MET:HB3	1:B:263:PRO:HD3	1.95	0.48
1:B:287:LYS:HG2	1:B:289:TYR:CZ	2.48	0.47
1:A:211:ARG:NH1	3:A:4292:NAP:P2B	2.87	0.47
1:A:161:LEU:HD23	1:A:194:MET:CE	2.45	0.47
1:B:106:CYS:C	1:B:108:ASN:H	2.22	0.47
1:B:191:ILE:H	1:B:191:ILE:CD1	2.19	0.47
1:A:62:MET:HG3	1:A:63:PRO:HA	1.97	0.47
1:A:67:GLY:HA3	1:A:135:TYR:CE1	2.49	0.47
1:B:2:SER:C	1:B:4:PRO:HD2	2.38	0.47
1:B:168:VAL:C	1:B:171:PRO:HD2	2.40	0.47
1:B:304:ILE:CD1	1:B:305:LYS:H	2.16	0.47
1:A:156:HIS:CD2	1:A:157:LEU:HD13	2.50	0.47
1:B:44:GLY:H	1:B:69:GLU:HB2	1.79	0.47
1:A:211:ARG:HA	1:A:230:THR:HB	1.96	0.47
1:B:253:SER:HB3	1:B:256:ASP:OD1	2.15	0.47
1:A:295:SER:HB2	1:B:295:SER:HB2	1.97	0.47
1:B:8:GLU:HA	1:B:25:LYS:HA	1.97	0.46
1:B:50:ILE:CD1	1:B:337:PHE:HA	2.44	0.46
1:B:137:ASN:HD22	1:B:137:ASN:N	2.12	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ILE:O	1:A:62:MET:HG3	2.15	0.46
1:B:3:TYR:CD1	1:B:4:PRO:HD3	2.50	0.46
1:B:336:ALA:HA	1:B:349:PHE:CE1	2.51	0.46
1:A:121:TYR:O	1:A:122:SER:HB2	2.16	0.46
1:A:273:VAL:HA	1:A:297:SER:O	2.16	0.46
1:A:285:SER:HA	1:B:284:LEU:O	2.15	0.46
1:A:3:TYR:HB2	1:A:138:TYR:CE1	2.51	0.46
1:A:82:ASN:N	1:A:82:ASN:ND2	2.63	0.46
1:B:111:GLU:OE1	1:B:304:ILE:HD12	2.14	0.46
1:B:96:GLN:CD	1:B:145:PHE:CE2	2.94	0.46
1:B:211:ARG:O	1:B:229:ALA:HB1	2.16	0.46
1:B:335:GLU:HG2	1:B:339:ARG:HH21	1.81	0.46
1:A:266:MET:HB2	1:A:294:VAL:HG21	1.96	0.46
1:B:211:ARG:HB2	1:B:231:LEU:CG	2.45	0.46
1:A:3:TYR:CD1	1:A:3:TYR:C	2.89	0.46
1:B:329:GLY:O	1:B:330:GLU:C	2.58	0.46
1:B:334:HIS:CE1	1:B:338:GLU:OE1	2.69	0.46
1:B:251:ALA:HB1	3:B:4293:NAP:H2A	1.93	0.46
1:B:175:ASN:ND2	1:B:273:VAL:HG21	2.30	0.46
1:B:259:PHE:HA	1:B:263:PRO:HD3	1.98	0.46
1:A:137:ASN:N	1:A:137:ASN:ND2	2.64	0.45
1:A:172:LEU:HD23	1:A:246:LEU:HD21	1.98	0.45
1:B:22:LYS:HA	1:B:330:GLU:OE2	2.16	0.45
1:B:51:HIS:HD1	1:B:51:HIS:H	1.64	0.45
1:B:65:VAL:HG23	1:B:132:GLN:OE1	2.16	0.45
1:B:71:VAL:HG13	1:B:149:ILE:HD12	1.98	0.45
1:A:20:ASN:N	1:A:21:PRO:HD3	2.32	0.45
1:B:339:ARG:HD2	1:B:344:ASP:OD1	2.17	0.45
1:A:32:TYR:HB2	1:A:35:ASP:OD2	2.17	0.45
1:A:136:ALA:O	1:A:354:TYR:CE1	2.70	0.45
1:B:279:GLU:C	1:B:281:HIS:N	2.74	0.45
1:B:314:VAL:HA	1:B:319:ILE:HG12	1.97	0.45
1:A:188:LEU:HB3	1:A:215:LYS:HD3	1.98	0.45
1:B:211:ARG:NH1	3:B:4293:NAP:O1X	2.39	0.45
1:A:275:ILE:HD13	1:A:275:ILE:HA	1.72	0.45
1:B:154:PRO:HG2	1:B:157:LEU:HB2	1.99	0.45
1:B:159:ALA:HB3	1:B:160:PRO:HD3	1.98	0.45
1:A:91:VAL:HB	1:A:146:VAL:HG12	1.99	0.45
1:B:257:ILE:O	1:B:259:PHE:N	2.50	0.45
1:B:47:GLY:O	1:B:50:ILE:HB	2.17	0.45
1:B:137:ASN:N	1:B:137:ASN:ND2	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:VAL:HG11	1:B:155:SER:HB3	1.99	0.44
1:A:96:GLN:OE1	1:A:112:PRO:HA	2.17	0.44
1:B:42:ALA:HB3	1:B:71:VAL:HB	1.99	0.44
1:B:82:ASN:N	1:B:82:ASN:HD22	2.15	0.44
1:B:346:ARG:O	1:B:347:TYR:CB	2.65	0.44
1:B:197:LEU:HD23	1:B:222:MET:HB3	1.99	0.44
1:B:198:ILE:HD13	1:B:310:LEU:HD11	2.00	0.44
1:A:70:ILE:HD11	1:A:134:GLY:HA2	1.98	0.44
1:A:123:GLN:O	1:A:130:VAL:HA	2.17	0.44
1:A:277:ILE:HD11	1:A:298:TYR:OH	2.18	0.44
1:B:304:ILE:CD1	1:B:304:ILE:H	2.30	0.44
1:A:290:GLY:C	1:A:292:LYS:N	2.68	0.44
1:B:340:MET:O	1:B:342:LYS:N	2.48	0.44
1:A:46:CYS:O	1:A:49:ASP:HB2	2.18	0.44
1:B:97:VAL:HG23	1:B:141:VAL:CG1	2.28	0.44
1:B:212:SER:HA	1:B:232:GLU:OE2	2.18	0.44
1:B:228:ILE:HG12	1:B:240:TYR:CZ	2.52	0.44
1:A:289:TYR:CE2	1:B:280:GLN:HG2	2.53	0.44
1:B:243:THR:HG23	1:B:244:PHE:CD2	2.52	0.44
1:A:328:VAL:HG21	1:A:354:TYR:CD2	2.53	0.44
1:B:163:CYS:SG	1:B:164:GLY:N	2.91	0.44
1:B:213:SER:N	1:B:232:GLU:OE2	2.51	0.44
1:A:8:GLU:O	1:A:136:ALA:HB1	2.18	0.43
1:A:161:LEU:HA	1:A:194:MET:CE	2.48	0.43
1:A:307:LEU:O	1:A:307:LEU:HD12	2.17	0.43
1:B:106:CYS:C	1:B:108:ASN:N	2.76	0.43
1:A:240:TYR:O	1:A:241:PHE:C	2.61	0.43
1:B:93:VAL:O	1:B:162:LEU:HD13	2.17	0.43
1:B:184:GLY:O	1:B:247:ILE:HA	2.18	0.43
1:B:316:GLU:C	1:B:318:ASP:N	2.74	0.43
1:B:48:SER:HB2	1:B:57:TRP:CH2	2.53	0.43
1:B:266:MET:HE3	1:B:270:GLY:HA3	2.00	0.43
1:A:29:LYS:HA	1:A:129:TYR:CE1	2.53	0.43
1:B:246:LEU:HD12	1:B:247:ILE:N	2.33	0.43
1:A:86:LYS:HD3	1:A:86:LYS:HA	1.82	0.43
1:A:276:SER:HA	3:A:4292:NAP:H1D	2.00	0.43
1:B:82:ASN:HD22	1:B:82:ASN:C	2.26	0.43
1:B:82:ASN:HD22	1:B:82:ASN:H	1.66	0.43
1:B:7:PHE:HB3	1:B:137:ASN:HD21	1.83	0.43
1:A:191:ILE:HG13	3:A:4292:NAP:O2N	2.19	0.43
1:B:163:CYS:SG	3:B:4293:NAP:H5N	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:LYS:HA	1:A:129:TYR:CD1	2.53	0.43
1:B:39:LYS:HA	1:B:137:ASN:O	2.18	0.43
1:B:70:ILE:O	1:B:162:LEU:CD1	2.67	0.43
1:B:76:LYS:HB2	1:B:76:LYS:NZ	2.34	0.43
1:A:40:ILE:HG22	1:A:41:GLU:N	2.33	0.42
1:B:6:LYS:HB2	1:B:26:TYR:O	2.19	0.42
1:B:157:LEU:HD12	1:B:323:VAL:HG21	2.01	0.42
1:B:167:THR:HG22	1:B:275:ILE:HD13	1.99	0.42
1:B:188:LEU:HB3	1:B:215:LYS:CD	2.49	0.42
1:B:211:ARG:HG2	1:B:211:ARG:NH1	2.33	0.42
1:A:214:ARG:HD3	1:A:214:ARG:O	2.19	0.42
1:B:345:VAL:HG22	1:B:349:PHE:CZ	2.54	0.42
1:A:248:VAL:HA	1:A:273:VAL:O	2.20	0.42
1:B:80:LYS:O	1:B:81:SER:C	2.61	0.42
1:B:327:PRO:O	1:B:329:GLY:N	2.52	0.42
1:B:328:VAL:HG22	1:B:351:LEU:C	2.44	0.42
1:B:12:ILE:HG22	1:B:13:GLN:N	2.35	0.42
1:A:92:GLY:CA	1:A:149:ILE:HD11	2.49	0.42
1:A:243:THR:HG23	1:A:244:PHE:CD2	2.55	0.42
1:A:253:SER:HB3	3:A:4292:NAP:C6A	2.49	0.42
1:A:262:MET:HB3	1:A:263:PRO:HD3	2.01	0.42
1:B:127:ASP:C	1:B:127:ASP:OD1	2.62	0.42
1:B:211:ARG:HG2	1:B:211:ARG:H	1.63	0.42
1:B:211:ARG:CA	1:B:230:THR:HB	2.50	0.42
1:B:275:ILE:O	1:B:275:ILE:HG23	2.20	0.42
1:A:283:MET:HE3	1:B:286:LEU:C	2.45	0.42
1:B:211:ARG:HH12	3:B:4293:NAP:P2B	2.43	0.42
1:A:286:LEU:HD12	1:A:286:LEU:C	2.45	0.42
1:A:38:ILE:HB	1:A:139:VAL:HG12	2.02	0.42
1:A:304:ILE:CD1	1:A:305:LYS:N	2.83	0.42
1:B:280:GLN:C	1:B:281:HIS:ND1	2.78	0.42
1:B:326:LEU:HB2	1:B:351:LEU:HD23	2.02	0.42
1:A:292:LYS:O	1:A:293:ALA:HB3	2.20	0.41
1:B:29:LYS:HG3	1:B:30:PRO:O	2.20	0.41
1:B:252:SER:HB3	3:B:4293:NAP:N1A	2.30	0.41
1:B:264:LYS:NZ	1:B:264:LYS:CB	2.78	0.41
1:B:345:VAL:HG13	1:B:349:PHE:CE2	2.55	0.41
1:A:45:VAL:HB	1:A:349:PHE:HB2	2.02	0.41
1:A:327:PRO:O	1:A:332:GLY:HA3	2.20	0.41
1:B:31:PHE:HA	1:B:140:ARG:HD3	2.02	0.41
1:B:167:THR:CG2	1:B:275:ILE:HG12	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:LEU:HD21	1:B:222:MET:HB3	2.02	0.41
1:B:348:ARG:HA	1:B:348:ARG:HD3	1.69	0.41
1:A:278:PRO:O	1:A:279:GLU:C	2.63	0.41
1:A:70:ILE:O	1:A:162:LEU:HD13	2.19	0.41
1:A:137:ASN:HD22	1:A:137:ASN:H	1.62	0.41
1:B:22:LYS:HE2	1:B:22:LYS:HA	2.01	0.41
1:B:45:VAL:HB	1:B:349:PHE:HB2	2.02	0.41
1:B:133:GLY:HA3	1:B:139:VAL:CG2	2.49	0.41
1:A:196:THR:O	1:A:199:SER:HB3	2.20	0.41
1:A:292:LYS:HE2	1:B:113:TYR:HD2	1.84	0.41
1:A:49:ASP:HA	1:A:66:VAL:HG11	2.02	0.41
1:A:261:ILE:O	1:A:262:MET:C	2.63	0.41
1:A:3:TYR:O	1:A:3:TYR:CG	2.70	0.41
1:A:258:ASP:OD1	1:A:260:ASN:HB2	2.20	0.41
1:A:300:ALA:C	1:A:301:LEU:O	2.64	0.41
1:B:301:LEU:HB2	1:B:302:GLY:H	1.51	0.41
1:A:64:LEU:HD21	1:A:122:SER:OG	2.21	0.41
1:A:183:VAL:HG12	1:A:184:GLY:N	2.36	0.41
1:B:137:ASN:ND2	1:B:138:TYR:N	2.66	0.41
1:B:240:TYR:HB3	1:B:243:THR:HG21	2.02	0.41
1:B:16:GLU:C	1:B:18:TRP:H	2.27	0.41
1:B:136:ALA:O	1:B:354:TYR:CE1	2.74	0.41
1:B:274:SER:O	1:B:298:TYR:HA	2.21	0.41
1:B:301:LEU:CD2	3:B:4293:NAP:O7N	2.44	0.41
1:B:1:MET:C	1:B:25:LYS:HZ3	2.29	0.41
1:A:133:GLY:HA3	1:A:139:VAL:HG23	2.03	0.40
1:A:163:CYS:O	1:A:164:GLY:C	2.64	0.40
1:A:218:ASP:O	1:A:222:MET:HG3	2.21	0.40
1:A:236:TRP:CD1	1:A:236:TRP:H	2.38	0.40
1:A:316:GLU:C	1:A:318:ASP:N	2.80	0.40
1:B:82:ASN:C	1:B:82:ASN:ND2	2.79	0.40
1:B:117:PHE:CD2	1:B:117:PHE:C	2.99	0.40
1:B:171:PRO:HG3	1:B:299:SER:CB	2.51	0.40
1:B:218:ASP:CG	1:B:346:ARG:HH22	2.28	0.40
1:B:218:ASP:OD1	1:B:346:ARG:NH1	2.53	0.40
1:B:323:VAL:HG12	1:B:324:GLU:N	2.35	0.40
1:B:129:TYR:CD2	1:B:129:TYR:C	2.99	0.40
1:B:213:SER:O	1:B:214:ARG:C	2.64	0.40
1:B:258:ASP:OD1	1:B:261:ILE:HG13	2.22	0.40
1:A:2:SER:O	1:A:4:PRO:N	2.54	0.40
1:A:121:TYR:HA	1:A:131:SER:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:ALA:CB	1:B:352:VAL:HG12	2.51	0.40
1:A:137:ASN:ND2	1:A:138:TYR:N	2.69	0.40
1:A:249:VAL:HB	1:A:274:SER:HA	2.03	0.40
1:A:359:SER:OG	1:A:360:ASP:N	2.55	0.40
1:B:292:LYS:O	1:B:293:ALA:HB3	2.21	0.40
1:A:4:PRO:HG2	1:A:5:GLU:OE1	2.22	0.40
1:A:110:ASN:C	1:A:112:PRO:CD	2.95	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:GLU:OE2	1:B:338:GLU:OE2[2_756]	1.91	0.29

5.3 Torsion angles

5.3.1 Protein backbone

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	358/360 (99%)	302 (84%)	45 (13%)	11 (3%)	3	19
1	B	358/360 (99%)	290 (81%)	51 (14%)	17 (5%)	2	11
All	All	716/720 (99%)	592 (83%)	96 (13%)	28 (4%)	2	14

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	TYR
1	A	83	SER
1	A	235	ASP
1	B	251	ALA
1	B	283	MET
1	B	330	GLU

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Mol	Chain	Res	Type
1	A	82	ASN
1	A	133	GLY
1	B	155	SER
1	B	230	THR
1	B	318	ASP
1	B	328	VAL
1	B	347	TYR
1	A	241	PHE
1	A	289	TYR
1	A	291	LEU
1	A	347	TYR
1	B	3	TYR
1	B	341	GLU
1	B	357	GLU
1	A	4	PRO
1	A	301	LEU
1	B	11	ALA
1	B	214	ARG
1	B	18	TRP
1	B	101	LEU
1	B	164	GLY
1	B	66	VAL

5.3.2 Protein sidechains ⓘ

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	303/303 (100%)	287 (95%)	16 (5%)	20	54
1	B	303/303 (100%)	283 (93%)	20 (7%)	15	46
All	All	606/606 (100%)	570 (94%)	36 (6%)	18	50

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	SER

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Mol	Chain	Res	Type
1	A	59	ASN
1	A	62	MET
1	A	75	VAL
1	A	82	ASN
1	A	90	ARG
1	A	97	VAL
1	A	137	ASN
1	A	168	VAL
1	A	213	SER
1	A	275	ILE
1	A	301	LEU
1	A	304	ILE
1	A	315	SER
1	A	328	VAL
1	A	348	ARG
1	B	48	SER
1	B	76	LYS
1	B	82	ASN
1	B	90	ARG
1	B	111	GLU
1	B	126	GLU
1	B	137	ASN
1	B	157	LEU
1	B	173	VAL
1	B	191	ILE
1	B	261	ILE
1	B	264	LYS
1	B	267	LYS
1	B	275	ILE
1	B	284	LEU
1	B	296	ILE
1	B	301	LEU
1	B	304	ILE
1	B	319	ILE
1	B	348	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	82	ASN
1	A	108	ASN

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Mol	Chain	Res	Type
1	A	137	ASN
1	A	226	HIS
1	A	308	ASN
1	B	13	GLN
1	B	20	ASN
1	B	82	ASN
1	B	123	GLN
1	B	137	ASN
1	B	144	HIS
1	B	308	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
3	NAP	B	4293	-	50,52,52	2.02	12 (24%)	71,80,80	1.50	15 (21%)
3	NAP	A	4292	-	50,52,52	2.02	12 (24%)	71,80,80	1.50	15 (21%)

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	NAP	B	4293	-	-	5/35/67/67	0/5/5/5
3	NAP	A	4292	-	-	5/35/67/67	0/5/5/5

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	4292	NAP	C2N-N1N	6.63	1.42	1.35
3	B	4293	NAP	C2N-N1N	6.57	1.42	1.35
3	B	4293	NAP	C4N-C3N	5.10	1.47	1.39
3	A	4292	NAP	C4N-C3N	5.08	1.47	1.39
3	B	4293	NAP	O4D-C1D	4.74	1.47	1.40
3	A	4292	NAP	O4D-C1D	4.72	1.47	1.40
3	B	4293	NAP	C6N-N1N	3.89	1.44	1.35
3	A	4292	NAP	C6N-N1N	3.88	1.44	1.35
3	B	4293	NAP	C2A-N1A	3.80	1.40	1.33
3	A	4292	NAP	C2A-N1A	3.78	1.40	1.33
3	A	4292	NAP	C2A-N3A	3.39	1.39	1.33
3	B	4293	NAP	C2A-N3A	3.35	1.39	1.33
3	B	4293	NAP	C4A-N3A	2.85	1.39	1.34
3	A	4292	NAP	C4A-N3A	2.83	1.39	1.34
3	B	4293	NAP	PN-O3	-2.66	1.56	1.59
3	A	4292	NAP	PN-O3	-2.62	1.56	1.59
3	A	4292	NAP	O4B-C4B	2.57	1.50	1.45
3	B	4293	NAP	O4B-C4B	2.55	1.50	1.45
3	A	4292	NAP	PA-O1A	-2.44	1.42	1.50
3	B	4293	NAP	PA-O1A	-2.44	1.42	1.50
3	B	4293	NAP	C5N-C4N	2.10	1.42	1.38
3	A	4292	NAP	C5N-C4N	2.09	1.42	1.38
3	A	4292	NAP	C3N-C7N	2.08	1.53	1.50
3	B	4293	NAP	C3N-C7N	2.06	1.53	1.50

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	4292	NAP	O4B-C1B-N9A	4.94	117.58	108.09
3	B	4293	NAP	O4B-C1B-N9A	4.94	117.58	108.09
3	A	4292	NAP	N3A-C2A-N1A	-3.48	123.31	128.58
3	B	4293	NAP	N3A-C2A-N1A	-3.46	123.34	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	4293	NAP	O2A-PA-O1A	3.05	126.64	112.44
3	A	4292	NAP	O2A-PA-O1A	3.05	126.63	112.44
3	A	4292	NAP	C3N-C7N-N7N	2.97	121.40	117.74
3	B	4293	NAP	C3N-C7N-N7N	2.96	121.39	117.74
3	B	4293	NAP	C1B-N9A-C8A	2.77	133.25	127.09
3	A	4292	NAP	C1B-N9A-C8A	2.76	133.22	127.09
3	A	4292	NAP	O7N-C7N-C3N	-2.60	116.42	119.60
3	B	4293	NAP	O7N-C7N-C3N	-2.59	116.43	119.60
3	A	4292	NAP	C2N-C3N-C4N	2.54	121.21	118.26
3	B	4293	NAP	C2N-C3N-C4N	2.51	121.18	118.26
3	B	4293	NAP	O3X-P2B-O2X	2.51	117.22	107.80
3	A	4292	NAP	O3X-P2B-O2X	2.51	117.21	107.80
3	A	4292	NAP	O5B-C5B-C4B	-2.37	100.91	108.99
3	B	4293	NAP	O5B-C5B-C4B	-2.37	100.92	108.99
3	A	4292	NAP	C2B-C1B-N9A	-2.31	109.95	113.75
3	B	4293	NAP	C2B-C1B-N9A	-2.29	109.99	113.75
3	B	4293	NAP	C6N-N1N-C1D	-2.22	115.36	119.73
3	B	4293	NAP	O4B-C1B-C2B	-2.22	102.76	106.59
3	A	4292	NAP	C6N-N1N-C1D	-2.22	115.37	119.73
3	A	4292	NAP	O4B-C1B-C2B	-2.21	102.78	106.59
3	B	4293	NAP	C4A-N9A-C1B	-2.13	121.65	126.63
3	A	4292	NAP	C4A-N9A-C1B	-2.12	121.68	126.63
3	B	4293	NAP	C5A-C4A-N3A	-2.06	123.88	126.72
3	A	4292	NAP	C5A-C4A-N3A	-2.05	123.90	126.72
3	A	4292	NAP	C5N-C4N-C3N	-2.03	118.37	120.36
3	B	4293	NAP	C5N-C4N-C3N	-2.02	118.38	120.36

There are no chirality outliers.

All (10) torsion outliers are listed below:

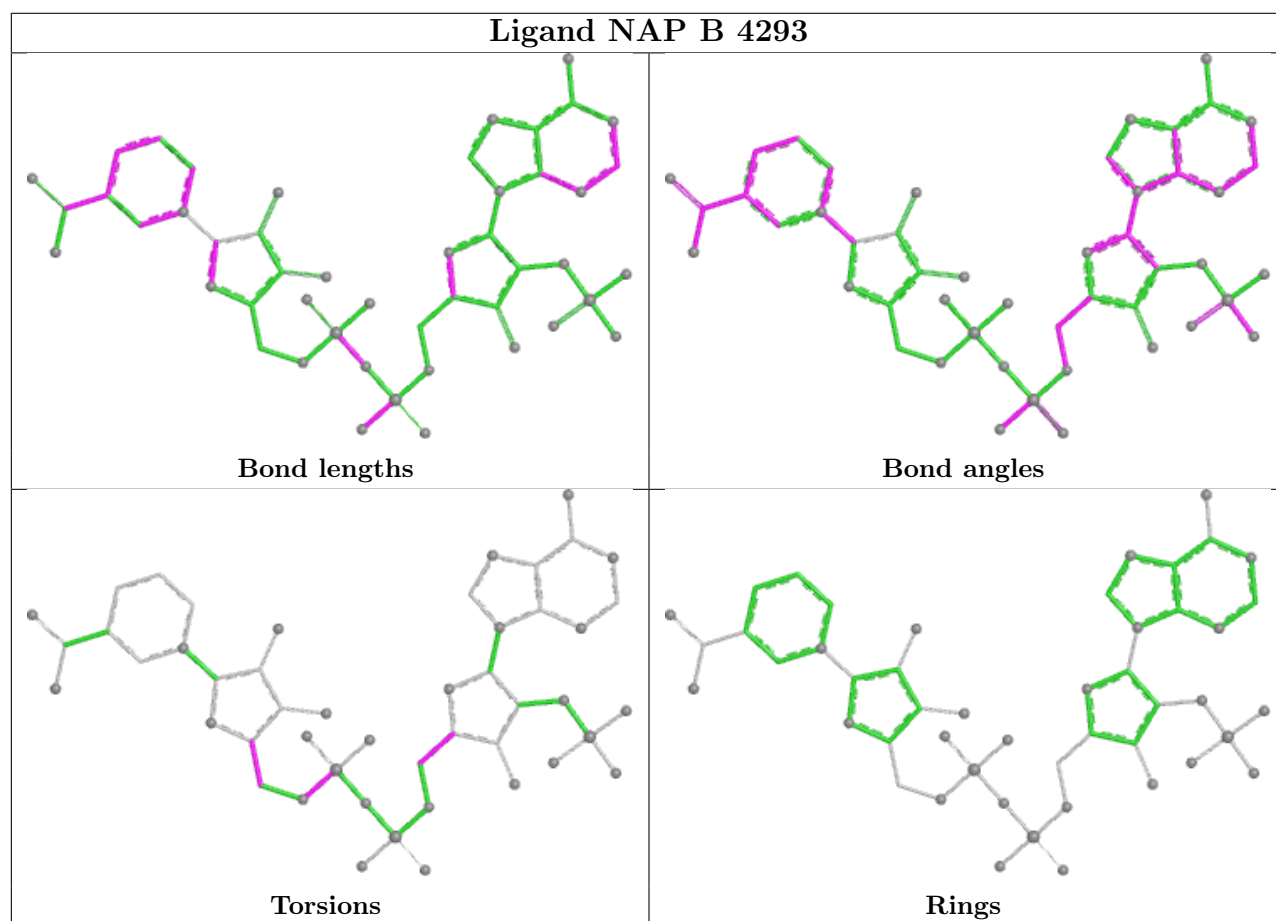
Mol	Chain	Res	Type	Atoms
3	A	4292	NAP	C5D-O5D-PN-O2N
3	B	4293	NAP	C5D-O5D-PN-O2N
3	A	4292	NAP	C3D-C4D-C5D-O5D
3	B	4293	NAP	C3D-C4D-C5D-O5D
3	A	4292	NAP	O4D-C4D-C5D-O5D
3	B	4293	NAP	O4D-C4D-C5D-O5D
3	A	4292	NAP	C5D-O5D-PN-O3
3	B	4293	NAP	C5D-O5D-PN-O3
3	A	4292	NAP	O4B-C4B-C5B-O5B
3	B	4293	NAP	O4B-C4B-C5B-O5B

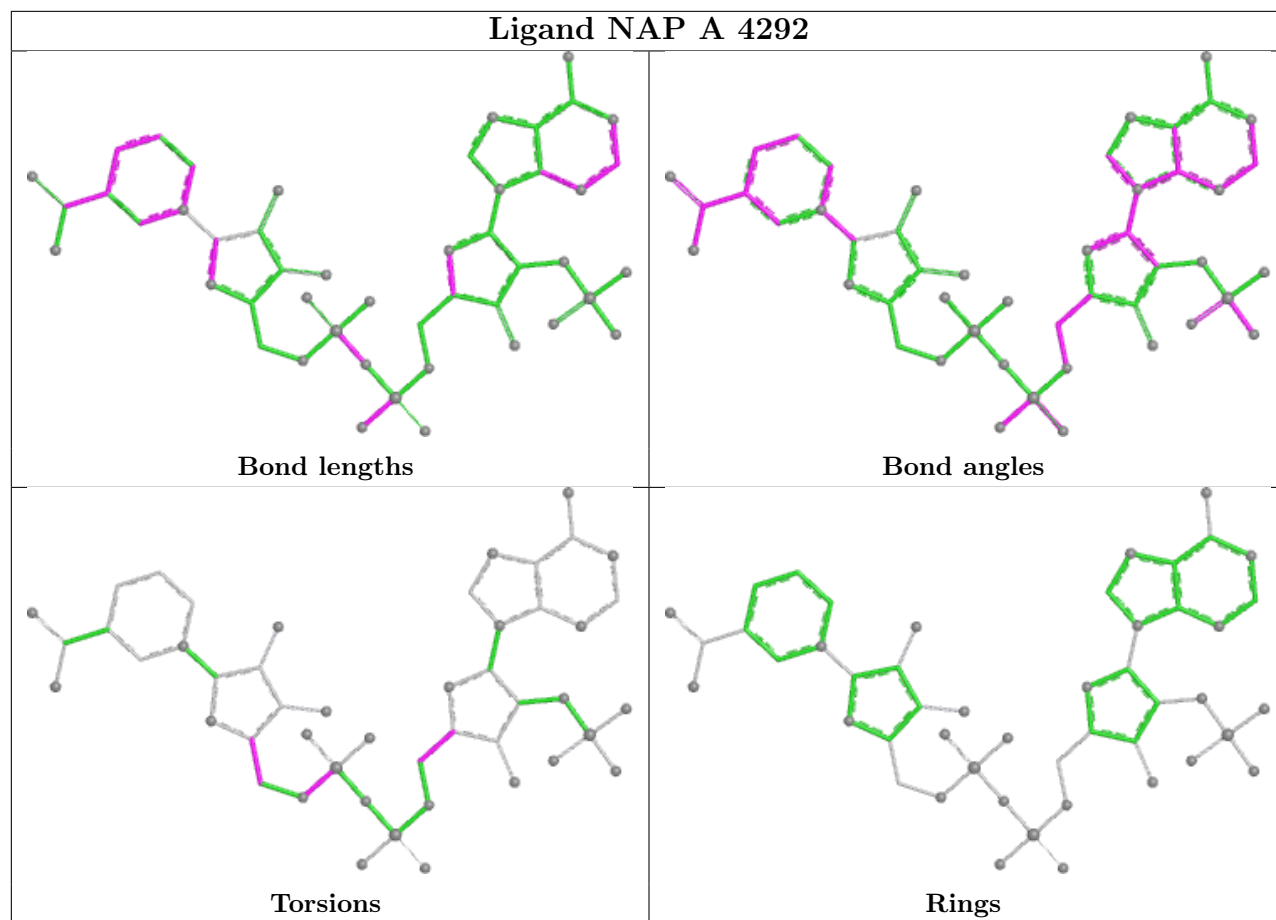
There are no ring outliers.

2 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	4293	NAP	27	0
3	A	4292	NAP	11	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	360/360 (100%)	-0.29	3 (0%) 82 64	23, 45, 71, 93	0
1	B	360/360 (100%)	0.10	3 (0%) 82 64	36, 67, 89, 100	0
All	All	720/720 (100%)	-0.09	6 (0%) 82 64	23, 54, 82, 100	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	2.7
1	B	255	THR	2.6
1	B	3	TYR	2.4
1	B	187	GLY	2.3
1	A	3	TYR	2.2
1	A	359	SER	2.2

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](#)

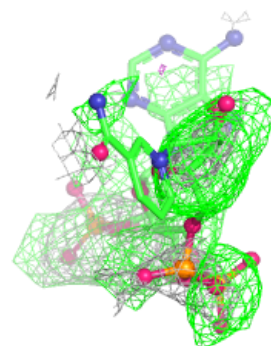
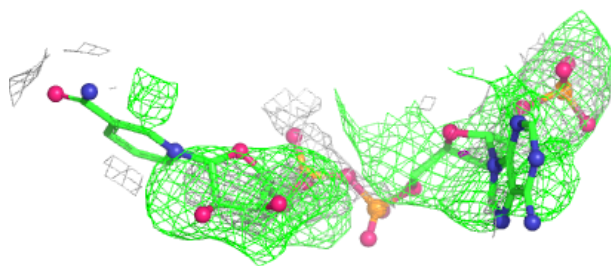
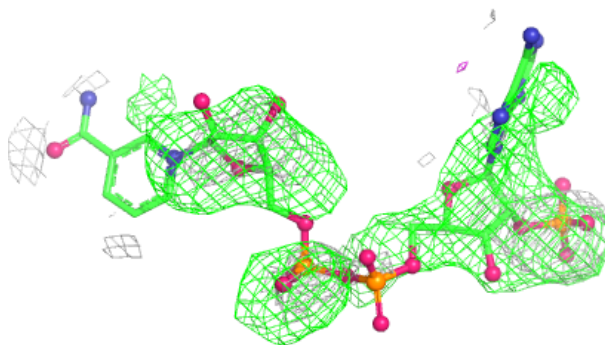
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAP	B	4293	48/48	0.83	0.51	50,50,50,50	48
3	NAP	A	4292	48/48	0.90	0.19	30,58,74,75	9
2	ZN	A	1502	1/1	0.99	0.06	32,32,32,32	0
2	ZN	B	1502	1/1	0.99	0.02	65,65,65,65	0
2	ZN	B	1501	1/1	1.00	0.03	52,52,52,52	0
2	ZN	A	1501	1/1	1.00	0.02	40,40,40,40	0

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

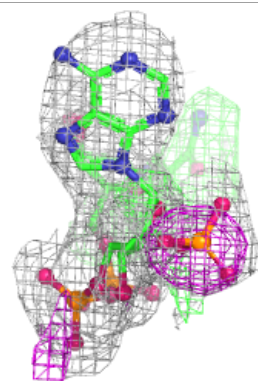
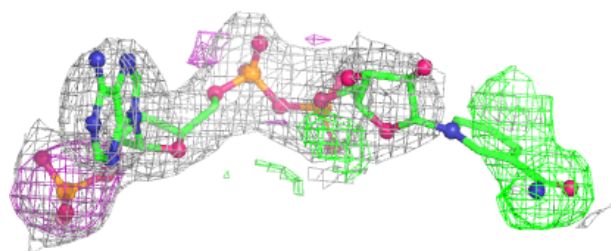
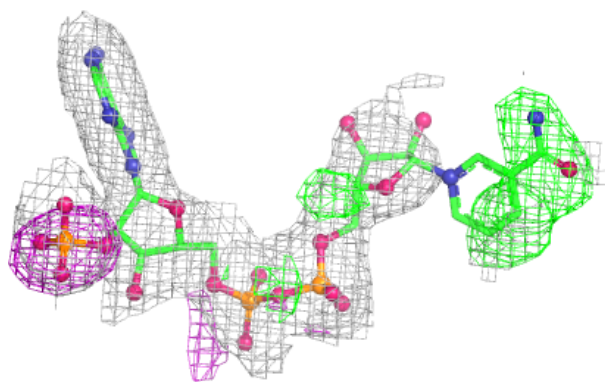
Electron density around NAP B 4293:

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



Electron density around NAP A 4292:

$2mF_o-DF_c$ (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.