



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 03:54 AM UTC

PDB ID : 1NXG / pdb_00001nxg
Title : The F383A variant of type II Citrate Synthase complexed with NADH
Authors : Maurus, R.; Nguyen, N.T.; Stokell, D.J.; Ayed, A.; Hultin, P.G.; Duckworth, H.W.; Brayer, G.D.
Deposited on : 2003-02-10
Resolution : 2.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

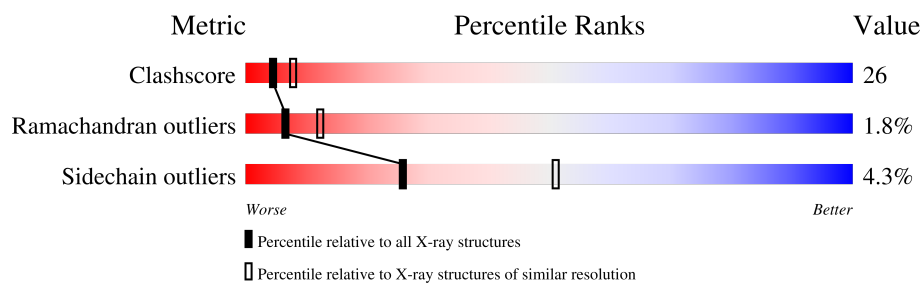
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	427	
1	B	427	

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
2	SO4	B	2001	-	X	-	-
2	SO4	B	2003	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7409 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 Citrate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	426	Total	C	N	O	S	0	0	0
			3359	2129	579	626	25			
1	B	426	Total	C	N	O	S	0	0	0
			3360	2129	579	627	25			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	383	ALA	PHE	engineered mutation	UNP P0ABH7
B	1383	ALA	PHE	engineered mutation	UNP P0ABH7

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



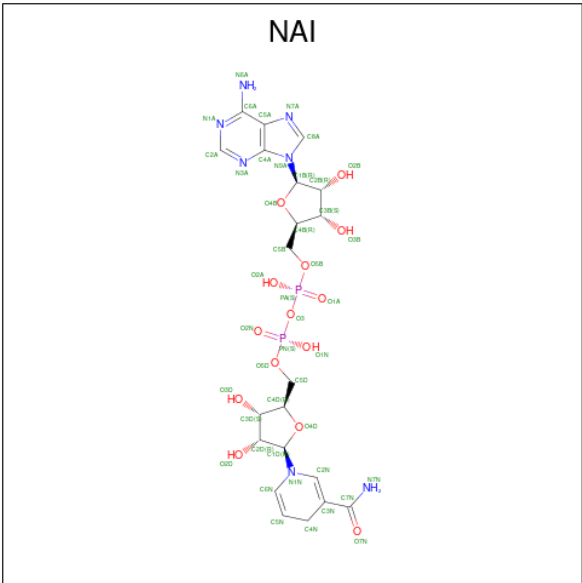
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	293	Total	O	0	0
			293	293		
4	B	289	Total	O	0	0
			289	289		

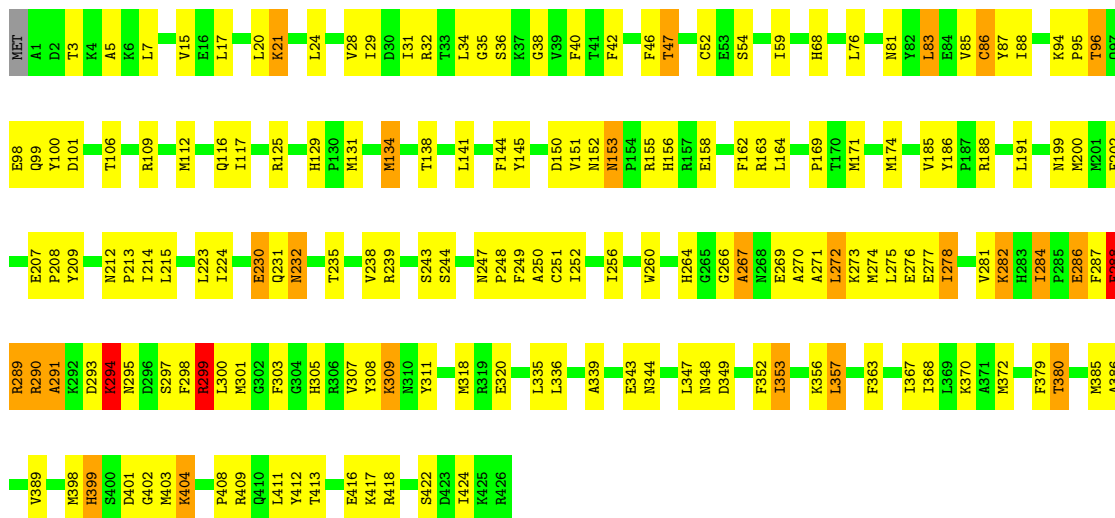
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

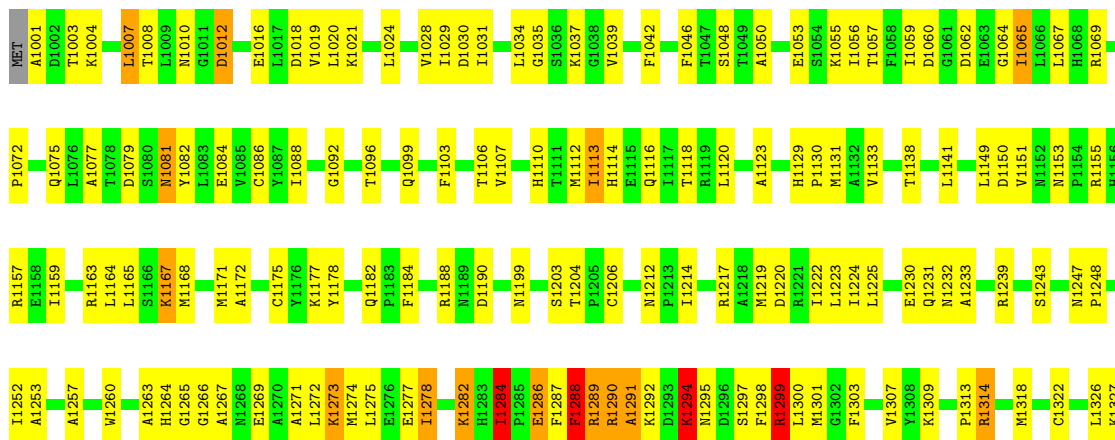
• Molecule 1: Citrate synthase

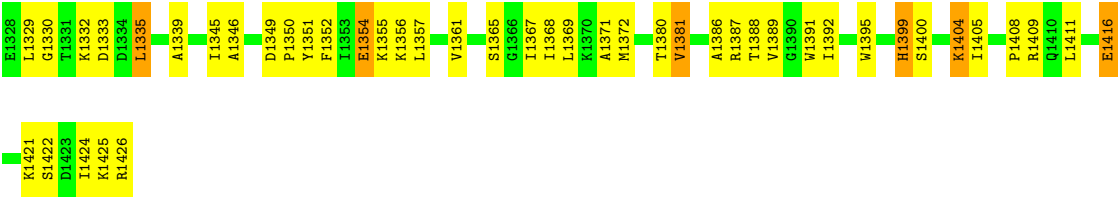
Chain A:  59% 34% 6% •



• Molecule 1: Citrate synthase

Chain B:  54% 40% 5% •





4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	164.89Å 164.89Å 158.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.187 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7409	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAI, SO4

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	0.49	0/3436	1.07	17/4645 (0.4%)
1	B	0.47	0/3437	1.08	24/4645 (0.5%)
All	All	0.48	0/6873	1.07	41/9290 (0.4%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1295	ASN	N-CA-C	-11.30	98.99	113.12
1	B	1291	ALA	N-CA-C	-9.53	102.74	114.75
1	A	291	ALA	N-CA-C	-9.39	102.92	114.75
1	A	295	ASN	N-CA-C	-9.09	100.99	112.90
1	B	1288	PHE	N-CA-C	-8.01	98.67	110.52
1	B	1048	SER	N-CA-C	-7.71	100.90	112.54
1	A	96	THR	N-CA-C	-7.08	101.00	110.55
1	B	1284	ILE	CA-C-N	7.03	128.63	119.84
1	B	1284	ILE	C-N-CA	7.03	128.63	119.84
1	A	288	PHE	N-CA-C	-6.93	99.02	110.17
1	B	1271	ALA	N-CA-C	-6.92	104.14	112.59
1	B	1167	LYS	N-CA-C	6.88	121.50	113.18
1	B	1222	ILE	N-CA-C	-6.82	103.83	110.72
1	B	1165	LEU	N-CA-C	-6.38	104.19	112.23
1	B	1212	ASN	CA-C-N	6.37	126.86	119.47
1	B	1212	ASN	C-N-CA	6.37	126.86	119.47
1	A	202	PHE	N-CA-C	6.36	121.57	113.55
1	A	47	THR	N-CA-C	-6.27	104.33	112.23
1	B	1307	VAL	N-CA-C	6.16	116.27	110.30
1	A	380	THR	N-CA-C	6.02	118.63	111.71
1	B	1289	ARG	N-CA-C	6.01	116.90	108.00
1	A	363	PHE	N-CA-C	-6.00	104.42	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	271	ALA	N-CA-C	-5.91	105.39	112.59
1	B	1416	GLU	N-CA-C	-5.84	102.99	110.53
1	B	1400	SER	N-CA-C	-5.82	105.44	113.18
1	A	299	ARG	N-CA-C	-5.78	105.99	113.16
1	A	309	LYS	N-CA-C	5.77	118.68	111.24
1	B	1065	ILE	N-CA-C	5.76	116.40	107.99
1	B	1299	ARG	N-CA-C	-5.66	106.14	113.16
1	A	416	GLU	N-CA-C	-5.65	102.10	110.46
1	B	1411	LEU	N-CA-C	-5.45	100.22	108.67
1	A	300	LEU	N-CA-C	-5.32	105.64	111.82
1	A	353	ILE	N-CA-C	-5.29	105.38	110.72
1	B	1056	ILE	N-CA-C	5.27	115.37	110.42
1	B	1314	ARG	N-CA-C	-5.25	107.00	113.41
1	B	1381	VAL	N-CA-C	-5.16	105.30	110.30
1	A	86	CYS	N-CA-C	-5.05	105.67	111.07
1	A	152	ASN	N-CA-C	5.04	119.48	113.23
1	A	289	ARG	N-CA-C	5.03	115.45	108.00
1	B	1399	HIS	N-CA-C	5.01	118.97	113.21
1	B	1079	ASP	N-CA-C	5.01	118.54	112.93

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	3359	0	3309	184	0
1	B	3360	0	3306	191	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	44	0	27	6	0
3	B	44	0	27	1	0
4	A	293	0	0	19	0
4	B	289	0	0	17	0
All	All	7409	0	6669	353	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 26.

All (353) 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:404:LYS:H	1:A:404:LYS:HD3	1.16	1.06
1:B:1081:ASN:ND2	1:B:1084:GLU:H	1.67	0.91
1:B:1272:LEU:HA	1:B:1275:LEU:HG	1.55	0.88
1:B:1287:PHE:HA	1:B:1291:ALA:HB3	1.58	0.86
1:A:275:LEU:HD22	1:A:289:ARG:HE	1.39	0.85
1:A:96:THR:H	1:A:99:GLN:HE21	1.24	0.85
1:A:298:PHE:HD2	1:A:301:MET:HE2	1.42	0.84
1:A:287:PHE:HA	1:A:291:ALA:HB3	1.61	0.82
1:A:282:LYS:HA	1:A:288:PHE:HD2	1.45	0.81
1:B:1129:HIS:CD2	1:B:1131:MET:H	1.99	0.81
1:A:309:LYS:NZ	1:A:356:LYS:HE3	1.96	0.80
1:B:1275:LEU:HD22	1:B:1289:ARG:HE	1.46	0.80
3:A:3000:NAI:O3	3:A:3000:NAI:H3B	1.83	0.79
1:A:404:LYS:HD3	1:A:404:LYS:N	1.96	0.79
1:B:1335:LEU:H	1:B:1335:LEU:HD23	1.49	0.77
1:A:305:HIS:HD2	1:A:307:VAL:H	1.31	0.76
1:A:158:GLU:HG2	4:A:3008:HOH:O	1.85	0.76
1:B:1282:LYS:HA	1:B:1288:PHE:HD2	1.51	0.76
1:A:274:MET:HG3	1:A:293:ASP:HA	1.67	0.75
1:A:52:CYS:SG	4:A:3233:HOH:O	2.43	0.75
1:A:269:GLU:HA	1:A:272:LEU:HB2	1.70	0.74
1:B:1318:MET:HE2	1:B:1368:ILE:HD12	1.68	0.74
1:A:101:ASP:HA	1:B:1426:ARG:HH21	1.52	0.74
1:A:298:PHE:CD2	1:A:301:MET:HE2	2.23	0.73
1:B:1129:HIS:HD2	1:B:1131:MET:H	1.32	0.73
1:B:1113:ILE:HA	4:B:49:HOH:O	1.87	0.73
1:B:1230:GLU:HG2	1:B:1231:GLN:H	1.55	0.72
1:A:85:VAL:HG21	1:A:224:ILE:HG23	1.70	0.72
1:B:1339:ALA:HB1	1:B:1367:ILE:HD13	1.71	0.71
1:B:1004:LYS:HB3	1:B:1021:LYS:HD3	1.71	0.71
1:B:1081:ASN:HD21	1:B:1084:GLU:H	1.34	0.71
1:B:1318:MET:HE3	1:B:1318:MET:HA	1.72	0.71
1:A:129:HIS:CD2	1:A:131:MET:H	2.09	0.71
1:B:1077:ALA:HA	1:B:1224:ILE:HG21	1.71	0.71
1:A:188:ARG:HB2	1:A:191:LEU:HD12	1.72	0.70
1:A:129:HIS:HD2	1:A:131:MET:H	1.37	0.70
1:B:1123:ALA:HB3	4:B:187:HOH:O	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:LEU:HA	1:A:275:LEU:HG	1.74	0.70
1:A:305:HIS:CD2	1:A:307:VAL:H	2.10	0.69
1:A:117:ILE:HG13	4:A:3062:HOH:O	1.93	0.68
1:A:404:LYS:H	1:A:404:LYS:CD	2.00	0.68
1:A:100:TYR:HE1	1:B:1426:ARG:HB3	1.58	0.68
1:A:145:TYR:HE1	3:A:3000:NAI:O2A	1.76	0.68
1:A:294:LYS:HB3	1:A:301:MET:HE1	1.75	0.68
1:A:249:PHE:HE2	4:A:3292:HOH:O	1.77	0.67
1:A:252:ILE:O	1:A:256:ILE:HG13	1.94	0.67
1:A:163:ARG:HB2	4:A:3292:HOH:O	1.93	0.67
1:B:1290:ARG:HB2	4:B:24:HOH:O	1.95	0.66
1:B:1188:ARG:HG2	4:B:165:HOH:O	1.95	0.66
1:B:1230:GLU:CG	1:B:1231:GLN:H	2.09	0.65
1:B:1388:THR:O	1:B:1392:ILE:HG13	1.96	0.65
1:A:47:THR:HG22	1:B:1409:ARG:HH11	1.62	0.64
1:A:100:TYR:CE1	1:B:1426:ARG:HB3	2.33	0.64
1:B:1289:ARG:HA	4:B:324:HOH:O	1.96	0.64
1:A:7:LEU:HB2	1:B:1010:ASN:O	1.98	0.63
1:A:129:HIS:HD2	1:A:131:MET:N	1.97	0.63
1:B:1303:PHE:HB3	1:B:1357:LEU:HB3	1.80	0.63
1:A:116:GLN:HB3	4:A:3062:HOH:O	1.99	0.63
1:A:244:SER:HB3	1:B:1233:ALA:HB1	1.80	0.63
1:B:1007:LEU:HD21	1:B:1029:ILE:HG23	1.81	0.62
1:B:1081:ASN:C	1:B:1081:ASN:HD22	2.05	0.62
1:B:1008:THR:HG22	1:B:1016:GLU:HG2	1.81	0.62
1:A:282:LYS:HA	1:A:288:PHE:CD2	2.33	0.62
1:B:1020:LEU:HD22	1:B:1020:LEU:H	1.65	0.62
1:A:112:MET:HE2	1:A:186:TYR:CD1	2.35	0.61
1:B:1248:PRO:HG3	1:B:1395:TRP:CH2	2.35	0.61
1:B:1284:ILE:HD12	1:B:1287:PHE:H	1.65	0.61
1:A:232:ASN:ND2	1:A:235:THR:H	1.99	0.61
1:B:1288:PHE:HD1	1:B:1288:PHE:O	1.84	0.61
1:A:298:PHE:HB2	1:A:301:MET:HG3	1.82	0.60
1:B:1062:ASP:HA	1:B:1309:LYS:HB2	1.83	0.60
1:A:164:LEU:HG	4:A:3292:HOH:O	2.01	0.60
1:A:269:GLU:O	1:A:273:LYS:HB2	2.02	0.60
1:B:1272:LEU:HA	1:B:1275:LEU:CG	2.31	0.59
1:B:1352:PHE:HA	1:B:1357:LEU:HD12	1.83	0.59
1:B:1103:PHE:O	1:B:1107:VAL:HG23	2.03	0.59
1:B:1141:LEU:HD11	1:B:1171:MET:HE1	1.85	0.59
1:A:418:ARG:O	1:B:1072:PRO:HD3	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1269:GLU:O	1:B:1273:LYS:HB2	2.03	0.58
1:B:1082:TYR:CD1	1:B:1223:LEU:HB3	2.39	0.58
1:A:169:PRO:HG2	4:A:3124:HOH:O	2.03	0.58
1:B:1220:ASP:O	1:B:1224:ILE:HG13	2.01	0.58
1:B:1248:PRO:HG3	1:B:1395:TRP:CZ2	2.38	0.58
1:A:398:MET:SD	1:A:403:MET:HG3	2.42	0.58
1:B:1368:ILE:O	1:B:1372:MET:HG3	2.03	0.58
1:A:153:ASN:HD22	1:A:155:ARG:H	1.52	0.58
1:A:199:ASN:HD22	1:A:209:TYR:HD1	1.50	0.58
1:A:284:ILE:HD12	1:A:287:PHE:H	1.69	0.57
1:B:1164:LEU:HD13	1:B:1252:ILE:HG13	1.86	0.57
1:B:1120:LEU:HD12	4:B:187:HOH:O	2.05	0.57
1:B:1269:GLU:HA	1:B:1272:LEU:HB2	1.86	0.57
1:B:1278:ILE:HD13	1:B:1278:ILE:O	2.05	0.57
1:A:54:SER:HB2	4:A:3233:HOH:O	2.05	0.57
1:A:131:MET:HE2	1:A:380:THR:HB	1.87	0.57
1:A:101:ASP:HA	1:B:1426:ARG:NH2	2.19	0.57
1:B:1069:ARG:CD	1:B:1092:GLY:HA2	2.35	0.57
1:B:1072:PRO:HG2	1:B:1075:GLN:HE21	1.69	0.57
1:A:370:LYS:HG3	1:A:379:PHE:HZ	1.69	0.57
1:B:1069:ARG:HD2	1:B:1092:GLY:HA2	1.85	0.57
1:A:7:LEU:HD21	1:A:29:ILE:HG12	1.87	0.56
1:A:275:LEU:HD22	1:A:289:ARG:NE	2.16	0.56
1:B:1035:GLY:C	1:B:1037:LYS:H	2.13	0.56
1:B:1177:LYS:NZ	1:B:1182:GLN:HE21	2.04	0.56
1:A:267:ALA:HB3	1:A:270:ALA:HB2	1.88	0.56
1:A:212:ASN:HD22	1:A:213:PRO:N	2.03	0.56
1:A:305:HIS:CD2	1:A:307:VAL:HB	2.41	0.56
1:A:96:THR:HG23	1:A:99:GLN:NE2	2.21	0.55
1:A:21:LYS:HE2	1:A:21:LYS:H	1.72	0.55
1:A:232:ASN:C	1:A:232:ASN:HD22	2.13	0.55
1:B:1294:LYS:HB2	1:B:1297:SER:HA	1.87	0.55
1:A:230:GLU:CG	1:A:231:GLN:H	2.19	0.55
1:A:281:VAL:HG11	4:A:3270:HOH:O	2.05	0.55
1:A:289:ARG:HG2	1:A:290:ARG:H	1.71	0.55
1:A:408:PRO:HA	1:B:1046:PHE:O	2.07	0.55
1:B:1062:ASP:HA	1:B:1309:LYS:CB	2.37	0.55
1:A:278:ILE:HD13	1:A:278:ILE:O	2.07	0.55
1:A:131:MET:HG2	1:A:260:TRP:CD2	2.43	0.54
1:A:352:PHE:HA	1:A:357:LEU:HD23	1.89	0.54
1:A:153:ASN:ND2	1:A:155:ARG:H	2.05	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:SER:C	1:A:38:GLY:H	2.14	0.54
1:A:299:ARG:HA	1:A:303:PHE:O	2.07	0.54
1:B:1219:MET:O	1:B:1223:LEU:HD13	2.07	0.54
1:B:1012:ASP:HA	4:B:78:HOH:O	2.07	0.54
1:B:1020:LEU:CD2	1:B:1030:ASP:HB2	2.39	0.53
1:B:1204:THR:HG22	1:B:1206:CYS:H	1.73	0.53
1:A:289:ARG:HG2	1:A:290:ARG:N	2.24	0.53
1:B:1303:PHE:CE1	1:B:1346:ALA:HB2	2.44	0.53
1:A:20:LEU:HB2	1:A:28:VAL:HG23	1.90	0.53
1:B:1019:VAL:HA	1:B:1028:VAL:O	2.08	0.53
1:A:153:ASN:HD21	1:A:155:ARG:HD3	1.73	0.53
1:B:1289:ARG:HG2	1:B:1290:ARG:H	1.73	0.53
1:A:94:LYS:NZ	1:B:1421:LYS:O	2.41	0.53
1:A:309:LYS:HZ3	1:A:356:LYS:HE3	1.73	0.52
1:A:212:ASN:HD22	1:A:212:ASN:C	2.17	0.52
1:A:47:THR:O	1:B:1409:ARG:NE	2.42	0.52
1:B:1057:THR:HG23	1:B:1067:LEU:O	2.08	0.52
1:B:1151:VAL:HB	1:B:1399:HIS:CE1	2.45	0.52
1:B:1294:LYS:HZ2	1:B:1297:SER:HA	1.74	0.52
1:A:223:LEU:CD2	1:A:386:ALA:HB2	2.40	0.52
1:A:212:ASN:ND2	1:A:214:ILE:H	2.07	0.52
1:A:278:ILE:O	1:A:282:LYS:HB3	2.10	0.52
1:A:145:TYR:CE1	3:A:3000:NAI:O2A	2.61	0.52
1:B:1177:LYS:NZ	1:B:1182:GLN:NE2	2.58	0.52
3:A:3000:NAI:H42N	4:A:3008:HOH:O	2.10	0.52
1:B:1230:GLU:HG2	1:B:1231:GLN:N	2.24	0.52
1:B:1034:LEU:HD22	1:B:1039:VAL:CG1	2.40	0.51
1:A:153:ASN:HD22	1:A:153:ASN:C	2.19	0.51
1:A:318:MET:HA	1:A:318:MET:HE3	1.92	0.51
1:B:1243:SER:HB2	1:B:1405:ILE:HA	1.92	0.51
1:B:1345:ILE:HG23	1:B:1349:ASP:HB2	1.92	0.51
1:A:169:PRO:HA	1:A:385:MET:SD	2.50	0.51
3:A:3000:NAI:H51A	3:A:3000:NAI:H8A	1.93	0.51
1:A:275:LEU:HA	1:A:289:ARG:HG3	1.93	0.51
1:B:1069:ARG:NE	1:B:1092:GLY:HA2	2.26	0.51
1:A:68:HIS:HB2	1:A:76:LEU:HD11	1.91	0.51
1:A:212:ASN:HB3	1:A:215:LEU:HG	1.91	0.51
1:B:1153:ASN:C	1:B:1153:ASN:HD22	2.19	0.51
1:B:1275:LEU:HD22	4:B:24:HOH:O	2.09	0.51
1:B:1020:LEU:HD21	1:B:1030:ASP:HB2	1.92	0.51
1:B:1422:SER:OG	1:B:1424:ILE:HG12	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1081:ASN:HD22	1:B:1084:GLU:H	1.57	0.50
1:B:1351:TYR:O	1:B:1355:LYS:HD3	2.11	0.50
1:A:223:LEU:HD22	1:A:386:ALA:HB2	1.94	0.50
1:A:335:LEU:HD23	1:A:335:LEU:H	1.76	0.50
1:B:1096:THR:H	1:B:1099:GLN:HB2	1.76	0.50
1:B:1204:THR:CG2	1:B:1206:CYS:SG	3.00	0.50
1:A:288:PHE:CD1	1:A:288:PHE:O	2.64	0.50
1:B:1081:ASN:HD21	1:B:1084:GLU:N	2.04	0.50
1:B:1327:LYS:O	1:B:1327:LYS:HG2	2.11	0.50
1:A:153:ASN:HD22	1:A:155:ARG:N	2.10	0.50
1:A:3:THR:C	1:A:5:ALA:H	2.20	0.50
1:B:1133:VAL:HG21	4:B:96:HOH:O	2.12	0.50
1:B:1287:PHE:HA	1:B:1291:ALA:CB	2.37	0.50
1:B:1288:PHE:O	1:B:1288:PHE:CD1	2.65	0.50
1:B:1289:ARG:HG2	1:B:1290:ARG:N	2.27	0.50
1:A:288:PHE:O	1:A:288:PHE:HD1	1.94	0.50
1:B:1167:LYS:HE3	3:B:3001:NAI:O2N	2.11	0.50
1:B:1365:SER:O	1:B:1369:LEU:HG	2.12	0.50
1:A:318:MET:HE2	1:A:368:ILE:CD1	2.42	0.49
1:A:309:LYS:HZ2	1:A:356:LYS:HE3	1.77	0.49
1:A:343:GLU:O	1:A:347:LEU:HG	2.13	0.49
1:A:46:PHE:O	1:B:1408:PRO:HA	2.12	0.49
1:B:1060:ASP:HB3	1:B:1065:ILE:HB	1.95	0.49
1:A:134:MET:HE3	1:A:134:MET:HA	1.94	0.49
1:A:277:GLU:O	1:A:281:VAL:HG22	2.12	0.49
1:B:1007:LEU:HD22	1:B:1029:ILE:HG12	1.95	0.49
1:B:1287:PHE:O	1:B:1292:LYS:HB2	2.12	0.49
1:B:1354:GLU:HB3	1:B:1355:LYS:HD2	1.95	0.49
1:B:1387:ARG:HG3	1:B:1391:TRP:CE2	2.48	0.49
1:B:1086:CYS:SG	1:B:1389:VAL:HG21	2.52	0.49
1:B:1260:TRP:HA	4:B:321:HOH:O	2.13	0.49
1:A:336:LEU:HG	4:A:3151:HOH:O	2.12	0.48
1:A:409:ARG:HB2	1:B:1050:ALA:HA	1.94	0.48
1:A:174:MET:HG3	4:A:3030:HOH:O	2.12	0.48
1:A:272:LEU:HA	1:A:275:LEU:CG	2.42	0.48
1:A:349:ASP:HB3	1:A:352:PHE:CD1	2.48	0.48
1:B:1164:LEU:O	1:B:1168:MET:HG2	2.13	0.48
1:A:88:ILE:HG12	1:A:94:LYS:HA	1.95	0.48
1:A:207:GLU:HB2	1:A:208:PRO:HD2	1.96	0.48
1:A:150:ASP:HB3	1:A:156:HIS:CD2	2.48	0.48
1:B:1322:CYS:O	1:B:1326:LEU:HD13	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:VAL:HB	1:A:399:HIS:CE1	2.48	0.48
1:A:230:GLU:HG2	1:A:231:GLN:H	1.78	0.48
1:B:1204:THR:HG22	1:B:1206:CYS:SG	2.54	0.48
1:A:81:ASN:HA	1:A:224:ILE:HD11	1.94	0.47
1:A:29:ILE:O	1:A:31:ILE:HG23	2.14	0.47
1:A:35:GLY:HA2	1:A:40:PHE:CE2	2.49	0.47
1:B:1294:LYS:NZ	1:B:1298:PHE:H	2.12	0.47
1:A:15:VAL:HA	4:A:3205:HOH:O	2.13	0.47
1:B:1243:SER:O	1:B:1405:ILE:HG12	2.14	0.47
1:B:1326:LEU:HG	1:B:1333:ASP:OD1	2.15	0.47
1:B:1303:PHE:HE1	1:B:1346:ALA:HB2	1.79	0.47
1:A:318:MET:HE2	1:A:368:ILE:HD12	1.97	0.47
1:B:1214:ILE:HG12	1:B:1217:ARG:NH1	2.29	0.47
1:B:1269:GLU:HB2	1:B:1273:LYS:CA	2.44	0.47
1:B:1300:LEU:HD13	1:B:1300:LEU:O	2.15	0.47
1:B:1264:HIS:ND1	1:B:1265:GLY:N	2.63	0.47
1:A:399:HIS:NE2	1:A:403:MET:SD	2.88	0.47
1:B:1277:GLU:HA	4:B:246:HOH:O	2.14	0.47
1:A:7:LEU:HD12	1:A:17:LEU:HB2	1.97	0.46
1:A:266:GLY:O	1:A:267:ALA:HB2	2.16	0.46
1:A:413:THR:HB	1:B:1055:LYS:NZ	2.31	0.46
1:B:1288:PHE:CD1	1:B:1288:PHE:C	2.93	0.46
1:B:1113:ILE:HD13	1:B:1118:THR:OG1	2.15	0.46
1:B:1149:LEU:HD22	1:B:1247:ASN:HB2	1.96	0.46
1:A:239:ARG:HD3	4:A:3022:HOH:O	2.15	0.46
1:B:1069:ARG:HG3	1:B:1088:ILE:HG22	1.98	0.46
1:B:1275:LEU:HA	1:B:1289:ARG:HG3	1.98	0.46
1:B:1272:LEU:CA	1:B:1275:LEU:HG	2.36	0.46
1:A:290:ARG:HG2	1:A:290:ARG:O	2.14	0.46
1:B:1112:MET:HE2	1:B:1112:MET:HB3	1.76	0.46
1:A:212:ASN:HD22	1:A:213:PRO:CD	2.29	0.46
1:A:417:LYS:HD3	1:A:417:LYS:C	2.41	0.46
1:A:96:THR:OG1	1:A:99:GLN:HG3	2.16	0.46
1:A:185:VAL:HB	1:A:200:MET:HA	1.98	0.46
1:A:370:LYS:HG3	1:A:379:PHE:CZ	2.49	0.46
1:B:1355:LYS:HD2	1:B:1355:LYS:N	2.31	0.46
1:A:243:SER:O	1:A:403:MET:HE3	2.15	0.46
1:A:289:ARG:CG	1:A:290:ARG:H	2.28	0.46
1:A:344:ASN:OD1	1:A:348:ASN:ND2	2.49	0.46
1:B:1404:LYS:HE3	4:B:154:HOH:O	2.15	0.46
1:A:24:LEU:HD11	1:B:1416:GLU:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:PHE:HB3	4:A:3144:HOH:O	2.15	0.45
1:B:1188:ARG:HB3	1:B:1190:ASP:OD1	2.15	0.45
1:B:1157:ARG:NH2	4:B:337:HOH:O	2.49	0.45
1:A:290:ARG:HA	1:A:293:ASP:CB	2.47	0.45
1:A:411:LEU:HD23	1:B:1028:VAL:HG22	1.98	0.45
1:A:59:ILE:HG21	1:A:307:VAL:CG1	2.46	0.45
1:B:1290:ARG:O	1:B:1290:ARG:HG2	2.15	0.45
1:B:1282:LYS:HA	1:B:1288:PHE:CD2	2.41	0.45
1:B:1266:GLY:O	1:B:1267:ALA:HB2	2.17	0.44
1:A:294:LYS:NZ	1:A:301:MET:HE3	2.32	0.44
1:B:1269:GLU:HB2	1:B:1273:LYS:HB2	1.98	0.44
1:A:247:ASN:HB2	1:A:248:PRO:HD2	2.00	0.44
1:B:1064:GLY:O	1:B:1313:PRO:HB3	2.18	0.44
1:B:1159:ILE:O	1:B:1163:ARG:HG3	2.17	0.44
1:B:1225:LEU:HD12	1:B:1318:MET:SD	2.58	0.44
1:A:138:THR:HG21	1:A:252:ILE:HG22	1.99	0.44
1:A:308:TYR:HB3	1:A:311:TYR:O	2.17	0.44
1:B:1167:LYS:O	1:B:1171:MET:HB2	2.17	0.44
1:A:7:LEU:HD12	1:A:7:LEU:C	2.43	0.44
1:A:42:PHE:HB2	1:B:1028:VAL:CG2	2.47	0.44
1:A:370:LYS:HE3	1:A:379:PHE:HE1	1.82	0.44
1:B:1231:GLN:HE22	1:B:1239:ARG:NE	2.15	0.44
1:A:32:ARG:HD2	1:B:1042:PHE:CD1	2.53	0.44
1:A:412:TYR:HD1	1:B:1024:LEU:HB2	1.83	0.44
1:B:1298:PHE:HD2	1:B:1301:MET:HE2	1.82	0.44
1:A:286:GLU:OE1	1:A:291:ALA:HB2	2.18	0.44
1:A:117:ILE:HG12	1:A:144:PHE:CD1	2.53	0.44
1:A:411:LEU:O	1:B:1053:GLU:HG3	2.18	0.43
1:A:272:LEU:HD12	1:A:272:LEU:N	2.33	0.43
1:A:288:PHE:HD1	1:A:288:PHE:H	1.66	0.43
1:B:1034:LEU:HD22	1:B:1039:VAL:HG11	2.00	0.43
1:B:1230:GLU:CG	1:B:1231:GLN:N	2.79	0.43
1:A:275:LEU:CD2	1:A:289:ARG:HE	2.21	0.43
1:B:1069:ARG:HG3	1:B:1088:ILE:CG2	2.49	0.43
1:B:1106:THR:O	1:B:1110:HIS:HD2	2.02	0.43
1:A:150:ASP:H	1:A:156:HIS:HD2	1.64	0.43
1:A:287:PHE:HA	1:A:291:ALA:CB	2.41	0.43
1:A:290:ARG:HA	1:A:293:ASP:HB2	2.00	0.43
1:A:412:TYR:CE1	1:B:1024:LEU:HD12	2.52	0.43
1:B:1003:THR:O	1:B:1019:VAL:HG11	2.18	0.43
1:B:1386:ALA:O	1:B:1389:VAL:HG12	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1129:HIS:CD2	1:B:1130:PRO:HD2	2.54	0.43
1:A:36:SER:C	1:A:38:GLY:N	2.77	0.43
1:A:320:GLU:HB3	4:A:3165:HOH:O	2.19	0.43
1:B:1149:LEU:HD22	1:B:1247:ASN:CB	2.49	0.43
1:B:1273:LYS:NZ	1:B:1273:LYS:HB3	2.33	0.43
1:B:1082:TYR:CE2	1:B:1086:CYS:SG	3.12	0.43
1:A:418:ARG:HH11	1:B:1069:ARG:HB3	1.84	0.42
1:B:1114:HIS:CE1	1:B:1116:GLN:HB2	2.54	0.42
1:A:17:LEU:HD21	1:A:34:LEU:HD21	2.01	0.42
1:A:286:GLU:O	1:A:291:ALA:HB3	2.20	0.42
1:B:1059:ILE:HG22	1:B:1060:ASP:N	2.34	0.42
1:A:32:ARG:HD2	1:B:1042:PHE:CE1	2.54	0.42
1:A:106:THR:HG23	1:A:109:ARG:HH21	1.83	0.42
1:A:250:ALA:HB1	1:B:1257:ALA:HB1	2.02	0.42
1:A:273:LYS:O	1:A:276:GLU:HB3	2.20	0.42
1:B:1294:LYS:HE2	1:B:1357:LEU:HD13	2.01	0.42
1:B:1007:LEU:HD11	1:B:1029:ILE:HG21	2.01	0.42
1:B:1059:ILE:HA	1:B:1065:ILE:O	2.19	0.42
1:B:1138:THR:HG22	1:B:1253:ALA:HB2	2.01	0.42
1:B:1153:ASN:ND2	1:B:1155:ARG:H	2.17	0.42
1:B:1356:LYS:HD3	4:B:558:HOH:O	2.19	0.42
1:A:86:CYS:HA	1:A:389:VAL:HG21	2.02	0.42
1:B:1329:LEU:HD13	1:B:1371:ALA:O	2.19	0.42
1:B:1131:MET:HE2	1:B:1380:THR:CG2	2.50	0.42
1:B:1314:ARG:HB2	1:B:1361:VAL:HB	2.02	0.42
1:A:21:LYS:H	1:A:21:LYS:CE	2.31	0.42
1:A:305:HIS:HD2	1:A:307:VAL:N	2.06	0.42
1:B:1034:LEU:HD22	1:B:1039:VAL:HG13	2.02	0.42
1:B:1286:GLU:OE1	1:B:1291:ALA:HB2	2.20	0.42
1:A:7:LEU:CD1	1:A:17:LEU:HB2	2.50	0.41
1:A:20:LEU:HB2	1:A:28:VAL:CG2	2.50	0.41
1:A:87:TYR:CD1	1:A:95:PRO:HB3	2.54	0.41
1:B:1199:ASN:O	1:B:1203:SER:HB3	2.20	0.41
1:B:1274:MET:CB	4:B:324:HOH:O	2.68	0.41
1:A:339:ALA:HB1	1:A:367:ILE:HD13	2.03	0.41
1:B:1018:ASP:HB2	1:B:1030:ASP:HB3	2.01	0.41
1:A:368:ILE:CG2	1:A:372:MET:HE3	2.51	0.41
1:B:1029:ILE:O	1:B:1031:ILE:HG23	2.21	0.41
1:B:1178:TYR:HB2	1:B:1184:PHE:CE1	2.55	0.41
1:A:212:ASN:HD22	1:A:214:ILE:H	1.66	0.41
1:B:1172:ALA:O	1:B:1175:CYS:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1072:PRO:HB2	1:B:1075:GLN:HG3	2.03	0.41
1:A:288:PHE:CD1	1:A:288:PHE:C	2.99	0.41
1:A:303:PHE:HB3	1:A:357:LEU:HB3	2.03	0.41
1:B:1272:LEU:HD12	1:B:1272:LEU:N	2.35	0.41
1:A:250:ALA:CB	1:B:1257:ALA:HB1	2.51	0.41
1:B:1113:ILE:HD12	1:B:1113:ILE:O	2.21	0.41
1:B:1269:GLU:HB2	1:B:1273:LYS:HA	2.02	0.41
1:B:1426:ARG:NH1	4:B:53:HOH:O	2.53	0.41
1:A:83:LEU:O	1:A:86:CYS:HB2	2.21	0.41
1:A:98:GLU:HB2	4:A:3166:HOH:O	2.20	0.41
1:B:1001:ALA:HA	1:B:1004:LYS:CG	2.51	0.41
1:B:1231:GLN:HE22	1:B:1239:ARG:CD	2.34	0.41
1:A:141:LEU:HD11	1:A:171:MET:HE1	2.03	0.41
1:A:412:TYR:CD1	1:B:1024:LEU:HB2	2.56	0.41
1:B:1274:MET:SD	1:B:1277:GLU:OE2	2.79	0.41
1:A:35:GLY:HA2	1:A:40:PHE:CD2	2.55	0.40
1:B:1231:GLN:HE22	1:B:1239:ARG:HD2	1.86	0.40
1:B:1263:ALA:HB2	4:B:321:HOH:O	2.20	0.40
1:B:1299:ARG:HA	1:B:1303:PHE:O	2.21	0.40
1:A:274:MET:C	1:A:276:GLU:H	2.28	0.40
1:A:294:LYS:HB2	1:A:297:SER:HA	2.02	0.40
1:B:1131:MET:SD	1:B:1381:VAL:HG22	2.61	0.40
1:A:125:ARG:NH2	4:A:3083:HOH:O	2.54	0.40
1:A:260:TRP:CE2	1:A:264:HIS:HD2	2.40	0.40
1:A:274:MET:HB2	1:A:293:ASP:OD2	2.21	0.40
3:A:3000:NAI:H3B	3:A:3000:NAI:PA	2.60	0.40
1:A:238:VAL:HG13	1:A:251:CYS:HB3	2.03	0.40
1:A:348:ASN:HA	1:A:353:ILE:HD11	2.04	0.40
1:B:1349:ASP:HA	1:B:1350:PRO:HD3	1.98	0.40

There are no symmetry-related clashes.

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	424/427 (99%)	373 (88%)	43 (10%)	8 (2%)	6	11
1	B	424/427 (99%)	369 (87%)	48 (11%)	7 (2%)	7	13
All	All	848/854 (99%)	742 (88%)	91 (11%)	15 (2%)	6	12

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	ALA
1	A	294	LYS
1	A	422	SER
1	B	1294	LYS
1	B	1330	GLY
1	A	290	ARG
1	A	401	ASP
1	B	1150	ASP
1	B	1286	GLU
1	B	1290	ARG
1	A	286	GLU
1	A	402	GLY
1	B	1012	ASP
1	B	1332	LYS
1	A	424	ILE

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	360/361 (100%)	344 (96%)	16 (4%)	25	50
1	B	360/361 (100%)	345 (96%)	15 (4%)	26	52
All	All	720/722 (100%)	689 (96%)	31 (4%)	26	51

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LYS

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Mol	Chain	Res	Type
1	A	83	LEU
1	A	134	MET
1	A	153	ASN
1	A	230	GLU
1	A	232	ASN
1	A	272	LEU
1	A	278	ILE
1	A	282	LYS
1	A	284	ILE
1	A	288	PHE
1	A	294	LYS
1	A	299	ARG
1	A	357	LEU
1	A	399	HIS
1	A	404	LYS
1	B	1007	LEU
1	B	1081	ASN
1	B	1113	ILE
1	B	1232	ASN
1	B	1273	LYS
1	B	1278	ILE
1	B	1282	LYS
1	B	1284	ILE
1	B	1288	PHE
1	B	1294	LYS
1	B	1299	ARG
1	B	1335	LEU
1	B	1354	GLU
1	B	1404	LYS
1	B	1425	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	99	GLN
1	A	129	HIS
1	A	153	ASN
1	A	156	HIS
1	A	199	ASN
1	A	212	ASN
1	A	232	ASN

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Mol	Chain	Res	Type
1	A	305	HIS
1	A	323	HIS
1	A	348	ASN
1	B	1026	GLN
1	B	1081	ASN
1	B	1097	GLN
1	B	1110	HIS
1	B	1129	HIS
1	B	1153	ASN
1	B	1156	HIS
1	B	1182	GLN
1	B	1231	GLN
1	B	1232	ASN
1	B	1264	HIS
1	B	1305	HIS
1	B	1323	HIS
1	B	1399	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	NAI	B	3001	-	47,48,48	2.17	10 (21%)	64,73,73	1.71	5 (7%)
2	SO4	A	2004	-	4,4,4	0.66	0	6,6,6	0.36	0
2	SO4	A	2006	-	4,4,4	0.70	0	6,6,6	0.44	0
2	SO4	B	2003	-	4,4,4	0.70	0	6,6,6	3.63	4 (66%)
2	SO4	B	2001	-	4,4,4	0.57	0	6,6,6	3.57	4 (66%)
3	NAI	A	3000	-	47,48,48	2.01	7 (14%)	64,73,73	2.00	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	NAI	A	3000	-	-	8/29/72/72	0/5/5/5
3	NAI	B	3001	-	-	9/29/72/72	0/5/5/5

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3001	NAI	PN-O3	10.90	1.71	1.59
3	A	3000	NAI	PN-O3	9.67	1.69	1.59
3	B	3001	NAI	C2N-C3N	-4.06	1.24	1.35
3	A	3000	NAI	C2N-C3N	-3.91	1.24	1.35
3	A	3000	NAI	PN-O2N	-3.73	1.37	1.50
3	A	3000	NAI	C7N-C3N	-3.66	1.40	1.48
3	B	3001	NAI	PN-O2N	-3.34	1.39	1.50
3	A	3000	NAI	C4N-C5N	-3.17	1.40	1.49
3	B	3001	NAI	C6N-N1N	2.79	1.44	1.37
3	B	3001	NAI	PA-O3	-2.77	1.56	1.59
3	A	3000	NAI	C2A-N1A	2.68	1.38	1.33
3	A	3000	NAI	C6N-N1N	2.67	1.43	1.37
3	B	3001	NAI	C4N-C5N	-2.45	1.42	1.49
3	B	3001	NAI	C7N-C3N	-2.43	1.43	1.48
3	B	3001	NAI	C2A-N1A	2.22	1.37	1.33
3	B	3001	NAI	C1D-N1N	-2.11	1.40	1.46
3	B	3001	NAI	C2D-C1D	-2.03	1.47	1.53

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	3000	NAI	C1D-N1N-C2N	-8.87	106.52	121.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3001	NAI	C6N-N1N-C2N	8.10	127.98	119.32
3	A	3000	NAI	C6N-N1N-C2N	7.84	127.71	119.32
2	B	2003	SO4	O4-S-O1	-5.90	78.70	109.56
2	B	2001	SO4	O4-S-O1	-5.62	80.16	109.56
3	B	3001	NAI	C1D-N1N-C2N	-5.50	112.08	121.14
3	B	3001	NAI	O7N-C7N-C3N	-4.80	111.86	120.90
3	A	3000	NAI	C3N-C2N-N1N	-4.59	116.46	123.20
2	B	2001	SO4	O3-S-O1	-4.53	85.86	109.56
2	B	2003	SO4	O4-S-O2	-4.48	86.15	109.56
3	A	3000	NAI	O7N-C7N-C3N	-4.47	112.47	120.90
2	B	2003	SO4	O4-S-O3	-3.96	86.70	108.54
3	B	3001	NAI	C3N-C2N-N1N	-3.75	117.70	123.20
3	B	3001	NAI	C3N-C7N-N7N	3.66	124.16	117.67
2	B	2001	SO4	O2-S-O1	-3.37	85.33	109.06
2	B	2001	SO4	O4-S-O2	3.24	126.50	109.56
3	A	3000	NAI	O5B-C5B-C4B	3.20	119.88	108.99
3	A	3000	NAI	O4B-C4B-C5B	-2.89	100.07	109.33
3	A	3000	NAI	C3N-C7N-N7N	2.71	122.49	117.67
2	B	2003	SO4	O2-S-O1	2.52	126.83	109.06
3	A	3000	NAI	C1D-N1N-C6N	2.36	125.77	120.77
3	A	3000	NAI	O4B-C1B-N9A	-2.20	103.87	108.09

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	3000	NAI	C4B-C5B-O5B-PA
3	A	3000	NAI	C5D-O5D-PN-O3
3	A	3000	NAI	C5D-O5D-PN-O1N
3	A	3000	NAI	C5D-O5D-PN-O2N
3	B	3001	NAI	C5D-O5D-PN-O3
3	B	3001	NAI	C5D-O5D-PN-O1N
3	B	3001	NAI	C5D-O5D-PN-O2N
3	B	3001	NAI	O4D-C4D-C5D-O5D
3	A	3000	NAI	O4D-C4D-C5D-O5D
3	B	3001	NAI	C3D-C4D-C5D-O5D
3	B	3001	NAI	O4D-C1D-N1N-C2N
3	B	3001	NAI	PN-O3-PA-O2A
3	A	3000	NAI	C3D-C4D-C5D-O5D
3	A	3000	NAI	O4D-C1D-N1N-C2N
3	B	3001	NAI	PN-O3-PA-O1A
3	A	3000	NAI	O4B-C4B-C5B-O5B

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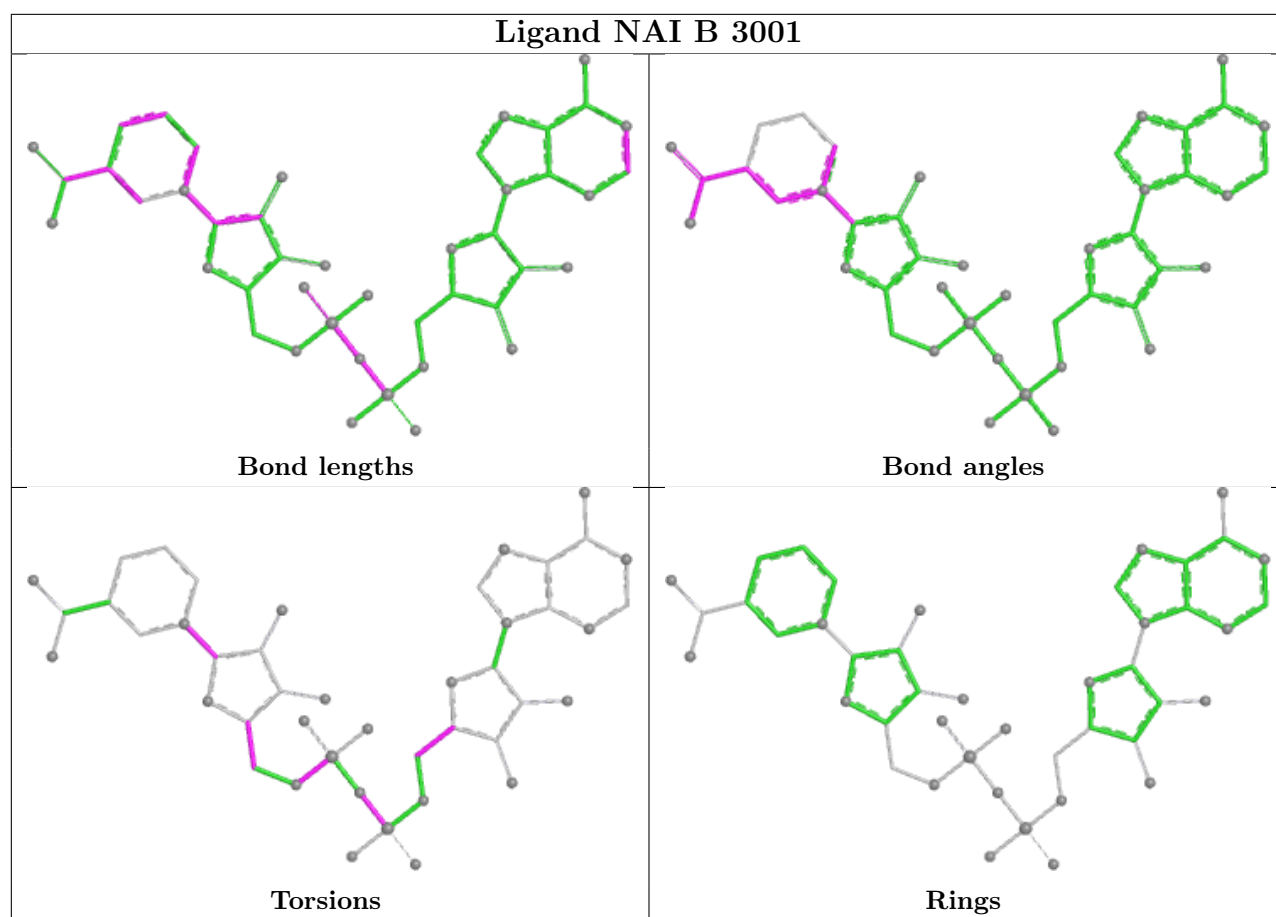
Mol	Chain	Res	Type	Atoms
3	B	3001	NAI	O4B-C4B-C5B-O5B

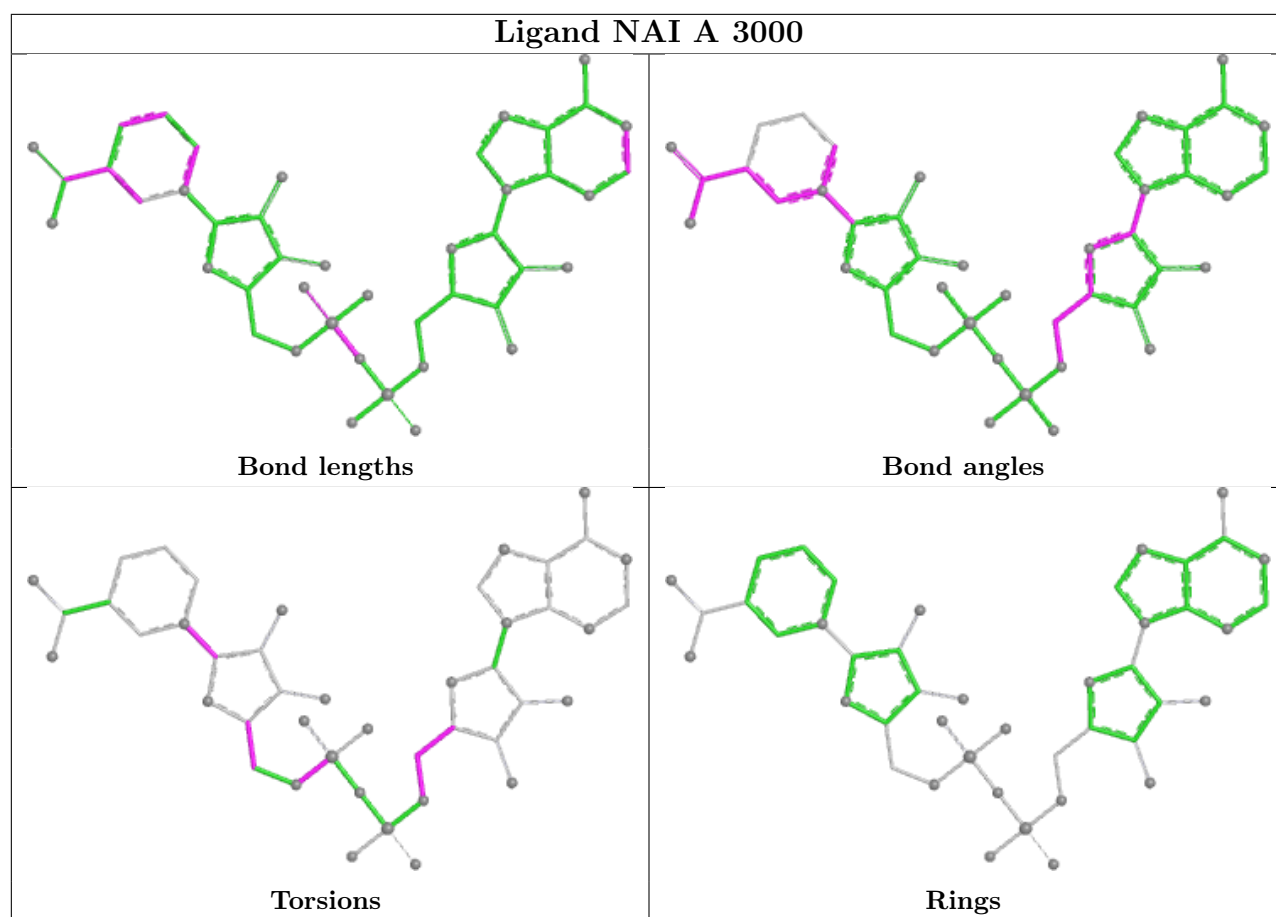
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	3001	NAI	1	0
3	A	3000	NAI	6	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.