



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:16 AM GMT

PDB ID : 3DXK
Title : Structure of Bos Taurus Arp2/3 Complex with Bound Inhibitor CK0944636
Authors : Nolen, B.J.; Tomasevic, N.; Russell, A.; Pierce, D.W.; Jia, Z.; Hartman, J.; Sakowicz, R.; Pollard, T.D.
Deposited on : 2008-07-24
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

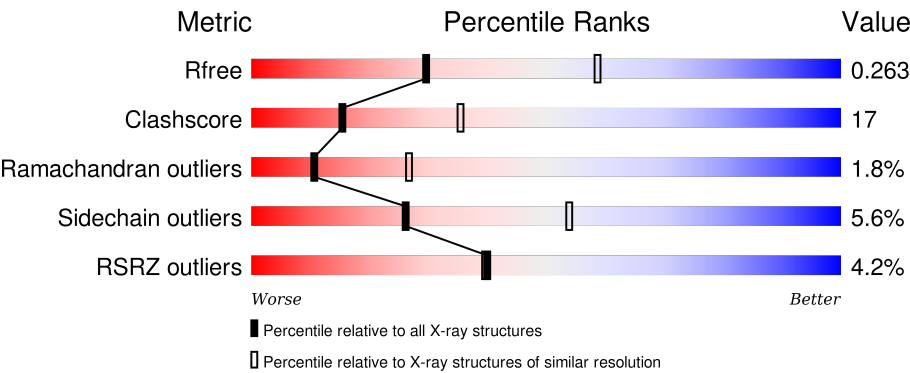
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	418	<div><div>3%</div><div>65%27%5%</div></div>
2	B	394	<div><div>4%</div><div>27%21%46%</div></div>
3	C	372	<div><div>3%</div><div>61%28%9%</div></div>
4	D	300	<div><div>%</div><div>71%23%6%</div></div>
5	E	178	<div><div>9%</div><div>54%37%6%</div></div>

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Mol	Chain	Length	Quality of chain
6	F	168	<div><div><div>%</div><div><div></div><div>77%</div><div>20%</div><div>••</div></div></div></div>
7	G	151	<div><div><div>10%</div><div><div></div><div>56%</div><div>28%</div><div>••</div><div>11%</div></div></div></div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 13598 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 Actin-related protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	0	0	0
			3176	2042	531	588	15			

- Molecule 2 is a protein called Actin-related protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	211	Total	C	N	O	S	0	0	0
			1655	1058	289	304	4			

- Molecule 3 is a protein called Actin-related protein 2/3 complex subunit 1B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	339	Total	C	N	O	S	0	0	0
			2618	1665	458	476	19			

- Molecule 4 is a protein called Actin-related protein 2/3 complex subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	282	Total	C	N	O	S	0	0	0
			2279	1449	395	427	8			

- Molecule 5 is a protein called Actin-related protein 2/3 complex subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	174	Total	C	N	O	S	0	0	0
			1415	908	236	262	9			

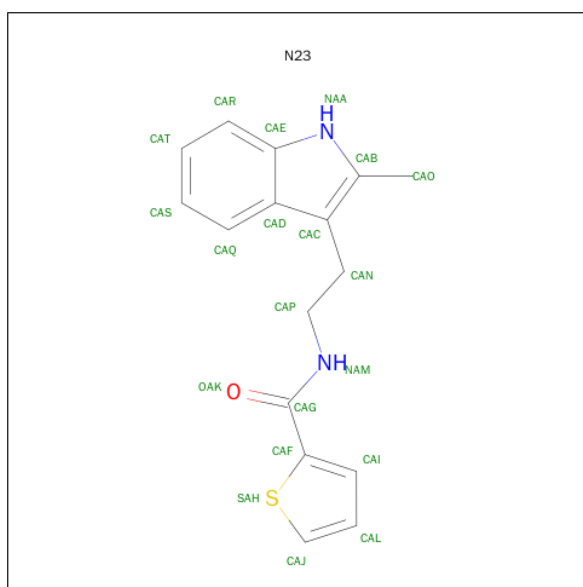
- Molecule 6 is a protein called Actin-related protein 2/3 complex subunit 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	167	Total	C	N	O	S	0	0	0
			1371	875	239	248	9			

- Molecule 7 is a protein called Actin-related protein 2/3 complex subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	134	Total	C	N	O	S	0	0	0
			1022	641	180	198	3			

- Molecule 8 is N-[2-(2-METHYL-1H-INDOL-3-YL)ETHYL]THIOPHENE-2-CARBOXAMIDE (three-letter code: N23) (formula: C₁₆H₁₆N₂OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	1	Total	C	N	O	S	0	0
			20	16	2	1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	8	Total	O	0	0
			8	8		
9	B	3	Total	O	0	0
			3	3		
9	C	7	Total	O	0	0
			7	7		
9	D	10	Total	O	0	0
			10	10		

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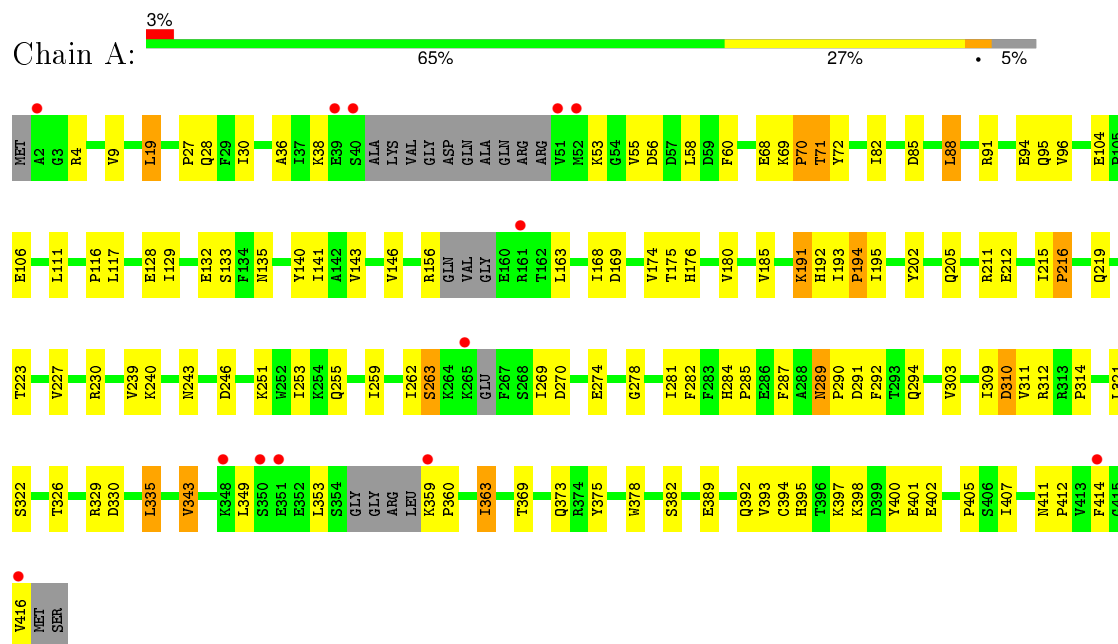
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	E	1	Total 1	O 1	0	0
9	F	12	Total 12	O 12	0	0
9	G	1	Total 1	O 1	0	0

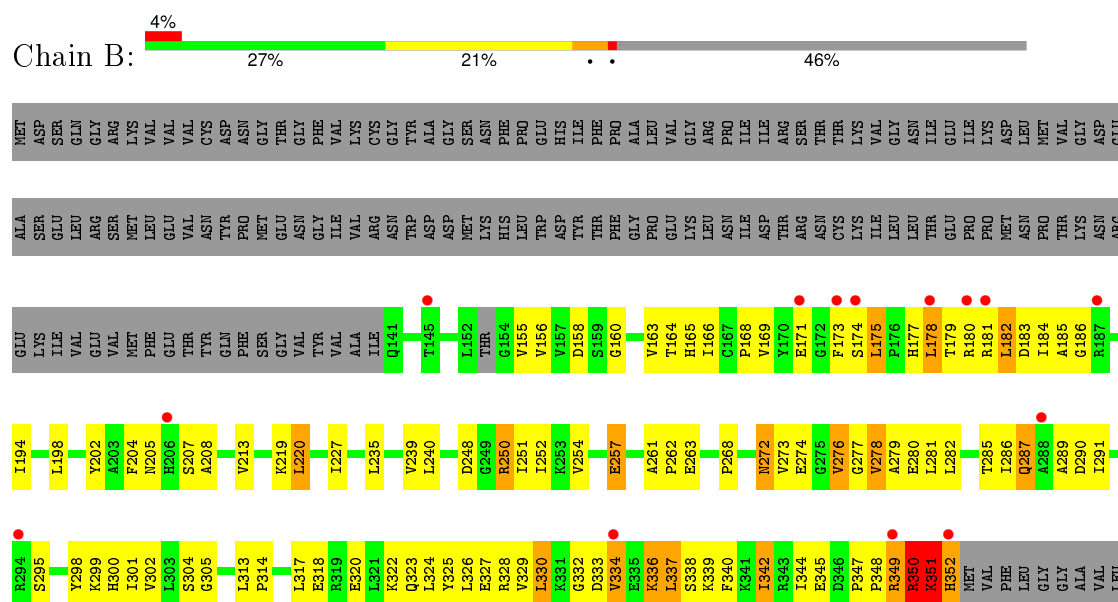
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Actin-related protein 3

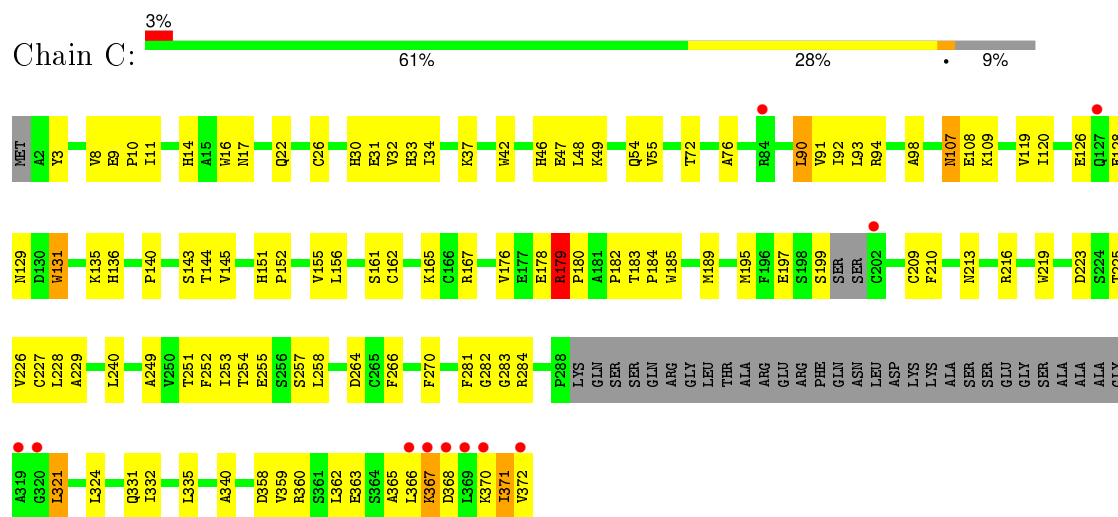


• Molecule 2: Actin-related protein 2

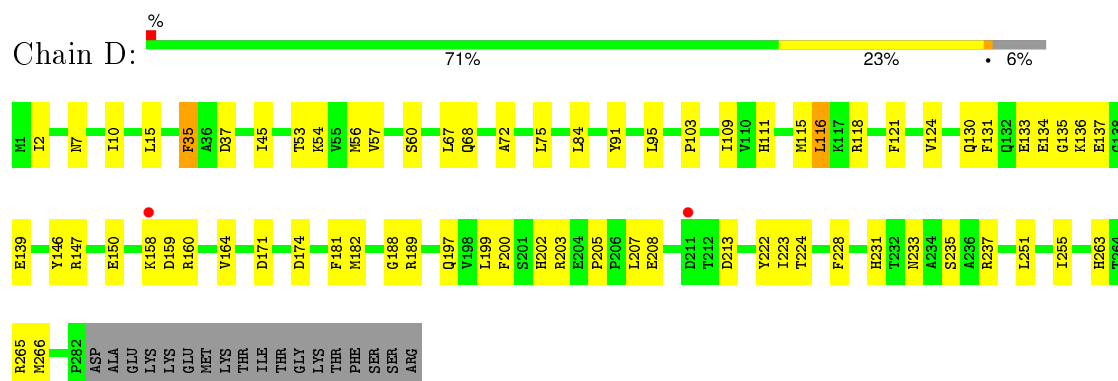


ALA ASP ILE MET LYS LYS ASP ASN PHE TRP MET THR ARG GLN GLU TYR GLN GLY VAL ARG VAL LEU GLU LYS LEU GLY VAL THR VAL ARG

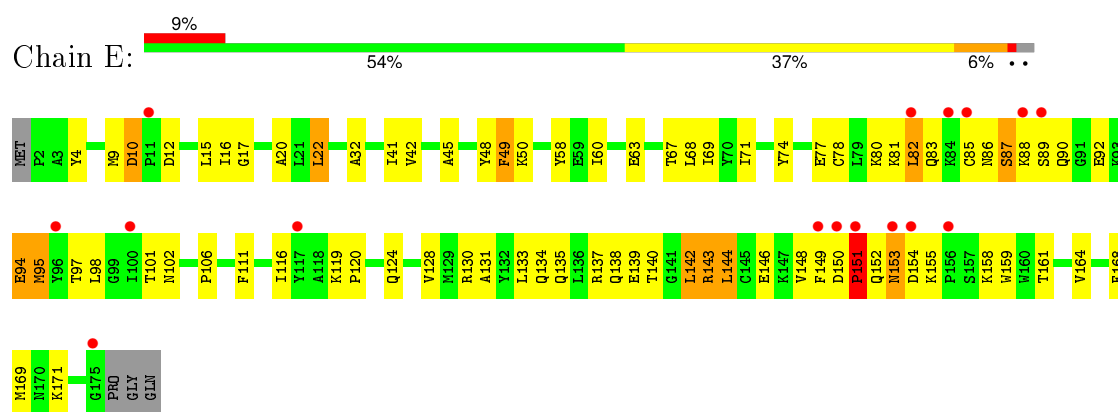
• Molecule 3: Actin-related protein 2/3 complex subunit 1B



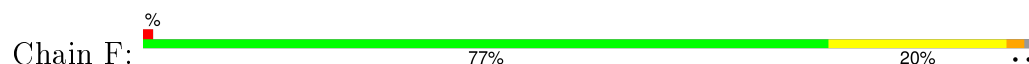
• Molecule 4: Actin-related protein 2/3 complex subunit 2

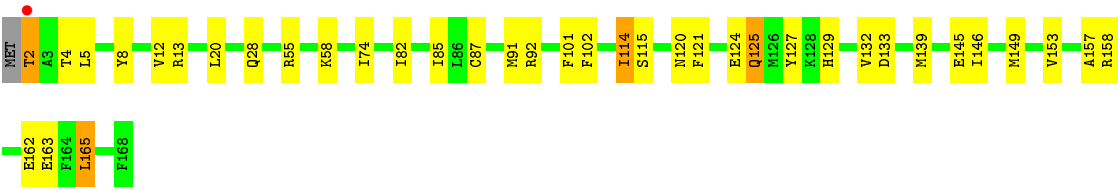


• Molecule 5: Actin-related protein 2/3 complex subunit 3

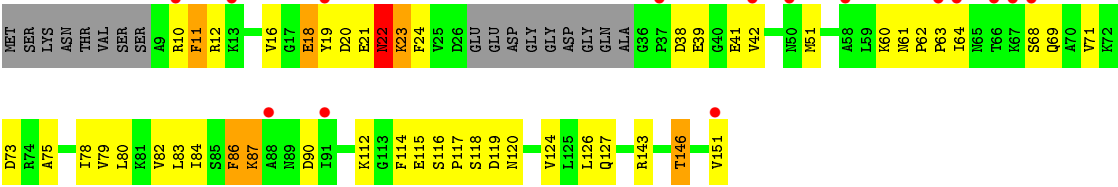


• Molecule 6: Actin-related protein 2/3 complex subunit 4





● Molecule 7: Actin-related protein 2/3 complex subunit 5



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.60 Å 130.66 Å 203.96 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 46.38 – 2.72	Depositor EDS
% Data completeness (in resolution range)	97.8 (30.00-2.70) 89.1 (46.38-2.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.73 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.235 , 0.262 0.225 , 0.263	Depositor DCC
R_{free} test set	3647 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	62.2	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 78822 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13598	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.68% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: N23

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	0.42	0/3255	0.62	0/4414
2	B	0.40	0/1685	0.66	1/2276 (0.0%)
3	C	0.40	0/2686	0.67	1/3647 (0.0%)
4	D	0.43	0/2328	0.61	0/3143
5	E	0.34	0/1449	0.60	1/1954 (0.1%)
6	F	0.44	0/1393	0.66	0/1868
7	G	0.34	0/1034	0.55	0/1390
All	All	0.40	0/13830	0.63	3/18692 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	11	ILE	N-CA-C	-5.82	95.29	111.00
2	B	350	ARG	NE-CZ-NH2	-5.39	117.60	120.30
5	E	17	GLY	N-CA-C	-5.24	100.01	113.10

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	3176	0	3124	99	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1655	0	1659	121	0
3	C	2618	0	2563	82	0
4	D	2279	0	2248	47	0
5	E	1415	0	1416	70	0
6	F	1371	0	1410	36	0
7	G	1022	0	1042	35	0
8	B	20	0	16	1	0
9	A	8	0	0	1	0
9	B	3	0	0	0	0
9	C	7	0	0	0	0
9	D	10	0	0	0	0
9	E	1	0	0	0	0
9	F	12	0	0	0	0
9	G	1	0	0	0	0
All	All	13598	0	13478	468	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (468) 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)
3:C:183:THR:HG22	3:C:185:TRP:H	1.28	0.97
2:B:182:LEU:HD22	2:B:184:ILE:HG12	1.46	0.95
5:E:152:GLN:HB2	5:E:155:LYS:HD2	1.49	0.94
2:B:166:ILE:HD12	2:B:281:LEU:HD22	1.50	0.90
3:C:189:MET:HG2	3:C:195:MET:HE1	1.54	0.90
2:B:205:ASN:HD22	2:B:208:ALA:H	1.14	0.89
1:A:191:LYS:HE2	1:A:303:VAL:HG22	1.52	0.88
1:A:412:PRO:HB2	1:A:414:PHE:CE2	2.09	0.88
2:B:165:HIS:CD2	2:B:181:ARG:HG2	2.09	0.88
6:F:4:THR:HG23	6:F:55:ARG:HE	1.39	0.87
2:B:329:VAL:O	2:B:330:LEU:HD13	1.77	0.84
1:A:55:VAL:HG13	1:A:58:LEU:HD12	1.59	0.84
2:B:165:HIS:HD2	2:B:181:ARG:HG2	1.40	0.84
2:B:330:LEU:HB3	2:B:336:LYS:HD3	1.57	0.83
2:B:327:GLU:HG3	2:B:328:ARG:N	1.94	0.83
1:A:135:ASN:HB3	1:A:397:LYS:NZ	1.94	0.83
2:B:330:LEU:CB	2:B:336:LYS:HD3	2.11	0.80
3:C:31:GLU:HG2	3:C:49:LYS:HG2	1.64	0.80
2:B:330:LEU:HG	2:B:336:LYS:CD	2.13	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:177:HIS:O	2:B:178:LEU:HB2	1.82	0.77
2:B:329:VAL:C	2:B:330:LEU:HD13	2.05	0.77
1:A:216:PRO:HB2	1:A:219:GLN:HB2	1.67	0.76
5:E:97:THR:O	5:E:101:THR:HG23	1.87	0.75
2:B:347:PRO:HB2	2:B:349:ARG:NH2	2.01	0.74
3:C:358:ASP:OD1	3:C:360:ARG:HG2	1.86	0.74
3:C:189:MET:HG2	3:C:195:MET:CE	2.18	0.74
2:B:175:LEU:HD12	2:B:178:LEU:HD12	1.69	0.74
2:B:349:ARG:HH21	2:B:349:ARG:HG2	1.52	0.74
3:C:107:ASN:ND2	3:C:109:LYS:H	1.85	0.74
1:A:19:LEU:HD13	1:A:96:VAL:HG13	1.70	0.73
2:B:330:LEU:HG	2:B:336:LYS:HD3	1.69	0.73
4:D:164:VAL:HG22	4:D:224:THR:HG23	1.71	0.72
5:E:90:GLN:HG2	5:E:94:GLU:OE2	1.90	0.72
2:B:336:LYS:HE2	2:B:337:LEU:N	2.04	0.72
1:A:55:VAL:O	1:A:55:VAL:HG12	1.90	0.71
6:F:145:GLU:O	6:F:149:MET:HG3	1.90	0.71
5:E:102:ASN:HD21	5:E:130:ARG:HE	1.36	0.71
4:D:147:ARG:HB2	4:D:150:GLU:HB2	1.71	0.71
3:C:14:HIS:H	3:C:331:GLN:HE22	1.39	0.70
3:C:282:GLY:HA2	3:C:370:LYS:HE3	1.72	0.70
2:B:282:LEU:HD21	2:B:301:ILE:HD13	1.73	0.70
2:B:325:TYR:CD2	2:B:330:LEU:HD21	2.26	0.69
2:B:330:LEU:CG	2:B:336:LYS:HD3	2.22	0.69
2:B:336:LYS:HA	2:B:339:LYS:NZ	2.07	0.69
2:B:261:ALA:HB3	2:B:262:PRO:HD3	1.74	0.68
3:C:155:VAL:HG21	3:C:180:PRO:HG3	1.74	0.68
3:C:144:THR:O	3:C:161:SER:HB2	1.93	0.68
2:B:326:LEU:HA	2:B:330:LEU:HD23	1.75	0.68
2:B:155:VAL:HG21	2:B:286:ILE:HD11	1.76	0.67
1:A:116:PRO:O	1:A:117:LEU:HB2	1.92	0.67
4:D:263:HIS:HD2	4:D:266:MET:CE	2.06	0.67
2:B:164:THR:HB	2:B:182:LEU:HB3	1.75	0.67
1:A:212:GLU:OE2	1:A:270:ASP:N	2.27	0.67
1:A:246:ASP:OD1	5:E:50:LYS:HE3	1.95	0.67
7:G:87:LYS:N	7:G:87:LYS:HD3	2.10	0.66
2:B:336:LYS:HE2	2:B:336:LYS:C	2.15	0.66
7:G:23:LYS:HG2	7:G:24:PHE:H	1.59	0.66
1:A:129:ILE:O	1:A:133:SER:HB2	1.96	0.66
1:A:55:VAL:CG1	1:A:58:LEU:HD12	2.25	0.66
2:B:198:LEU:HD23	2:B:202:TYR:O	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ASN:HB3	1:A:397:LYS:HZ1	1.59	0.65
1:A:30:ILE:HD13	1:A:375:TYR:CZ	2.31	0.65
2:B:329:VAL:HG13	2:B:330:LEU:CD2	2.27	0.65
5:E:82:LEU:HD13	5:E:95:MET:SD	2.36	0.65
3:C:179:ARG:HH11	3:C:179:ARG:HG3	1.61	0.65
2:B:295:SER:O	2:B:299:LYS:HG2	1.96	0.65
7:G:87:LYS:H	7:G:87:LYS:HD3	1.62	0.65
2:B:287:GLN:NE2	2:B:287:GLN:HA	2.12	0.64
1:A:359:LYS:N	1:A:360:PRO:HD3	2.13	0.64
6:F:92:ARG:HG2	6:F:92:ARG:HH11	1.62	0.64
3:C:72:THR:HA	3:C:98:ALA:HB1	1.80	0.64
1:A:389:GLU:O	1:A:393:VAL:HG13	1.98	0.63
2:B:287:GLN:HE21	2:B:287:GLN:HA	1.62	0.63
3:C:47:GLU:OE2	3:C:49:LYS:HE3	1.98	0.63
3:C:144:THR:H	6:F:28:GLN:NE2	1.96	0.63
4:D:222:TYR:C	4:D:223:ILE:HD12	2.19	0.63
3:C:107:ASN:HD22	3:C:107:ASN:C	2.02	0.63
4:D:7:ASN:HD21	4:D:111:HIS:CE1	2.17	0.63
2:B:205:ASN:ND2	2:B:208:ALA:H	1.94	0.63
1:A:215:ILE:HD11	1:A:269:ILE:HD13	1.81	0.63
5:E:134:GLN:O	5:E:138:GLN:HG3	1.98	0.63
4:D:228:PHE:H	4:D:231:HIS:HD2	1.45	0.63
5:E:86:ASN:O	5:E:87:SER:HB3	1.98	0.62
6:F:4:THR:HG23	6:F:55:ARG:NE	2.12	0.62
4:D:121:PHE:O	4:D:124:VAL:HG12	1.99	0.62
1:A:289:ASN:ND2	1:A:291:ASP:H	1.97	0.62
3:C:370:LYS:O	3:C:371:ILE:HB	1.99	0.62
2:B:178:LEU:O	2:B:285:THR:HG23	2.00	0.62
3:C:17:ASN:ND2	3:C:22:GLN:HG3	2.15	0.62
1:A:38:LYS:HE2	1:A:72:TYR:CZ	2.35	0.62
4:D:68:GLN:HA	4:D:72:ALA:HB3	1.82	0.62
4:D:189:ARG:NH1	4:D:197:GLN:HB2	2.15	0.62
5:E:133:LEU:HB3	5:E:137:ARG:HH12	1.65	0.61
4:D:130:GLN:HA	4:D:130:GLN:OE1	1.99	0.61
3:C:119:VAL:HG22	3:C:120:ILE:N	2.15	0.61
1:A:289:ASN:HD22	1:A:290:PRO:N	1.99	0.61
3:C:76:ALA:HB2	3:C:93:LEU:HD11	1.82	0.61
2:B:163:VAL:HG22	2:B:164:THR:N	2.16	0.61
1:A:397:LYS:HG2	1:A:401:GLU:OE2	2.01	0.61
2:B:287:GLN:HE21	2:B:287:GLN:CA	2.11	0.61
2:B:337:LEU:HG	2:B:338:SER:H	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:351:LYS:HE3	2:B:351:LYS:N	2.16	0.61
4:D:205:PRO:HB3	4:D:222:TYR:CZ	2.36	0.61
1:A:289:ASN:HD22	1:A:289:ASN:C	2.02	0.60
5:E:95:MET:HA	5:E:95:MET:HE3	1.81	0.60
1:A:135:ASN:HB3	1:A:397:LYS:HZ2	1.64	0.60
4:D:135:GLY:O	4:D:137:GLU:HG3	2.01	0.60
3:C:321:LEU:HD11	6:F:129:HIS:CE1	2.36	0.60
4:D:182:MET:HG3	4:D:200:PHE:CD1	2.37	0.60
1:A:223:THR:O	1:A:227:VAL:HG23	2.01	0.60
2:B:334:VAL:HG21	7:G:19:TYR:CE1	2.35	0.60
7:G:10:ARG:O	7:G:11:PHE:HB3	2.02	0.60
5:E:139:GLU:O	5:E:142:LEU:HD23	2.02	0.59
7:G:75:ALA:O	7:G:79:VAL:HG23	2.01	0.59
6:F:120:ASN:O	6:F:124:GLU:HG3	2.03	0.59
3:C:107:ASN:HD22	3:C:108:GLU:N	1.99	0.59
4:D:159:ASP:O	4:D:160:ARG:HB3	2.02	0.59
5:E:60:ILE:CD1	5:E:116:ILE:HG21	2.32	0.59
6:F:2:THR:OG1	6:F:5:LEU:HB3	2.03	0.59
3:C:26:CYS:SG	3:C:55:VAL:HB	2.43	0.59
3:C:213:ASN:ND2	3:C:255:GLU:CD	2.56	0.58
6:F:4:THR:HG23	6:F:55:ARG:HH21	1.69	0.58
7:G:143:ARG:HA	7:G:146:THR:OG1	2.02	0.58
2:B:325:TYR:O	2:B:329:VAL:HG12	2.03	0.58
2:B:219:LYS:HG2	2:B:220:LEU:HD13	1.86	0.58
1:A:393:VAL:HG11	1:A:414:PHE:CD2	2.39	0.58
2:B:175:LEU:N	2:B:175:LEU:HD23	2.19	0.58
1:A:143:VAL:HG13	1:A:146:VAL:HG23	1.86	0.58
1:A:140:TYR:HB2	1:A:394:CYS:SG	2.43	0.58
5:E:15:LEU:HD22	5:E:63:GLU:HG3	1.85	0.58
3:C:17:ASN:HD21	3:C:22:GLN:HG3	1.68	0.57
2:B:173:PHE:CG	2:B:174:SER:N	2.72	0.57
5:E:150:ASP:O	5:E:152:GLN:N	2.38	0.57
5:E:87:SER:HA	5:E:153:ASN:OD1	2.05	0.57
3:C:365:ALA:C	3:C:366:LEU:HD12	2.24	0.57
7:G:118:SER:O	7:G:120:ASN:N	2.38	0.57
5:E:150:ASP:OD1	5:E:151:PRO:HD2	2.05	0.57
3:C:155:VAL:HG21	3:C:180:PRO:CG	2.35	0.57
1:A:329:ARG:O	1:A:330:ASP:HB2	2.05	0.57
2:B:333:ASP:O	2:B:336:LYS:HG3	2.06	0.56
3:C:252:PHE:HA	3:C:258:LEU:HD23	1.88	0.56
1:A:239:VAL:HG13	5:E:4:TYR:CE2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:ILE:O	1:A:263:SER:CB	2.53	0.56
1:A:202:TYR:O	1:A:205:GLN:HB3	2.06	0.56
5:E:152:GLN:CB	5:E:155:LYS:HD2	2.31	0.56
1:A:309:ILE:HA	1:A:312:ARG:NE	2.21	0.56
5:E:144:LEU:HD22	5:E:148:VAL:HG23	1.88	0.56
3:C:359:VAL:O	3:C:363:GLU:HG3	2.06	0.56
6:F:4:THR:CG2	6:F:55:ARG:HE	2.16	0.56
2:B:330:LEU:HG	2:B:336:LYS:CE	2.36	0.55
5:E:82:LEU:HD23	5:E:148:VAL:HG21	1.89	0.55
5:E:95:MET:HA	5:E:95:MET:CE	2.36	0.55
3:C:254:THR:HA	3:C:340:ALA:O	2.07	0.55
1:A:400:TYR:CE1	1:A:405:PRO:HA	2.41	0.55
4:D:263:HIS:HD2	4:D:266:MET:HE2	1.71	0.55
4:D:202:HIS:O	4:D:203:ARG:HB3	2.05	0.55
2:B:177:HIS:O	2:B:178:LEU:CB	2.54	0.54
3:C:144:THR:H	6:F:28:GLN:HE22	1.55	0.54
2:B:158:ASP:HA	2:B:304:SER:O	2.08	0.54
2:B:329:VAL:HG13	2:B:330:LEU:HD22	1.90	0.54
4:D:118:ARG:HD3	4:D:118:ARG:C	2.28	0.54
2:B:160:GLY:O	2:B:185:ALA:HB1	2.07	0.53
2:B:336:LYS:HA	2:B:339:LYS:HZ3	1.73	0.53
5:E:58:TYR:CD1	5:E:168:PHE:HZ	2.27	0.53
1:A:395:HIS:HE1	1:A:411:ASN:OD1	1.90	0.53
1:A:253:ILE:HD12	5:E:159:TRP:CH2	2.43	0.53
2:B:334:VAL:HG13	2:B:334:VAL:O	2.08	0.53
7:G:78:ILE:O	7:G:82:VAL:HG23	2.08	0.53
5:E:124:GLN:O	5:E:128:VAL:HG23	2.08	0.53
1:A:343:VAL:HG22	1:A:363:ILE:HD12	1.91	0.53
2:B:349:ARG:HG2	2:B:349:ARG:NH2	2.24	0.53
3:C:216:ARG:NH1	3:C:255:GLU:O	2.40	0.53
1:A:395:HIS:CE1	1:A:411:ASN:OD1	2.62	0.52
1:A:85:ASP:OD2	1:A:88:LEU:HD22	2.10	0.52
3:C:249:ALA:HB1	3:C:332:ILE:HG22	1.92	0.52
7:G:114:PHE:HZ	7:G:126:LEU:HD23	1.72	0.52
2:B:330:LEU:HD22	2:B:330:LEU:N	2.25	0.52
2:B:334:VAL:HG21	7:G:19:TYR:CZ	2.44	0.52
5:E:50:LYS:NZ	5:E:159:TRP:O	2.42	0.52
1:A:28:GLN:HG2	4:D:10:ILE:HD13	1.90	0.52
1:A:412:PRO:HB2	1:A:414:PHE:HE2	1.65	0.52
1:A:104:GLU:OE1	1:A:106:GLU:HG3	2.10	0.52
3:C:144:THR:N	6:F:28:GLN:NE2	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:35:PHE:N	4:D:35:PHE:CD2	2.78	0.52
5:E:9:MET:SD	5:E:63:GLU:HG2	2.50	0.51
3:C:223:ASP:OD1	3:C:225:THR:HG23	2.10	0.51
4:D:223:ILE:N	4:D:223:ILE:HD12	2.24	0.51
4:D:158:LYS:O	4:D:158:LYS:HD3	2.11	0.51
5:E:85:CYS:SG	5:E:149:PHE:HZ	2.33	0.51
1:A:94:GLU:HB3	9:A:422:HOH:O	2.10	0.51
1:A:143:VAL:HG13	1:A:146:VAL:CG2	2.41	0.51
1:A:343:VAL:HG22	1:A:363:ILE:CD1	2.41	0.51
6:F:158:ARG:O	6:F:162:GLU:HG3	2.10	0.51
5:E:42:VAL:HB	5:E:143:ARG:NH1	2.26	0.51
4:D:133:GLU:O	4:D:133:GLU:HG2	2.11	0.51
5:E:20:ALA:HB3	5:E:22:LEU:HD22	1.93	0.51
1:A:321:LEU:HD13	1:A:326:THR:HB	1.93	0.51
2:B:313:LEU:HB3	2:B:314:PRO:HD3	1.92	0.51
2:B:327:GLU:HG3	2:B:328:ARG:HG3	1.93	0.51
4:D:7:ASN:OD1	4:D:115:MET:HG2	2.11	0.50
1:A:174:VAL:HG12	1:A:175:THR:N	2.26	0.50
6:F:87:CYS:O	6:F:91:MET:HG2	2.11	0.50
2:B:325:TYR:O	2:B:330:LEU:HD22	2.11	0.50
3:C:167:ARG:HG2	3:C:197:GLU:HG3	1.93	0.50
4:D:67:LEU:O	4:D:72:ALA:HB2	2.11	0.50
1:A:91:ARG:O	1:A:94:GLU:HB2	2.12	0.50
4:D:95:LEU:HD11	4:D:116:LEU:HG	1.93	0.50
2:B:351:LYS:CA	2:B:351:LYS:HE3	2.41	0.50
7:G:38:ASP:O	7:G:42:VAL:HG23	2.12	0.50
2:B:286:ILE:HG21	2:B:298:TYR:CE2	2.46	0.50
2:B:347:PRO:CB	2:B:349:ARG:NH2	2.74	0.50
1:A:239:VAL:HG23	1:A:240:LYS:N	2.25	0.50
6:F:8:TYR:O	6:F:12:VAL:HG23	2.12	0.50
2:B:302:VAL:O	2:B:302:VAL:HG13	2.11	0.50
7:G:114:PHE:CZ	7:G:126:LEU:HD23	2.47	0.50
3:C:94:ARG:HH11	3:C:94:ARG:HB3	1.76	0.50
4:D:137:GLU:OE2	4:D:158:LYS:HE2	2.11	0.50
3:C:281:PHE:CZ	3:C:283:GLY:HA2	2.46	0.50
2:B:322:LYS:HD2	7:G:16:VAL:HG11	1.93	0.49
6:F:92:ARG:HG2	6:F:92:ARG:NH1	2.27	0.49
2:B:325:TYR:CD1	2:B:329:VAL:HG11	2.47	0.49
3:C:367:LYS:HD3	3:C:368:ASP:N	2.28	0.49
3:C:183:THR:HG23	3:C:184:PRO:HD2	1.94	0.49
2:B:194:ILE:HG12	2:B:213:VAL:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:20:ASP:OD1	7:G:22:ASN:HB2	2.12	0.49
2:B:322:LYS:HG2	2:B:342:ILE:HD11	1.94	0.49
4:D:251:LEU:O	4:D:255:ILE:HG13	2.13	0.49
1:A:215:ILE:HD11	1:A:269:ILE:CD1	2.42	0.49
1:A:174:VAL:HG13	1:A:193:ILE:O	2.12	0.49
3:C:209:CYS:SG	3:C:251:THR:HA	2.53	0.49
4:D:181:PHE:CE2	6:F:157:ALA:HA	2.48	0.49
4:D:171:ASP:O	4:D:174:ASP:HB2	2.13	0.49
2:B:330:LEU:HG	2:B:336:LYS:HE3	1.95	0.49
1:A:278:GLY:O	1:A:281:ILE:HG12	2.12	0.49
1:A:156:ARG:HH11	1:A:156:ARG:HG2	1.77	0.49
1:A:311:VAL:C	1:A:314:PRO:HD2	2.34	0.49
1:A:55:VAL:CG1	1:A:55:VAL:O	2.61	0.48
6:F:146:ILE:HA	6:F:149:MET:CE	2.43	0.48
2:B:251:ILE:HG22	2:B:252:ILE:N	2.28	0.48
7:G:87:LYS:HE3	7:G:90:ASP:CG	2.34	0.48
3:C:34:ILE:HB	3:C:46:HIS:HB2	1.94	0.48
5:E:74:TYR:O	5:E:77:GLU:HB2	2.14	0.48
6:F:121:PHE:O	6:F:125:GLN:HG2	2.12	0.48
5:E:153:ASN:O	5:E:155:LYS:HG3	2.13	0.48
2:B:340:PHE:CE2	2:B:342:ILE:HD11	2.48	0.48
7:G:151:VAL:HG23	7:G:151:VAL:O	2.14	0.48
3:C:92:ILE:HD12	3:C:92:ILE:N	2.29	0.48
5:E:152:GLN:O	5:E:155:LYS:HB2	2.14	0.48
1:A:369:THR:HA	1:A:373:GLN:OE1	2.14	0.48
4:D:2:ILE:HG21	6:F:163:GLU:HG2	1.96	0.48
7:G:124:VAL:O	7:G:127:GLN:HB2	2.13	0.48
1:A:169:ASP:HA	1:A:322:SER:O	2.14	0.48
6:F:127:TYR:HB3	6:F:129:HIS:CE1	2.49	0.48
2:B:156:VAL:HG22	2:B:302:VAL:CG1	2.43	0.48
3:C:32:VAL:HB	3:C:48:LEU:HB2	1.94	0.48
1:A:389:GLU:HA	1:A:392:GLN:NE2	2.29	0.48
7:G:80:LEU:O	7:G:84:ILE:HD13	2.13	0.47
3:C:240:LEU:HD23	3:C:270:PHE:CE2	2.49	0.47
1:A:398:LYS:O	1:A:402:GLU:HG3	2.14	0.47
5:E:158:LYS:HE3	5:E:159:TRP:CZ2	2.49	0.47
1:A:193:ILE:HG23	1:A:292:PHE:CE2	2.49	0.47
2:B:268:PRO:HB2	2:B:273:VAL:O	2.14	0.47
1:A:274:GLU:N	1:A:274:GLU:OE1	2.37	0.47
1:A:163:LEU:HG	1:A:416:VAL:CG2	2.44	0.47
2:B:276:VAL:HB	2:B:280:GLU:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:60:SER:HB2	4:D:91:TYR:CD2	2.49	0.47
3:C:144:THR:OG1	6:F:28:GLN:NE2	2.48	0.47
5:E:86:ASN:HB3	5:E:154:ASP:OD2	2.15	0.47
1:A:111:LEU:HD23	1:A:111:LEU:C	2.35	0.47
5:E:139:GLU:O	5:E:143:ARG:HB2	2.15	0.47
3:C:253:ILE:HD11	3:C:257:SER:HB3	1.96	0.47
1:A:285:PRO:HG2	1:A:294:GLN:O	2.15	0.47
3:C:14:HIS:H	3:C:331:GLN:NE2	2.08	0.47
1:A:270:ASP:OD2	5:E:158:LYS:NZ	2.47	0.47
2:B:299:LYS:O	2:B:300:HIS:ND1	2.48	0.47
3:C:254:THR:HG1	3:C:257:SER:HB2	1.79	0.47
2:B:291:ILE:HG22	2:B:291:ILE:O	2.15	0.47
3:C:107:ASN:C	3:C:107:ASN:ND2	2.65	0.47
6:F:125:GLN:HE21	6:F:125:GLN:HB3	1.44	0.47
3:C:165:LYS:NZ	3:C:199:SER:HA	2.30	0.47
2:B:239:VAL:HG23	2:B:240:LEU:HD13	1.97	0.47
1:A:194:PRO:C	1:A:195:ILE:HD12	2.35	0.47
3:C:358:ASP:O	3:C:362:LEU:HG	2.15	0.46
2:B:347:PRO:O	2:B:350:ARG:HG3	2.15	0.46
3:C:179:ARG:HG3	3:C:179:ARG:NH1	2.29	0.46
3:C:178:GLU:O	3:C:179:ARG:C	2.54	0.46
7:G:38:ASP:HB3	7:G:41:GLU:HB3	1.98	0.46
2:B:323:GLN:NE2	7:G:12:ARG:HA	2.31	0.46
2:B:349:ARG:HH21	2:B:349:ARG:CG	2.24	0.46
3:C:145:VAL:HA	3:C:161:SER:HB3	1.98	0.46
2:B:272:ASN:O	2:B:273:VAL:HG23	2.15	0.46
6:F:4:THR:HG23	6:F:55:ARG:NH2	2.31	0.46
2:B:347:PRO:HG2	2:B:350:ARG:CG	2.45	0.46
2:B:182:LEU:CD2	2:B:184:ILE:HG12	2.30	0.46
2:B:336:LYS:HE2	2:B:337:LEU:CA	2.45	0.46
2:B:168:PRO:HD2	2:B:179:THR:OG1	2.16	0.46
4:D:188:GLY:HA3	6:F:165:LEU:HD23	1.97	0.46
1:A:36:ALA:HB1	1:A:72:TYR:HB3	1.97	0.46
1:A:262:ILE:O	1:A:263:SER:HB3	2.16	0.46
1:A:389:GLU:CD	1:A:414:PHE:HB3	2.36	0.46
2:B:337:LEU:O	2:B:339:LYS:N	2.43	0.46
4:D:233:ASN:OD1	4:D:235:SER:HB3	2.16	0.46
7:G:60:LYS:HG2	7:G:61:ASN:ND2	2.31	0.46
2:B:163:VAL:CG2	2:B:164:THR:N	2.79	0.46
1:A:259:ILE:N	1:A:259:ILE:HD12	2.31	0.45
1:A:176:HIS:HD2	1:A:192:HIS:CD2	2.33	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:ASP:N	1:A:310:ASP:OD1	2.41	0.45
2:B:325:TYR:HA	2:B:329:VAL:HG12	1.98	0.45
1:A:38:LYS:HE2	1:A:72:TYR:CE2	2.52	0.45
5:E:67:THR:O	5:E:71:ILE:HG13	2.15	0.45
7:G:51:MET:HB3	7:G:86:PHE:CE1	2.52	0.45
5:E:169:MET:O	5:E:171:LYS:HG2	2.16	0.45
5:E:153:ASN:OD1	5:E:154:ASP:N	2.48	0.45
2:B:326:LEU:HD23	2:B:326:LEU:C	2.37	0.45
7:G:116:SER:N	7:G:117:PRO:HD3	2.32	0.45
4:D:75:LEU:C	4:D:75:LEU:HD23	2.37	0.45
3:C:131:TRP:O	3:C:131:TRP:HE3	1.99	0.45
7:G:87:LYS:HE3	7:G:90:ASP:CB	2.47	0.45
7:G:62:PRO:HA	7:G:63:PRO:HD3	1.77	0.45
1:A:69:LYS:HB3	1:A:72:TYR:CD1	2.51	0.45
3:C:16:TRP:CE2	3:C:335:LEU:HD21	2.51	0.45
3:C:3:TYR:HB2	3:C:324:LEU:HG	1.99	0.45
2:B:350:ARG:O	2:B:351:LYS:C	2.56	0.45
1:A:211:ARG:NH1	5:E:159:TRP:CZ3	2.85	0.45
3:C:119:VAL:CG2	3:C:120:ILE:N	2.80	0.45
4:D:164:VAL:CG2	4:D:224:THR:HG23	2.44	0.45
3:C:151:HIS:CB	3:C:156:LEU:HB2	2.47	0.44
1:A:128:GLU:O	1:A:132:GLU:HB2	2.16	0.44
2:B:337:LEU:C	2:B:339:LYS:H	2.19	0.44
2:B:280:GLU:HA	2:B:324:LEU:HD11	1.98	0.44
1:A:343:VAL:CG2	1:A:363:ILE:HD12	2.48	0.44
2:B:205:ASN:ND2	2:B:207:SER:H	2.14	0.44
2:B:326:LEU:CD2	2:B:332:GLY:HA2	2.48	0.44
2:B:169:VAL:HG13	2:B:173:PHE:C	2.37	0.44
5:E:87:SER:OG	5:E:90:GLN:CB	2.66	0.44
2:B:337:LEU:HD12	2:B:340:PHE:HB3	2.00	0.44
2:B:178:LEU:HD22	2:B:289:ALA:HB2	1.98	0.44
6:F:149:MET:O	6:F:153:VAL:HG23	2.18	0.44
5:E:78:CYS:HA	5:E:95:MET:HE1	2.00	0.44
5:E:74:TYR:CE1	5:E:137:ARG:HD2	2.53	0.44
1:A:369:THR:HB	1:A:373:GLN:OE1	2.16	0.44
1:A:176:HIS:CD2	1:A:192:HIS:CD2	3.06	0.44
2:B:180:ARG:HH11	2:B:180:ARG:HG3	1.83	0.44
5:E:83:GLN:NE2	5:E:164:VAL:HG11	2.33	0.44
1:A:27:PRO:HG3	1:A:378:TRP:CD2	2.53	0.44
2:B:329:VAL:CG1	2:B:330:LEU:HD22	2.48	0.43
2:B:337:LEU:CG	2:B:338:SER:H	2.27	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:197:GLN:NE2	4:D:199:LEU:HD11	2.33	0.43
5:E:16:ILE:O	5:E:16:ILE:HG23	2.18	0.43
3:C:90:LEU:CD2	3:C:91:VAL:H	2.31	0.43
2:B:351:LYS:HA	2:B:351:LYS:HE3	1.99	0.43
3:C:257:SER:OG	3:C:372:VAL:N	2.47	0.43
1:A:180:VAL:HG22	1:A:185:VAL:HG22	2.01	0.43
5:E:150:ASP:C	5:E:152:GLN:N	2.72	0.43
2:B:163:VAL:HG22	2:B:164:THR:H	1.82	0.43
5:E:158:LYS:HE3	5:E:159:TRP:CE2	2.54	0.43
5:E:74:TYR:OH	5:E:98:LEU:HD12	2.18	0.43
3:C:90:LEU:HD22	3:C:91:VAL:N	2.33	0.43
4:D:103:PRO:HG3	4:D:109:ILE:HD13	1.99	0.43
2:B:182:LEU:HD12	2:B:278:VAL:HB	1.99	0.43
2:B:227:ILE:HD11	2:B:263:GLU:OE2	2.17	0.43
4:D:45:ILE:HG12	4:D:57:VAL:HG22	2.01	0.43
5:E:140:THR:O	5:E:144:LEU:HB2	2.19	0.43
1:A:71:THR:OG1	1:A:71:THR:O	2.34	0.43
5:E:41:ILE:O	5:E:41:ILE:HG13	2.19	0.43
2:B:347:PRO:O	2:B:350:ARG:CD	2.66	0.43
4:D:53:THR:O	4:D:54:LYS:HD2	2.19	0.43
5:E:154:ASP:CG	5:E:154:ASP:O	2.57	0.43
7:G:42:VAL:HG21	7:G:78:ILE:HG21	2.01	0.43
5:E:128:VAL:O	5:E:131:ALA:HB3	2.19	0.43
4:D:131:PHE:CD1	4:D:139:GLU:HG3	2.54	0.43
6:F:82:ILE:O	6:F:85:ILE:HG13	2.19	0.43
1:A:168:ILE:CD1	1:A:335:LEU:HD11	2.48	0.43
2:B:254:VAL:HG13	2:B:257:GLU:HG3	2.00	0.43
1:A:395:HIS:HB3	1:A:407:ILE:HD12	2.00	0.42
1:A:195:ILE:HG22	1:A:282:PHE:CE1	2.54	0.42
4:D:53:THR:C	4:D:54:LYS:HD2	2.39	0.42
5:E:119:LYS:HG3	5:E:120:PRO:HD2	2.00	0.42
7:G:68:SER:OG	7:G:71:VAL:HG12	2.19	0.42
3:C:126:GLU:C	3:C:128:GLU:H	2.22	0.42
2:B:329:VAL:HG13	2:B:330:LEU:CD1	2.49	0.42
3:C:151:HIS:CG	3:C:152:PRO:HD2	2.54	0.42
4:D:15:LEU:HA	4:D:15:LEU:HD23	1.79	0.42
1:A:9:VAL:HG21	1:A:382:SER:HA	2.02	0.42
5:E:139:GLU:HA	5:E:139:GLU:OE1	2.20	0.42
5:E:10:ASP:C	5:E:12:ASP:H	2.23	0.42
1:A:4:ARG:HG3	1:A:4:ARG:HH11	1.84	0.42
1:A:191:LYS:HE2	1:A:303:VAL:CG2	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:ASN:ND2	1:A:289:ASN:C	2.71	0.42
1:A:343:VAL:CG1	1:A:363:ILE:HG13	2.49	0.42
7:G:115:GLU:C	7:G:117:PRO:HD3	2.39	0.42
3:C:90:LEU:CD2	3:C:91:VAL:N	2.82	0.42
3:C:9:GLU:HB3	3:C:10:PRO:CD	2.50	0.42
3:C:321:LEU:CD1	6:F:129:HIS:CE1	3.03	0.42
3:C:90:LEU:HD23	3:C:91:VAL:H	1.83	0.42
1:A:53:LYS:O	1:A:56:ASP:OD2	2.38	0.42
5:E:154:ASP:OD1	5:E:154:ASP:O	2.38	0.42
2:B:337:LEU:O	2:B:338:SER:HB2	2.20	0.42
7:G:79:VAL:O	7:G:83:LEU:HG	2.19	0.42
5:E:60:ILE:HD11	5:E:116:ILE:HG21	2.00	0.42
5:E:81:LYS:O	5:E:83:GLN:N	2.52	0.42
4:D:7:ASN:OD1	4:D:115:MET:CG	2.68	0.42
2:B:169:VAL:HG13	2:B:173:PHE:O	2.19	0.42
2:B:277:GLY:O	2:B:278:VAL:C	2.58	0.42
3:C:226:VAL:HG13	3:C:240:LEU:HB3	2.01	0.42
3:C:72:THR:HA	3:C:98:ALA:CB	2.47	0.41
5:E:137:ARG:HG3	5:E:137:ARG:HH11	1.83	0.41
3:C:143:SER:OG	3:C:162:CYS:HB2	2.19	0.41
2:B:323:GLN:HG2	7:G:16:VAL:HG23	2.02	0.41
2:B:298:TYR:CD1	2:B:340:PHE:HE1	2.37	0.41
1:A:141:ILE:HB	1:A:411:ASN:ND2	2.34	0.41
1:A:60:PHE:CE1	1:A:95:GLN:HG3	2.55	0.41
5:E:106:PRO:HG3	5:E:111:PHE:CE2	2.55	0.41
1:A:359:LYS:O	1:A:359:LYS:CG	2.68	0.41
1:A:400:TYR:HE1	1:A:405:PRO:HA	1.83	0.41
3:C:367:LYS:HD3	3:C:368:ASP:H	1.86	0.41
2:B:248:ASP:OD1	2:B:250:ARG:HG3	2.20	0.41
5:E:87:SER:OG	5:E:90:GLN:HB2	2.20	0.41
5:E:133:LEU:HB3	5:E:137:ARG:NH1	2.33	0.41
1:A:88:LEU:HA	1:A:88:LEU:HD12	1.84	0.41
3:C:92:ILE:CD1	3:C:92:ILE:N	2.83	0.41
1:A:163:LEU:HA	1:A:163:LEU:HD23	1.91	0.41
5:E:89:SER:O	5:E:92:GLU:HB2	2.19	0.41
2:B:318:GLU:HG3	2:B:344:ILE:HD12	2.02	0.41
2:B:350:ARG:O	2:B:352:HIS:N	2.53	0.41
6:F:114:ILE:HG13	6:F:115:SER:N	2.36	0.41
5:E:45:ALA:HA	5:E:68:LEU:HD11	2.03	0.41
5:E:144:LEU:O	5:E:148:VAL:HG23	2.20	0.41
3:C:129:ASN:HB2	3:C:131:TRP:CZ3	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:129:ASN:ND2	3:C:131:TRP:CH2	2.89	0.41
5:E:150:ASP:C	5:E:152:GLN:H	2.24	0.41
4:D:146:TYR:CD1	4:D:147:ARG:HG3	2.56	0.41
7:G:18:GLU:HA	7:G:18:GLU:OE1	2.20	0.41
6:F:85:ILE:HD12	6:F:85:ILE:C	2.41	0.41
6:F:20:LEU:HD12	6:F:132:VAL:HG22	2.02	0.41
7:G:18:GLU:OE1	7:G:23:LYS:NZ	2.53	0.41
2:B:198:LEU:HD21	2:B:204:PHE:O	2.21	0.41
6:F:2:THR:O	6:F:2:THR:OG1	2.38	0.41
3:C:219:TRP:CE2	3:C:227:CYS:HB2	2.56	0.41
6:F:13:ARG:NH1	6:F:133:ASP:OD1	2.41	0.41
5:E:48:TYR:O	5:E:49:PHE:C	2.59	0.41
1:A:284:HIS:HB3	1:A:287:PHE:CE1	2.56	0.41
6:F:146:ILE:HA	6:F:149:MET:HE2	2.02	0.40
3:C:119:VAL:HG21	3:C:136:HIS:HB3	2.02	0.40
1:A:156:ARG:NH1	1:A:156:ARG:HG2	2.36	0.40
2:B:250:ARG:HB2	8:B:395:N23:HAL	2.03	0.40
6:F:74:ILE:HD13	6:F:139:MET:HG2	2.02	0.40
4:D:208:GLU:OE2	4:D:208:GLU:N	2.44	0.40
5:E:88:LYS:HG2	5:E:88:LYS:O	2.20	0.40
2:B:329:VAL:CG1	2:B:330:LEU:CD2	2.99	0.40
2:B:305:GLY:O	2:B:350:ARG:NH2	2.54	0.40
4:D:134:GLU:OE2	4:D:136:LYS:HE2	2.21	0.40
2:B:279:ALA:HB1	2:B:320:GLU:HB3	2.03	0.40
2:B:334:VAL:HG21	7:G:19:TYR:CD1	2.56	0.40
2:B:324:LEU:HD21	7:G:11:PHE:CE2	2.56	0.40
5:E:69:ILE:HG23	5:E:169:MET:HE1	2.02	0.40
6:F:58:LYS:HA	6:F:58:LYS:HE2	2.03	0.40
3:C:228:LEU:HD23	3:C:229:ALA:N	2.36	0.40
2:B:280:GLU:HA	2:B:324:LEU:CD1	2.51	0.40
3:C:165:LYS:HZ2	3:C:199:SER:HA	1.86	0.40
2:B:317:LEU:HB3	2:B:344:ILE:CD1	2.51	0.40
3:C:264:ASP:HB3	3:C:266:PHE:CE2	2.56	0.40
3:C:37:LYS:HD2	3:C:42:TRP:CZ2	2.56	0.40
3:C:135:LYS:HE3	3:C:135:LYS:HB3	1.98	0.40
2:B:298:TYR:O	2:B:342:ILE:HA	2.22	0.40
1:A:69:LYS:HB3	1:A:72:TYR:HD1	1.86	0.40
4:D:45:ILE:HA	4:D:56:MET:O	2.21	0.40
5:E:32:ALA:CB	5:E:135:GLN:OE1	2.69	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	387/418 (93%)	351 (91%)	32 (8%)	4 (1%)	19	45
2	B	207/394 (52%)	173 (84%)	24 (12%)	10 (5%)	3	5
3	C	333/372 (90%)	308 (92%)	22 (7%)	3 (1%)	21	49
4	D	280/300 (93%)	260 (93%)	17 (6%)	3 (1%)	17	42
5	E	172/178 (97%)	154 (90%)	13 (8%)	5 (3%)	6	14
6	F	165/168 (98%)	156 (94%)	8 (5%)	1 (1%)	30	59
7	G	130/151 (86%)	121 (93%)	5 (4%)	4 (3%)	5	12
All	All	1674/1981 (84%)	1523 (91%)	121 (7%)	30 (2%)	11	27

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	171	GLU
2	B	334	VAL
2	B	350	ARG
6	F	102	PHE
7	G	11	PHE
7	G	119	ASP
1	A	263	SER
2	B	278	VAL
2	B	351	LYS
5	E	82	LEU
5	E	87	SER
5	E	153	ASN
7	G	22	ASN
3	C	371	ILE
7	G	23	LYS
2	B	178	LEU
4	D	207	LEU
4	D	237	ARG

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Mol	Chain	Res	Type
5	E	151	PRO
2	B	272	ASN
2	B	337	LEU
3	C	176	VAL
4	D	213	ASP
1	A	70	PRO
2	B	186	GLY
5	E	49	PHE
2	B	348	PRO
1	A	216	PRO
1	A	194	PRO
3	C	179	ARG

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	346/363 (95%)	328 (95%)	18 (5%)	29	58
2	B	174/345 (50%)	156 (90%)	18 (10%)	9	20
3	C	283/313 (90%)	269 (95%)	14 (5%)	31	61
4	D	248/264 (94%)	243 (98%)	5 (2%)	63	87
5	E	156/159 (98%)	145 (93%)	11 (7%)	18	41
6	F	154/155 (99%)	149 (97%)	5 (3%)	46	77
7	G	110/123 (89%)	99 (90%)	11 (10%)	9	22
All	All	1471/1722 (85%)	1389 (94%)	82 (6%)	26	54

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	68	GLU
1	A	70	PRO
1	A	71	THR
1	A	82	ILE

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Mol	Chain	Res	Type
1	A	88	LEU
1	A	191	LYS
1	A	230	ARG
1	A	243	ASN
1	A	251	LYS
1	A	255	GLN
1	A	289	ASN
1	A	310	ASP
1	A	335	LEU
1	A	343	VAL
1	A	349	LEU
1	A	353	LEU
1	A	363	ILE
2	B	175	LEU
2	B	182	LEU
2	B	183	ASP
2	B	220	LEU
2	B	235	LEU
2	B	250	ARG
2	B	257	GLU
2	B	274	GLU
2	B	276	VAL
2	B	287	GLN
2	B	290	ASP
2	B	330	LEU
2	B	336	LYS
2	B	342	ILE
2	B	345	GLU
2	B	349	ARG
2	B	351	LYS
2	B	352	HIS
3	C	8	VAL
3	C	30	HIS
3	C	33	HIS
3	C	54	GLN
3	C	90	LEU
3	C	107	ASN
3	C	131	TRP
3	C	140	PRO
3	C	179	ARG
3	C	182	PRO
3	C	210	PHE

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Mol	Chain	Res	Type
3	C	284	ARG
3	C	321	LEU
3	C	367	LYS
4	D	35	PHE
4	D	37	ASP
4	D	84	LEU
4	D	116	LEU
4	D	265	ARG
5	E	10	ASP
5	E	22	LEU
5	E	80	LYS
5	E	94	GLU
5	E	95	MET
5	E	142	LEU
5	E	143	ARG
5	E	144	LEU
5	E	146	GLU
5	E	151	PRO
5	E	161	THR
6	F	2	THR
6	F	101	PHE
6	F	114	ILE
6	F	125	GLN
6	F	165	LEU
7	G	18	GLU
7	G	21	GLU
7	G	22	ASN
7	G	39	GLU
7	G	64	ILE
7	G	69	GLN
7	G	73	ASP
7	G	86	PHE
7	G	87	LYS
7	G	112	LYS
7	G	146	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	A	144	GLN
1	A	176	HIS

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Mol	Chain	Res	Type
1	A	192	HIS
1	A	205	GLN
1	A	243	ASN
1	A	255	GLN
1	A	289	ASN
1	A	306	ASN
1	A	318	ASN
1	A	392	GLN
1	A	395	HIS
2	B	165	HIS
2	B	205	ASN
2	B	284	ASN
2	B	287	GLN
2	B	323	GLN
3	C	46	HIS
3	C	65	ASN
3	C	107	ASN
3	C	331	GLN
4	D	111	HIS
4	D	140	ASN
4	D	202	HIS
4	D	231	HIS
4	D	263	HIS
5	E	83	GLN
5	E	102	ASN
5	E	167	GLN
6	F	28	GLN
6	F	120	ASN
6	F	125	GLN
7	G	61	ASN
7	G	96	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	$\# Z > 2$
8	N23	B	395	-	19,22,22	1.04	0	17,30,30	1.75	2 (11%)

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
8	N23	B	395	-	-	0/8/10/10	0/3/3/3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	395	N23	CAL-CAJ-SAH	-6.07	106.51	113.23
8	B	395	N23	CAO-CAB-CAC	-2.23	124.57	129.68

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	395	N23	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	397/418 (94%)	0.22	13 (3%) 50 50	37, 64, 110, 128	0
2	B	211/394 (53%)	0.52	14 (6%) 22 20	43, 81, 125, 140	0
3	C	339/372 (91%)	0.20	11 (3%) 51 51	42, 63, 94, 125	0
4	D	282/300 (94%)	0.00	2 (0%) 89 90	38, 62, 89, 106	0
5	E	174/178 (97%)	0.47	16 (9%) 11 9	60, 85, 126, 132	0
6	F	167/168 (99%)	0.08	1 (0%) 90 91	42, 55, 75, 103	0
7	G	134/151 (88%)	0.59	15 (11%) 7 5	52, 94, 114, 118	0
All	All	1704/1981 (86%)	0.26	72 (4%) 40 39	37, 67, 114, 140	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	352	HIS	8.9
3	C	372	VAL	6.1
1	A	51	VAL	6.1
2	B	173	PHE	6.0
7	G	151	VAL	5.2
1	A	265	LYS	5.0
2	B	178	LEU	4.9
5	E	85	CYS	4.7
1	A	414	PHE	4.5
6	F	2	THR	4.3
4	D	211	ASP	3.9
2	B	334	VAL	3.8
3	C	319	ALA	3.8
1	A	416	VAL	3.7
1	A	359	LYS	3.7
1	A	40	SER	3.4
2	B	181	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
3	C	367	LYS	3.3
5	E	154	ASP	3.3
5	E	175	GLY	3.2
7	G	63	PRO	3.2
1	A	351	GLU	3.2
5	E	100	ILE	3.1
2	B	349	ARG	3.1
7	G	13	LYS	3.1
2	B	294	ARG	3.0
5	E	156	PRO	3.0
7	G	64	ILE	3.0
5	E	84	LYS	3.0
1	A	39	GLU	2.9
2	B	180	ARG	2.9
5	E	117	TYR	2.9
5	E	151	PRO	2.9
3	C	368	ASP	2.8
5	E	153	ASN	2.8
5	E	96	TYR	2.8
7	G	68	SER	2.8
7	G	58	ALA	2.8
7	G	66	THR	2.7
5	E	82	LEU	2.7
7	G	67	LYS	2.7
2	B	171	GLU	2.7
2	B	206	HIS	2.7
3	C	366	LEU	2.6
2	B	174	SER	2.6
5	E	150	ASP	2.5
1	A	52	MET	2.5
7	G	91	ILE	2.5
5	E	89	SER	2.5
3	C	369	LEU	2.5
7	G	10	ARG	2.5
7	G	50	ASN	2.3
4	D	158	LYS	2.3
5	E	88	LYS	2.3
7	G	19	TYR	2.3
5	E	149	PHE	2.2
2	B	288	ALA	2.2
3	C	370	LYS	2.2
7	G	88	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
3	C	84	ARG	2.2
3	C	127	GLN	2.2
2	B	145	THR	2.2
2	B	187	ARG	2.1
3	C	202	CYS	2.1
1	A	348	LYS	2.1
1	A	161	ARG	2.1
3	C	320	GLY	2.1
7	G	42	VAL	2.1
1	A	2	ALA	2.1
7	G	37	PRO	2.0
1	A	350	SER	2.0
5	E	11	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
8	N23	B	395	20/20	0.94	0.20	1.76	69,71,84,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.