



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:25 PM GMT

PDB ID : 1FKP
Title : CRYSTAL STRUCTURE OF NNRTI RESISTANT K103N MUTANT HIV-1
REVERSE TRANSCRIPTASE IN COMPLEX WITH NEVIRAPINE
Authors : Ren, J.; Milton, J.; Weaver, K.L.; Short, S.A.; Stuart, D.I.; Stammers, D.K.
Deposited on : 2000-08-10
Resolution : 2.90 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

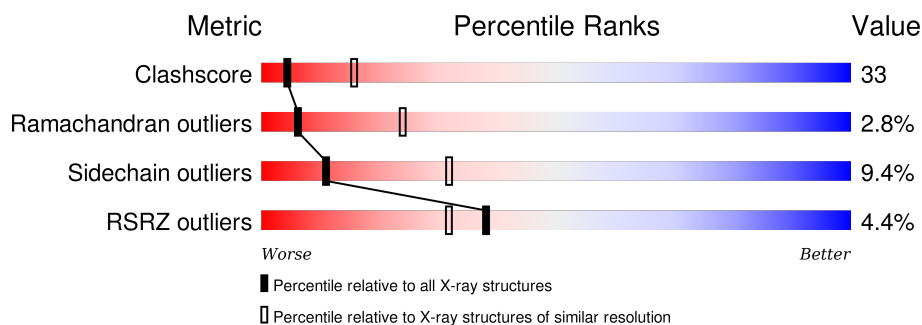
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	543	<div> <div>5%</div> <div>40%</div> <div>51%</div> <div>8%</div> </div>
2	B	440	<div> <div>4%</div> <div>44%</div> <div>43%</div> <div>6%</div> <div>7%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7814 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 HIV-1 RT, A-CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	542	Total	C	N	O	S	0	0	0
			4430	2865	738	819	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ASN	LEU	MUTATION	UNP P04585

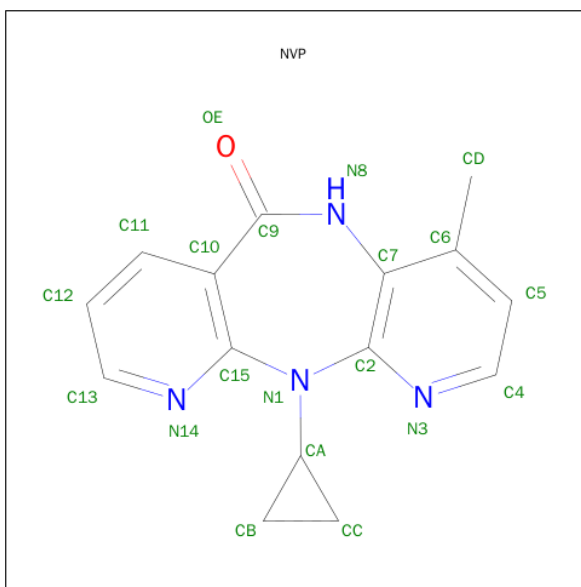
- Molecule 2 is a protein called HIV-1 RT, B-CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	408	Total	C	N	O	S	0	0	0
			3364	2183	560	614	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	103	ASN	LEU	MUTATION	UNP P04585

- Molecule 3 is 11-CYCLOPROPYL-5,11-DIHYDRO-4-METHYL-6H-DIPYRIDO[3,2-B:2',3'-E][1,4]DIAZEPIN-6-ONE (three-letter code: NVP) (formula: C₁₅H₁₄N₄O).

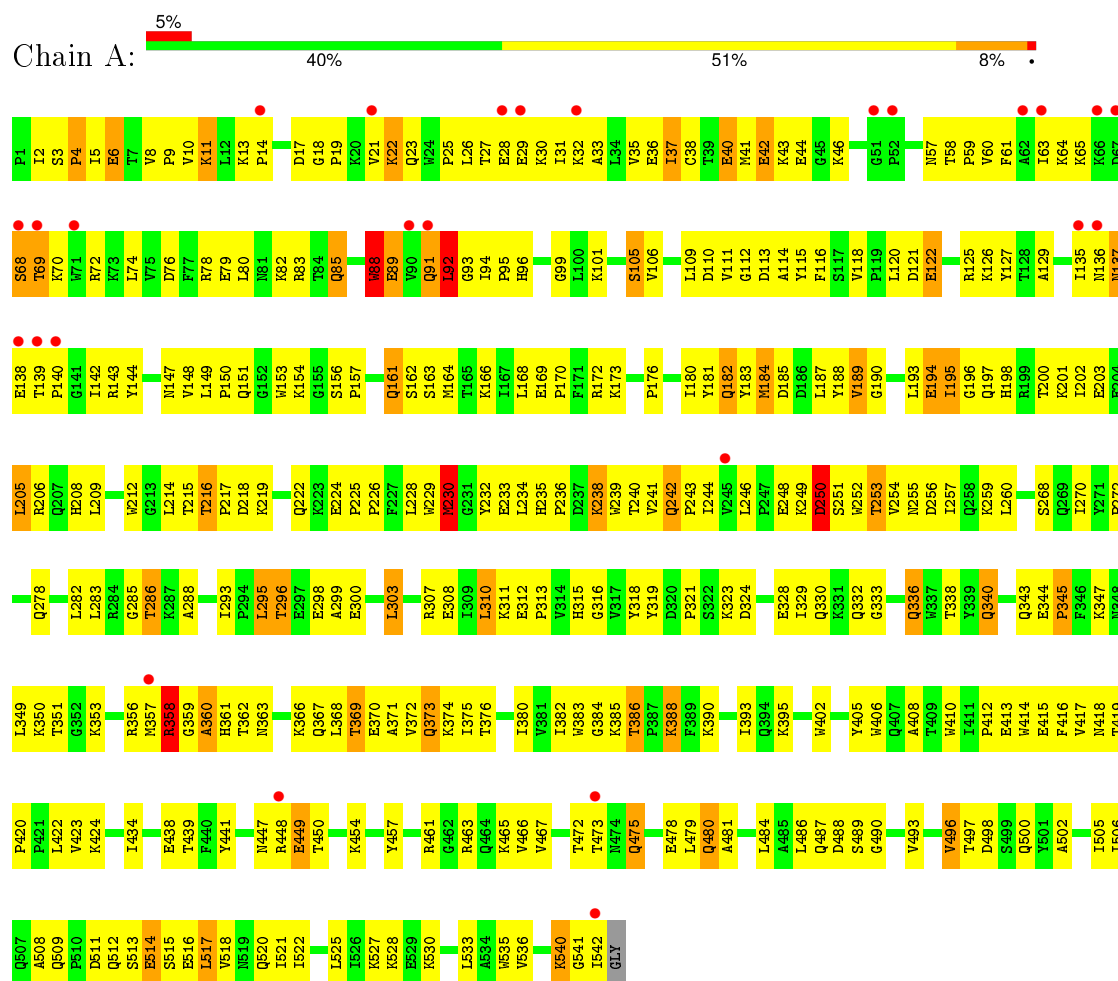


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			20	15	4	1		

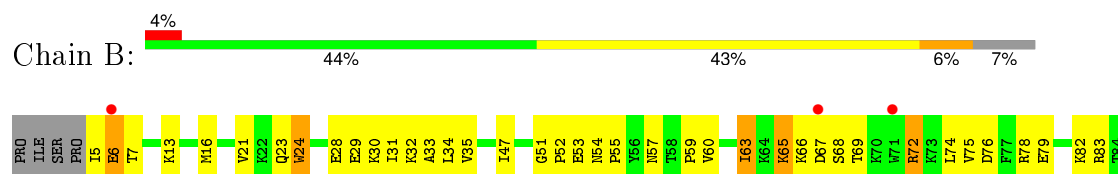
3 Residue-property plots

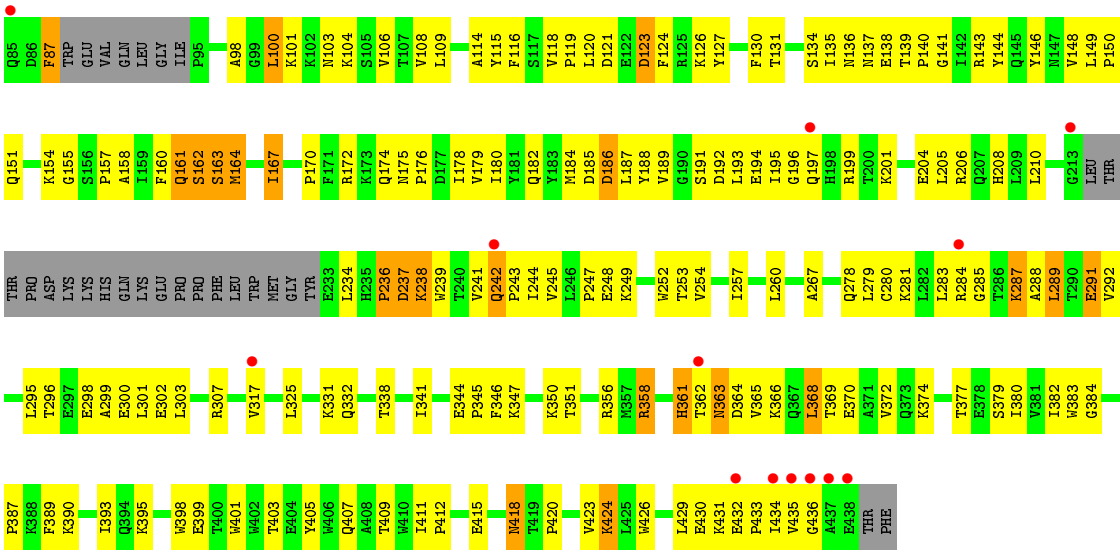
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: HIV-1 RT, A-CHAIN



• Molecule 2: HIV-1 RT, B-CHAIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	139.70 Å 109.80 Å 72.90 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 29.67 – 2.88	Depositor EDS
% Data completeness (in resolution range)	94.9 (30.00-2.90) 92.3 (29.67-2.88)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.90 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.215 , 0.281 0.195 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.735	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 91.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 24022 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7814	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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: CSD, NVP

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	0.50	0/4539	0.73	1/6170 (0.0%)
2	B	0.49	0/3456	0.73	0/4692
All	All	0.50	0/7995	0.73	1/10862 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	388	LYS	N-CA-C	-5.18	97.00	111.00

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	4430	0	4474	333	0
2	B	3364	0	3393	199	0
3	A	20	0	14	2	0
All	All	7814	0	7881	514	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 33.

All (514) 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:239:TRP:CZ2	1:A:316:GLY:HA3	2.00	0.96
1:A:65:LYS:HB3	1:A:72:ARG:HD3	1.46	0.96
1:A:358:ARG:HB3	1:A:358:ARG:NH1	1.81	0.96
1:A:125:ARG:HD3	1:A:147:ASN:HA	1.46	0.93
2:B:395:LYS:O	2:B:399:GLU:HG2	1.70	0.92
1:A:23:GLN:HE22	1:A:60:VAL:H	1.05	0.90
1:A:239:TRP:CE2	1:A:316:GLY:HA3	2.05	0.90
1:A:138:GLU:HG3	1:A:139:THR:H	1.37	0.88
1:A:61:PHE:HB2	1:A:74:LEU:HD21	1.55	0.88
2:B:356:ARG:NH2	2:B:361:HIS:ND1	2.22	0.88
2:B:7:THR:HG22	2:B:119:PRO:HG2	1.56	0.87
2:B:242:GLN:HE21	2:B:242:GLN:HA	1.38	0.86
1:A:136:ASN:OD1	1:A:138:GLU:HB3	1.80	0.82
1:A:61:PHE:HB2	1:A:74:LEU:CD2	2.09	0.82
1:A:448:ARG:HG3	1:A:449:GLU:HG2	1.59	0.82
1:A:59:PRO:HG2	1:A:76:ASP:HB3	1.60	0.81
2:B:420:PRO:HB2	2:B:423:VAL:HG13	1.63	0.80
1:A:17:ASP:O	1:A:83:ARG:HD3	1.82	0.80
1:A:125:ARG:NH1	1:A:147:ASN:HB3	1.97	0.79
1:A:79:GLU:HG3	1:A:83:ARG:NH1	1.98	0.79
1:A:540:LYS:HB3	1:A:542:ILE:HD12	1.66	0.78
1:A:129:ALA:HB1	1:A:143:ARG:NH2	2.00	0.77
2:B:157:PRO:HG2	2:B:184:MET:HA	1.66	0.77
1:A:161:GLN:HE21	1:A:161:GLN:C	1.89	0.76
2:B:158:ALA:O	2:B:161:GLN:HB2	1.85	0.76
1:A:502:ALA:O	1:A:506:ILE:HD12	1.87	0.75
1:A:278:GLN:HG2	1:A:298:GLU:HB3	1.66	0.75
1:A:26:LEU:HD22	1:A:30:LYS:HE2	1.69	0.75
2:B:247:PRO:HB2	2:B:249:LYS:NZ	2.03	0.74
1:A:122:GLU:CD	1:A:122:GLU:H	1.90	0.72
2:B:245:VAL:HG23	2:B:431:LYS:HB2	1.73	0.71
1:A:193:LEU:HB3	1:A:197:GLN:HE21	1.55	0.71
2:B:13:LYS:HE2	2:B:82:LYS:O	1.91	0.71
1:A:358:ARG:HH11	1:A:358:ARG:HB3	1.56	0.71
1:A:6:GLU:H	1:A:6:GLU:CD	1.94	0.71
1:A:244:ILE:CG2	1:A:310:LEU:HD13	2.21	0.71
1:A:206:ARG:NH2	1:A:218:ASP:HB3	2.06	0.70
1:A:180:ILE:HG12	1:A:189:VAL:HG13	1.74	0.70
2:B:100:LEU:HD23	2:B:100:LEU:H	1.57	0.70
2:B:182:GLN:HB2	2:B:187:LEU:HD12	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:GLU:HA	1:A:307:ARG:NH2	2.07	0.69
1:A:366:LYS:O	1:A:369:THR:HB	1.93	0.69
2:B:387:PRO:HG2	2:B:389:PHE:CE1	2.27	0.69
1:A:506:ILE:HG13	1:A:533:LEU:HD23	1.75	0.69
2:B:65:LYS:HZ3	2:B:72:ARG:HD3	1.59	0.68
1:A:234:LEU:N	1:A:234:LEU:HD12	2.08	0.68
1:A:126:LYS:HE3	1:A:127:TYR:CZ	2.29	0.68
1:A:270:ILE:O	1:A:272:PRO:HD3	1.94	0.67
2:B:170:PRO:O	2:B:174:GLN:HG2	1.94	0.67
2:B:197:GLN:O	2:B:201:LYS:HB2	1.95	0.67
2:B:5:ILE:CG1	2:B:6:GLU:N	2.58	0.67
1:A:410:TRP:CD1	2:B:363:ASN:HB2	2.29	0.67
2:B:155:GLY:O	2:B:158:ALA:HB3	1.95	0.67
1:A:497:THR:O	1:A:535:TRP:HA	1.95	0.67
1:A:235:HIS:ND1	1:A:238:LYS:HE3	2.09	0.67
1:A:139:THR:CG2	1:A:140:PRO:HD2	2.25	0.66
1:A:31:ILE:O	1:A:35:VAL:HG23	1.95	0.66
2:B:238:LYS:O	2:B:239:TRP:HD1	1.78	0.66
1:A:296:THR:HG23	1:A:299:ALA:HB2	1.77	0.65
1:A:358:ARG:HH22	1:A:512:GLN:HB2	1.60	0.65
1:A:358:ARG:CZ	1:A:358:ARG:HB3	2.25	0.65
2:B:237:ASP:O	2:B:239:TRP:N	2.28	0.65
1:A:393:ILE:HB	1:A:423:VAL:HG22	1.78	0.65
1:A:88:TRP:CE3	1:A:88:TRP:HA	2.32	0.65
1:A:22:LYS:H	1:A:22:LYS:HD3	1.62	0.64
1:A:184:MET:CE	1:A:184:MET:HA	2.28	0.64
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.79	0.64
2:B:426:TRP:O	2:B:429:LEU:HB2	1.97	0.64
2:B:242:GLN:NE2	2:B:242:GLN:HA	2.13	0.64
2:B:100:LEU:H	2:B:100:LEU:CD2	2.11	0.64
1:A:383:TRP:O	1:A:385:LYS:HG2	1.98	0.64
2:B:79:GLU:O	2:B:83:ARG:HG3	1.98	0.64
1:A:242:GLN:HB3	1:A:243:PRO:O	1.98	0.63
2:B:23:GLN:OE1	2:B:59:PRO:HA	1.99	0.63
1:A:181:TYR:CE1	1:A:183:TYR:HB2	2.34	0.63
2:B:100:LEU:HD23	2:B:100:LEU:N	2.14	0.63
1:A:253:THR:HG23	1:A:256:ASP:OD1	1.99	0.63
2:B:242:GLN:HB2	2:B:430:GLU:OE1	1.99	0.62
2:B:356:ARG:HH22	2:B:361:HIS:CE1	2.17	0.62
2:B:120:LEU:HD21	2:B:124:PHE:HD2	1.64	0.62
1:A:395:LYS:HD3	1:A:414:TRP:CZ2	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:180:ILE:HD11	2:B:189:VAL:HG13	1.82	0.62
1:A:517:LEU:O	1:A:520:GLN:HB2	1.98	0.62
1:A:541:GLY:HA3	2:B:284:ARG:CZ	2.30	0.61
1:A:40:GLU:O	1:A:43:LYS:HG2	2.00	0.61
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.82	0.61
1:A:358:ARG:NH2	1:A:512:GLN:HB2	2.16	0.61
2:B:24:TRP:CZ3	2:B:403:THR:HG21	2.35	0.61
2:B:390:LYS:HE2	2:B:415:GLU:OE2	1.99	0.61
2:B:65:LYS:NZ	2:B:72:ARG:HD3	2.14	0.61
2:B:154:LYS:O	2:B:157:PRO:HD2	2.01	0.61
1:A:40:GLU:OE2	1:A:43:LYS:NZ	2.32	0.61
1:A:255:ASN:O	1:A:259:LYS:HG3	2.00	0.61
1:A:22:LYS:H	1:A:22:LYS:CD	2.15	0.60
1:A:142:ILE:N	1:A:142:ILE:HD12	2.16	0.60
1:A:88:TRP:HE3	1:A:88:TRP:HA	1.65	0.60
1:A:68:SER:C	1:A:70:LYS:H	2.04	0.60
1:A:235:HIS:HB2	1:A:238:LYS:HD2	1.83	0.60
1:A:382:ILE:O	2:B:135:ILE:HG22	2.00	0.60
2:B:247:PRO:HB2	2:B:249:LYS:HZ1	1.67	0.60
1:A:22:LYS:HD3	1:A:22:LYS:N	2.17	0.60
1:A:195:ILE:HG13	1:A:196:GLY:N	2.17	0.60
2:B:52:PRO:HD2	2:B:53:GLU:OE1	2.01	0.60
2:B:295:LEU:HB3	2:B:300:GLU:HG2	1.84	0.60
1:A:195:ILE:HG13	1:A:196:GLY:H	1.66	0.59
2:B:254:VAL:HB	2:B:289:LEU:HA	1.83	0.59
2:B:279:LEU:HG	2:B:302:GLU:OE2	2.02	0.59
2:B:278:GLN:HB2	2:B:302:GLU:OE2	2.02	0.59
2:B:5:ILE:CG1	2:B:6:GLU:H	2.16	0.59
1:A:490:GLY:H	1:A:528:LYS:NZ	1.99	0.59
2:B:253:THR:O	2:B:257:ILE:HG12	2.03	0.59
1:A:229:TRP:O	1:A:230:MET:C	2.41	0.59
1:A:540:LYS:HG2	2:B:280:CYS:SG	2.42	0.59
2:B:151:GLN:HB3	2:B:185:ASP:OD1	2.02	0.59
1:A:257:ILE:O	1:A:260:LEU:HB3	2.02	0.58
2:B:5:ILE:HG12	2:B:6:GLU:N	2.17	0.58
1:A:486:LEU:CD1	1:A:521:ILE:HG23	2.32	0.58
1:A:38:CYS:O	1:A:42:GLU:HB2	2.04	0.58
1:A:246:LEU:HB2	1:A:307:ARG:NH1	2.18	0.58
2:B:135:ILE:O	2:B:138:GLU:HG2	2.03	0.58
2:B:5:ILE:HG13	2:B:6:GLU:H	1.67	0.58
1:A:244:ILE:HG23	1:A:310:LEU:HD13	1.83	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:PRO:HG3	1:A:313:PRO:HA	1.85	0.58
2:B:163:SER:O	2:B:167:ILE:HG13	2.04	0.58
1:A:42:GLU:CG	1:A:144:TYR:HE2	2.17	0.57
1:A:13:LYS:HB3	1:A:14:PRO:HD2	1.86	0.57
1:A:515:SER:HB3	1:A:518:VAL:CG2	2.34	0.57
2:B:380:ILE:O	2:B:384:GLY:N	2.37	0.57
1:A:173:LYS:O	1:A:176:PRO:HD3	2.04	0.57
1:A:139:THR:HG22	1:A:140:PRO:HD2	1.86	0.57
1:A:92:LEU:HD12	1:A:92:LEU:H	1.69	0.57
1:A:323:LYS:HE3	1:A:344:GLU:HG3	1.87	0.57
1:A:489:SER:OG	1:A:528:LYS:NZ	2.28	0.56
2:B:179:VAL:O	2:B:180:ILE:HD12	2.05	0.56
1:A:96:HIS:NE2	1:A:350:LYS:HE2	2.19	0.56
1:A:515:SER:HB3	1:A:518:VAL:HG23	1.88	0.56
1:A:457:TYR:HE1	1:A:463:ARG:HG2	1.69	0.56
1:A:410:TRP:NE1	2:B:363:ASN:HB2	2.21	0.56
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.88	0.56
2:B:139:THR:HB	2:B:140:PRO:HD2	1.88	0.56
1:A:244:ILE:HG21	1:A:310:LEU:HD13	1.88	0.56
1:A:226:PRO:HA	1:A:234:LEU:O	2.05	0.56
2:B:368:LEU:O	2:B:372:VAL:HG23	2.06	0.56
1:A:500:GLN:HA	1:A:500:GLN:OE1	2.05	0.56
2:B:424:LYS:O	2:B:424:LYS:HD2	2.05	0.56
1:A:515:SER:CB	1:A:518:VAL:HG23	2.36	0.56
1:A:254:VAL:HG22	1:A:293:ILE:HD11	1.88	0.56
1:A:116:PHE:O	1:A:148:VAL:HG21	2.05	0.55
1:A:27:THR:HG22	1:A:28:GLU:N	2.20	0.55
2:B:101:LYS:NZ	2:B:101:LYS:HB3	2.20	0.55
1:A:95:PRO:HA	2:B:136:ASN:O	2.06	0.55
1:A:540:LYS:HB3	1:A:542:ILE:CD1	2.36	0.55
2:B:157:PRO:CG	2:B:184:MET:HA	2.33	0.55
1:A:441:TYR:HA	1:A:496:VAL:HG13	1.89	0.55
1:A:76:ASP:OD1	1:A:78:ARG:HG3	2.07	0.55
2:B:278:GLN:HB2	2:B:302:GLU:CD	2.26	0.55
1:A:92:LEU:HD12	1:A:92:LEU:N	2.22	0.55
1:A:343:GLN:HG3	1:A:349:LEU:CD1	2.37	0.55
2:B:254:VAL:HG21	2:B:288:ALA:O	2.07	0.55
2:B:424:LYS:C	2:B:424:LYS:HD2	2.27	0.55
1:A:138:GLU:CG	1:A:139:THR:H	2.07	0.54
2:B:65:LYS:HA	2:B:407:GLN:HG2	1.87	0.54
1:A:42:GLU:HG3	1:A:144:TYR:HE2	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:GLU:HG3	1:A:139:THR:N	2.14	0.54
1:A:25:PRO:O	1:A:26:LEU:HD23	2.06	0.54
2:B:345:PRO:O	2:B:346:PHE:HB2	2.07	0.54
1:A:402:TRP:HZ3	2:B:364:ASP:OD2	1.90	0.54
2:B:420:PRO:HB2	2:B:423:VAL:CG1	2.37	0.54
2:B:98:ALA:HB1	2:B:101:LYS:HD3	1.89	0.54
2:B:172:ARG:HG3	2:B:172:ARG:HH11	1.72	0.54
1:A:417:VAL:HG13	1:A:419:THR:HG22	1.88	0.54
1:A:330:GLN:HB2	1:A:338:THR:OG1	2.07	0.54
1:A:37:ILE:CD1	1:A:72:ARG:HA	2.38	0.54
1:A:125:ARG:CD	1:A:147:ASN:HA	2.29	0.54
1:A:120:LEU:HD12	1:A:121:ASP:H	1.72	0.54
1:A:248:GLU:HA	1:A:248:GLU:OE2	2.06	0.53
1:A:252:TRP:CD1	1:A:295:LEU:HD12	2.43	0.53
2:B:295:LEU:HD22	2:B:300:GLU:OE2	2.08	0.53
2:B:433:PRO:CG	2:B:436:GLY:HA2	2.38	0.53
1:A:161:GLN:HE21	1:A:162:SER:N	2.05	0.53
1:A:226:PRO:HB3	1:A:235:HIS:NE2	2.23	0.53
1:A:89:GLU:HB2	1:A:92:LEU:HG	1.89	0.53
2:B:291:GLU:HG3	2:B:292:VAL:N	2.22	0.53
2:B:161:GLN:O	2:B:164:MET:HB3	2.08	0.53
2:B:365:VAL:O	2:B:369:THR:HG23	2.09	0.53
2:B:395:LYS:O	2:B:399:GLU:CG	2.52	0.52
2:B:13:LYS:HB2	2:B:16:MET:HG3	1.91	0.52
1:A:282:LEU:HB3	1:A:293:ILE:HG21	1.91	0.52
1:A:233:GLU:C	1:A:234:LEU:HD12	2.29	0.52
2:B:29:GLU:HG2	2:B:30:LYS:N	2.25	0.52
1:A:283:LEU:O	1:A:286:THR:HG23	2.10	0.52
1:A:206:ARG:NE	1:A:218:ASP:HA	2.24	0.52
1:A:58:THR:HG23	1:A:76:ASP:O	2.09	0.52
2:B:161:GLN:HA	2:B:161:GLN:NE2	2.25	0.52
1:A:8:VAL:O	1:A:10:VAL:HG23	2.09	0.52
1:A:226:PRO:HB3	1:A:235:HIS:CD2	2.45	0.52
1:A:41:MET:HA	1:A:44:GLU:HB2	1.91	0.52
1:A:198:HIS:CE1	1:A:202:ILE:HD11	2.45	0.52
1:A:109:LEU:N	1:A:109:LEU:HD12	2.25	0.52
1:A:505:ILE:O	1:A:509:GLN:N	2.41	0.52
2:B:120:LEU:HD12	2:B:150:PRO:HD3	1.91	0.51
2:B:63:ILE:HD13	2:B:72:ARG:O	2.10	0.51
1:A:295:LEU:HD22	1:A:300:GLU:CD	2.30	0.51
2:B:180:ILE:CD1	2:B:189:VAL:HG13	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:GLY:HA3	1:A:336:GLN:NE2	2.25	0.51
1:A:268:SER:HB2	1:A:353:LYS:HD3	1.92	0.51
2:B:182:GLN:HB2	2:B:187:LEU:CD1	2.39	0.51
2:B:175:ASN:N	2:B:176:PRO:HD3	2.25	0.51
2:B:332:GLN:OE1	2:B:424:LYS:HD3	2.10	0.51
1:A:516:GLU:O	1:A:520:GLN:HG2	2.11	0.51
1:A:332:GLN:HA	1:A:332:GLN:NE2	2.26	0.51
1:A:136:ASN:OD1	1:A:138:GLU:CB	2.55	0.51
1:A:356:ARG:HH11	1:A:356:ARG:HG3	1.76	0.51
2:B:65:LYS:O	2:B:68:SER:HB3	2.11	0.51
2:B:358:ARG:CD	2:B:358:ARG:H	2.24	0.51
2:B:248:GLU:HG2	2:B:307:ARG:NH2	2.26	0.51
2:B:195:ILE:HG12	2:B:199:ARG:CZ	2.41	0.51
1:A:169:GLU:N	1:A:170:PRO:HD2	2.26	0.51
1:A:83:ARG:HG3	1:A:83:ARG:HH11	1.75	0.51
1:A:6:GLU:N	1:A:6:GLU:CD	2.63	0.51
1:A:296:THR:HG23	1:A:299:ALA:CB	2.40	0.50
1:A:21:VAL:CG2	1:A:59:PRO:HD3	2.42	0.50
1:A:118:VAL:O	1:A:148:VAL:HG23	2.11	0.50
1:A:161:GLN:NE2	1:A:161:GLN:C	2.61	0.50
1:A:2:ILE:HG22	1:A:3:SER:N	2.26	0.50
1:A:233:GLU:HB3	1:A:240:THR:OG1	2.12	0.50
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.94	0.50
1:A:43:LYS:HZ3	1:A:44:GLU:HG3	1.76	0.50
1:A:358:ARG:CZ	1:A:358:ARG:CB	2.90	0.50
1:A:250:ASP:N	1:A:250:ASP:OD1	2.44	0.50
1:A:216:THR:OG1	1:A:217:PRO:HD2	2.12	0.50
2:B:78:ARG:O	2:B:82:LYS:HG3	2.12	0.50
1:A:136:ASN:O	1:A:138:GLU:N	2.45	0.50
1:A:367:GLN:O	1:A:368:LEU:C	2.50	0.50
2:B:241:VAL:HG13	2:B:351:THR:OG1	2.12	0.50
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.46	0.50
1:A:448:ARG:HG3	1:A:449:GLU:N	2.27	0.49
2:B:160:PHE:CD1	2:B:160:PHE:O	2.65	0.49
1:A:206:ARG:HG3	1:A:216:THR:HG23	1.93	0.49
2:B:78:ARG:HD3	2:B:411:ILE:O	2.11	0.49
1:A:344:GLU:O	1:A:345:PRO:C	2.50	0.49
1:A:254:VAL:HG21	1:A:288:ALA:O	2.12	0.49
1:A:139:THR:HG23	1:A:140:PRO:HD2	1.94	0.49
2:B:303:LEU:HD22	2:B:307:ARG:HE	1.77	0.49
2:B:78:ARG:NH1	2:B:412:PRO:O	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:PRO:HA	1:A:121:ASP:OD2	2.12	0.49
1:A:43:LYS:NZ	1:A:44:GLU:HG3	2.27	0.49
1:A:285:GLY:O	1:A:286:THR:C	2.51	0.49
1:A:283:LEU:HA	1:A:286:THR:HG23	1.94	0.49
2:B:53:GLU:CD	2:B:53:GLU:H	2.16	0.49
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.48	0.49
1:A:296:THR:O	1:A:299:ALA:HB3	2.12	0.49
1:A:10:VAL:HG11	1:A:153:TRP:CH2	2.47	0.49
1:A:329:ILE:HD11	1:A:375:ILE:HD12	1.94	0.49
2:B:296:THR:HB	2:B:298:GLU:OE2	2.13	0.49
1:A:151:GLN:NE2	1:A:151:GLN:HA	2.28	0.49
1:A:61:PHE:HB2	1:A:74:LEU:HD23	1.91	0.48
1:A:439:THR:HG21	2:B:289:LEU:HD13	1.95	0.48
1:A:283:LEU:HA	1:A:286:THR:CG2	2.43	0.48
2:B:72:ARG:HH11	2:B:72:ARG:HG3	1.78	0.48
1:A:234:LEU:CD1	1:A:234:LEU:N	2.76	0.48
2:B:418:ASN:N	2:B:418:ASN:OD1	2.46	0.48
1:A:212:TRP:C	1:A:214:LEU:H	2.16	0.48
1:A:319:TYR:O	1:A:321:PRO:HD3	2.13	0.48
1:A:106:VAL:HA	1:A:189:VAL:O	2.14	0.48
1:A:226:PRO:HB2	1:A:233:GLU:HG2	1.95	0.48
1:A:168:LEU:O	1:A:172:ARG:HG3	2.13	0.48
2:B:63:ILE:HD13	2:B:63:ILE:N	2.28	0.48
2:B:120:LEU:HD23	2:B:121:ASP:O	2.13	0.48
1:A:200:THR:O	1:A:203:GLU:HB3	2.14	0.48
2:B:332:GLN:HG3	2:B:338:THR:HG23	1.95	0.48
1:A:249:LYS:HD3	1:A:251:SER:O	2.14	0.48
2:B:374:LYS:HE2	2:B:374:LYS:HA	1.95	0.48
1:A:418:ASN:O	1:A:420:PRO:HD3	2.14	0.48
1:A:308:GLU:O	1:A:311:LYS:HB2	2.14	0.48
2:B:248:GLU:O	2:B:249:LYS:HD3	2.14	0.48
1:A:212:TRP:C	1:A:214:LEU:N	2.65	0.48
1:A:92:LEU:H	1:A:92:LEU:CD1	2.26	0.48
1:A:329:ILE:HG22	1:A:330:GLN:N	2.28	0.48
1:A:356:ARG:HH22	1:A:371:ALA:N	2.12	0.48
2:B:350:LYS:HG2	2:B:351:THR:N	2.28	0.48
1:A:465:LYS:HG2	1:A:466:VAL:H	1.79	0.47
2:B:75:VAL:HG12	2:B:76:ASP:N	2.30	0.47
1:A:243:PRO:CG	1:A:313:PRO:HB3	2.44	0.47
1:A:515:SER:OG	1:A:518:VAL:HG23	2.14	0.47
1:A:209:LEU:O	1:A:214:LEU:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:LEU:O	1:A:172:ARG:CG	2.61	0.47
2:B:109:LEU:O	2:B:186:ASP:HA	2.14	0.47
1:A:270:ILE:C	1:A:272:PRO:HD3	2.35	0.47
1:A:329:ILE:CG2	1:A:330:GLN:N	2.78	0.47
1:A:79:GLU:O	1:A:82:LYS:HB2	2.14	0.47
2:B:366:LYS:HG2	2:B:405:TYR:CD2	2.48	0.47
1:A:472:THR:OG1	1:A:473:THR:N	2.46	0.47
2:B:325:LEU:HD21	2:B:383:TRP:CE3	2.50	0.47
2:B:205:LEU:HD23	2:B:205:LEU:C	2.35	0.47
1:A:21:VAL:HG21	1:A:59:PRO:HD3	1.96	0.47
1:A:454:LYS:HA	1:A:467:VAL:O	2.14	0.47
2:B:247:PRO:HB2	2:B:249:LYS:HZ2	1.78	0.47
2:B:242:GLN:HE21	2:B:243:PRO:HD2	1.79	0.47
1:A:88:TRP:HD1	2:B:143:ARG:HD3	1.79	0.47
2:B:57:ASN:ND2	2:B:131:THR:OG1	2.48	0.47
1:A:390:LYS:HZ1	1:A:415:GLU:CD	2.18	0.47
2:B:170:PRO:HB2	2:B:208:HIS:HE1	1.79	0.47
2:B:244:ILE:HG21	2:B:426:TRP:CH2	2.50	0.47
2:B:126:LYS:HG3	2:B:127:TYR:N	2.30	0.47
1:A:239:TRP:HZ3	1:A:241:VAL:HG23	1.81	0.46
2:B:100:LEU:N	2:B:100:LEU:CD2	2.76	0.46
1:A:438:GLU:HA	1:A:461:ARG:HD2	1.97	0.46
1:A:347:LYS:HD3	1:A:347:LYS:HA	1.66	0.46
1:A:150:PRO:HG2	1:A:153:TRP:CB	2.45	0.46
1:A:363:ASN:HA	1:A:511:ASP:CG	2.34	0.46
1:A:522:ILE:HA	1:A:525:LEU:HD12	1.97	0.46
2:B:118:VAL:HB	2:B:149:LEU:CD1	2.45	0.46
1:A:333:GLY:HA3	1:A:336:GLN:HE22	1.78	0.46
2:B:32:LYS:O	2:B:33:ALA:C	2.52	0.46
1:A:318:TYR:CE2	3:A:999:NVP:H12	2.51	0.46
1:A:37:ILE:HD11	1:A:72:ARG:HA	1.98	0.46
2:B:363:ASN:ND2	2:B:366:LYS:HB2	2.30	0.46
2:B:108:VAL:O	2:B:109:LEU:HD12	2.16	0.46
2:B:382:ILE:HG22	2:B:383:TRP:CD2	2.51	0.46
1:A:419:THR:HG23	1:A:419:THR:O	2.16	0.46
2:B:238:LYS:O	2:B:239:TRP:CD1	2.65	0.46
1:A:484:LEU:O	1:A:487:GLN:HB2	2.16	0.46
1:A:13:LYS:CB	1:A:14:PRO:HD2	2.46	0.46
1:A:246:LEU:O	1:A:307:ARG:NH1	2.36	0.46
1:A:235:HIS:N	1:A:235:HIS:CD2	2.84	0.46
1:A:19:PRO:HG2	1:A:80:LEU:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ASP:O	1:A:83:ARG:CD	2.60	0.46
1:A:37:ILE:O	1:A:40:GLU:HB2	2.15	0.46
1:A:253:THR:O	1:A:256:ASP:HB2	2.16	0.46
2:B:379:SER:CB	2:B:387:PRO:HD3	2.46	0.45
1:A:393:ILE:HB	1:A:423:VAL:CG2	2.45	0.45
1:A:498:ASP:HA	1:A:536:VAL:O	2.16	0.45
2:B:108:VAL:HG22	2:B:188:TYR:CD2	2.51	0.45
1:A:37:ILE:HG22	1:A:38:CYS:N	2.31	0.45
1:A:136:ASN:O	1:A:137:ASN:C	2.53	0.45
2:B:303:LEU:HD22	2:B:307:ARG:NE	2.31	0.45
1:A:234:LEU:C	1:A:235:HIS:HD2	2.19	0.45
1:A:423:VAL:O	1:A:423:VAL:HG23	2.16	0.45
1:A:142:ILE:N	1:A:142:ILE:CD1	2.80	0.45
1:A:405:TYR:O	2:B:331:LYS:HD3	2.15	0.45
1:A:542:ILE:HD12	1:A:542:ILE:H	1.80	0.45
2:B:201:LYS:O	2:B:204:GLU:HB3	2.17	0.45
2:B:434:ILE:HG13	2:B:435:VAL:N	2.32	0.45
1:A:541:GLY:HA3	2:B:284:ARG:NH2	2.31	0.45
1:A:252:TRP:CD1	1:A:295:LEU:CD1	2.99	0.45
1:A:416:PHE:HE1	1:A:422:LEU:HD22	1.82	0.45
1:A:138:GLU:CG	1:A:139:THR:N	2.78	0.45
2:B:5:ILE:HG12	2:B:6:GLU:O	2.16	0.45
1:A:163:SER:O	1:A:166:LYS:HB2	2.16	0.45
2:B:47:ILE:HG22	2:B:146:TYR:HA	1.98	0.45
2:B:72:ARG:HH21	2:B:409:THR:HG22	1.81	0.45
1:A:42:GLU:HG3	1:A:144:TYR:CE2	2.50	0.45
1:A:94:ILE:HD11	1:A:183:TYR:CE2	2.51	0.45
2:B:104:LYS:HB2	2:B:192:ASP:HA	1.98	0.45
2:B:130:PHE:CZ	2:B:144:TYR:HB2	2.51	0.45
1:A:41:MET:HB2	1:A:46:LYS:HB2	1.98	0.44
1:A:161:GLN:HG2	1:A:182:GLN:OE1	2.17	0.44
1:A:243:PRO:HG3	1:A:313:PRO:CA	2.48	0.44
2:B:51:GLY:HA3	2:B:53:GLU:OE2	2.18	0.44
1:A:3:SER:OG	1:A:5:ILE:HG22	2.17	0.44
2:B:185:ASP:HB2	2:B:409:THR:HG21	1.99	0.44
1:A:359:GLY:O	1:A:360:ALA:C	2.56	0.44
1:A:249:LYS:CG	1:A:250:ASP:N	2.80	0.44
2:B:123:ASP:O	2:B:126:LYS:NZ	2.42	0.44
1:A:201:LYS:HA	1:A:201:LYS:HD3	1.74	0.44
1:A:28:GLU:HG2	1:A:32:LYS:HE3	2.00	0.44
2:B:195:ILE:CG2	2:B:196:GLY:N	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ASP:N	1:A:113:ASP:OD2	2.49	0.44
1:A:434:ILE:CD1	1:A:530:LYS:HB3	2.46	0.44
2:B:149:LEU:HA	2:B:150:PRO:HD3	1.87	0.44
2:B:87:PHE:CD1	2:B:87:PHE:N	2.86	0.44
2:B:31:ILE:O	2:B:35:VAL:HG23	2.18	0.44
1:A:376:THR:HG23	1:A:386:THR:HB	1.99	0.44
1:A:382:ILE:O	2:B:135:ILE:CG2	2.66	0.44
2:B:114:ALA:HB1	2:B:160:PHE:CZ	2.53	0.44
1:A:307:ARG:HG2	1:A:307:ARG:HH11	1.82	0.44
2:B:366:LYS:O	2:B:370:GLU:HG3	2.18	0.44
2:B:28:GLU:O	2:B:29:GLU:C	2.56	0.44
1:A:187:LEU:HA	1:A:187:LEU:HD12	1.81	0.44
1:A:42:GLU:CG	1:A:144:TYR:CE2	3.00	0.43
1:A:161:GLN:O	1:A:164:MET:HB3	2.18	0.43
1:A:356:ARG:NH1	1:A:356:ARG:HG3	2.33	0.43
2:B:104:LYS:CB	2:B:192:ASP:HA	2.48	0.43
1:A:475:GLN:O	1:A:479:LEU:HG	2.18	0.43
1:A:64:LYS:NZ	1:A:69:THR:HG23	2.33	0.43
1:A:412:PRO:O	1:A:413:GLU:C	2.55	0.43
1:A:340:GLN:CB	1:A:351:THR:HG22	2.48	0.43
2:B:172:ARG:HH21	2:B:180:ILE:HG22	1.83	0.43
1:A:357:MET:O	1:A:359:GLY:N	2.51	0.43
1:A:254:VAL:HG13	1:A:286:THR:HG21	2.00	0.43
1:A:480:GLN:NE2	1:A:484:LEU:HG	2.33	0.43
1:A:91:GLN:C	1:A:93:GLY:H	2.21	0.43
2:B:377:THR:O	2:B:380:ILE:HB	2.19	0.43
2:B:344:GLU:HB2	2:B:347:LYS:HD2	2.01	0.43
1:A:278:GLN:HG2	1:A:298:GLU:CB	2.44	0.43
1:A:94:ILE:HG22	1:A:95:PRO:O	2.18	0.43
1:A:508:ALA:O	1:A:509:GLN:C	2.57	0.43
1:A:19:PRO:HG3	1:A:80:LEU:HA	2.01	0.43
2:B:317:VAL:O	2:B:317:VAL:HG23	2.19	0.43
1:A:360:ALA:O	1:A:513:SER:HB2	2.18	0.43
1:A:463:ARG:NH1	1:A:488:ASP:O	2.52	0.43
1:A:151:GLN:HE21	1:A:151:GLN:CA	2.31	0.43
1:A:69:THR:HG22	1:A:69:THR:O	2.19	0.43
1:A:370:GLU:O	1:A:374:LYS:HG3	2.19	0.43
1:A:340:GLN:HA	1:A:351:THR:HG22	2.00	0.43
1:A:105:SER:O	1:A:190:GLY:HA2	2.17	0.43
2:B:245:VAL:CG2	2:B:431:LYS:HB2	2.47	0.43
2:B:201:LYS:HA	2:B:201:LYS:HD2	1.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:LEU:HD21	2:B:141:GLY:H	1.83	0.43
1:A:184:MET:HE2	1:A:184:MET:HA	2.00	0.43
2:B:101:LYS:CB	2:B:101:LYS:NZ	2.82	0.43
1:A:219:LYS:HA	1:A:222:GLN:NE2	2.33	0.43
1:A:293:ILE:N	1:A:293:ILE:HD12	2.34	0.43
1:A:150:PRO:HG2	1:A:153:TRP:HB3	2.00	0.43
1:A:324:ASP:OD2	1:A:324:ASP:N	2.51	0.43
2:B:115:TYR:OH	2:B:157:PRO:HG3	2.19	0.43
1:A:5:ILE:HG13	1:A:6:GLU:OE2	2.19	0.43
2:B:244:ILE:HG23	2:B:429:LEU:HB3	2.01	0.43
2:B:301:LEU:O	2:B:302:GLU:C	2.57	0.43
1:A:219:LYS:HD3	1:A:222:GLN:NE2	2.34	0.43
1:A:112:GLY:C	1:A:114:ALA:H	2.20	0.43
2:B:278:GLN:HB3	2:B:299:ALA:HA	2.01	0.43
1:A:151:GLN:HE21	1:A:151:GLN:HA	1.82	0.43
1:A:363:ASN:HB2	1:A:511:ASP:OD2	2.18	0.43
2:B:63:ILE:HD12	2:B:74:LEU:HB2	2.01	0.42
2:B:23:GLN:HG3	2:B:24:TRP:O	2.19	0.42
1:A:373:GLN:C	1:A:373:GLN:CD	2.78	0.42
2:B:175:ASN:HB3	2:B:178:ILE:HD12	2.00	0.42
1:A:249:LYS:HG3	1:A:250:ASP:N	2.34	0.42
1:A:232:TYR:CD1	1:A:232:TYR:N	2.87	0.42
1:A:57:ASN:HA	1:A:129:ALA:O	2.19	0.42
2:B:136:ASN:O	2:B:137:ASN:HB2	2.19	0.42
1:A:371:ALA:O	1:A:372:VAL:C	2.57	0.42
2:B:358:ARG:H	2:B:358:ARG:HD2	1.84	0.42
1:A:249:LYS:HE3	1:A:249:LYS:HB2	1.67	0.42
1:A:91:GLN:O	1:A:93:GLY:N	2.49	0.42
1:A:11:LYS:O	1:A:85:GLN:HB3	2.20	0.42
2:B:191:SER:HB2	2:B:193:LEU:HD13	2.00	0.42
1:A:194:GLU:HG3	1:A:197:GLN:HG3	2.02	0.42
1:A:240:THR:HG22	1:A:315:HIS:ND1	2.35	0.42
2:B:285:GLY:O	2:B:287:LYS:HG2	2.19	0.42
2:B:435:VAL:HG23	2:B:436:GLY:N	2.35	0.42
2:B:66:LYS:O	2:B:67:ASP:HB2	2.19	0.42
1:A:478:GLU:O	1:A:481:ALA:HB3	2.19	0.42
2:B:242:GLN:NE2	2:B:243:PRO:HD2	2.34	0.42
1:A:88:TRP:HB3	1:A:89:GLU:OE1	2.20	0.42
1:A:110:ASP:O	1:A:217:PRO:HD3	2.19	0.42
2:B:52:PRO:C	2:B:54:ASN:N	2.70	0.42
1:A:219:LYS:HA	1:A:222:GLN:CD	2.39	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:242:GLN:CA	2:B:242:GLN:HE21	2.11	0.42
1:A:13:LYS:HD2	1:A:82:LYS:O	2.19	0.42
1:A:541:GLY:HA3	2:B:284:ARG:NE	2.34	0.42
2:B:78:ARG:NH1	2:B:411:ILE:HG22	2.34	0.42
2:B:101:LYS:HE2	2:B:382:ILE:HG23	2.02	0.42
2:B:66:LYS:HD2	2:B:66:LYS:HA	1.94	0.42
2:B:116:PHE:HA	2:B:148:VAL:HG21	2.02	0.42
1:A:180:ILE:HA	1:A:188:TYR:O	2.20	0.42
1:A:33:ALA:O	1:A:36:GLU:N	2.53	0.42
1:A:514:GLU:HG3	1:A:514:GLU:H	1.52	0.42
1:A:61:PHE:CB	1:A:74:LEU:HD21	2.38	0.41
2:B:387:PRO:HG2	2:B:389:PHE:HE1	1.81	0.41
2:B:288:ALA:O	2:B:291:GLU:HB3	2.20	0.41
1:A:447:ASN:HB3	1:A:450:THR:OG1	2.20	0.41
1:A:58:THR:HG23	1:A:59:PRO:HD2	2.02	0.41
1:A:181:TYR:HB3	1:A:188:TYR:HB2	2.02	0.41
1:A:344:GLU:O	1:A:347:LYS:HB2	2.20	0.41
1:A:27:THR:HG22	1:A:29:GLU:H	1.85	0.41
1:A:328:GLU:HB3	1:A:340:GLN:OE1	2.20	0.41
1:A:303:LEU:HD23	1:A:303:LEU:O	2.20	0.41
2:B:280:CYS:C	2:B:284:ARG:HH21	2.23	0.41
2:B:34:LEU:HD23	2:B:34:LEU:HA	1.85	0.41
1:A:27:THR:HG22	1:A:28:GLU:H	1.84	0.41
1:A:111:VAL:HG22	1:A:185:ASP:O	2.20	0.41
1:A:380:ILE:O	1:A:384:GLY:N	2.48	0.41
1:A:408:ALA:O	2:B:393:ILE:HG13	2.21	0.41
1:A:95:PRO:HB3	2:B:136:ASN:O	2.21	0.41
1:A:356:ARG:HH21	1:A:374:LYS:HZ1	1.69	0.41
1:A:154:LYS:O	1:A:157:PRO:HD2	2.21	0.41
1:A:205:LEU:O	1:A:208:HIS:HB3	2.20	0.41
1:A:224:GLU:HA	1:A:225:PRO:HD3	1.94	0.41
1:A:115:TYR:O	1:A:149:LEU:HB2	2.19	0.41
1:A:136:ASN:C	1:A:138:GLU:N	2.72	0.41
1:A:63:ILE:HG22	1:A:74:LEU:HD22	2.03	0.41
1:A:513:SER:OG	1:A:518:VAL:HB	2.21	0.41
2:B:252:TRP:CZ3	2:B:260:LEU:HD22	2.55	0.41
1:A:206:ARG:HH21	1:A:218:ASP:HB3	1.85	0.41
2:B:51:GLY:CA	2:B:53:GLU:OE2	2.68	0.41
1:A:457:TYR:CE2	1:A:465:LYS:HB3	2.56	0.41
1:A:205:LEU:O	1:A:208:HIS:N	2.54	0.41
2:B:206:ARG:HG2	2:B:210:LEU:CD1	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:MET:HB3	1:A:46:LYS:HE2	2.03	0.41
1:A:101:LYS:O	3:A:999:NVP:H13	2.21	0.41
1:A:480:GLN:HE22	1:A:484:LEU:HG	1.86	0.41
2:B:106:VAL:O	2:B:234:LEU:HB2	2.20	0.41
1:A:410:TRP:HZ3	2:B:401:TRP:CE2	2.39	0.40
2:B:195:ILE:HG23	2:B:196:GLY:N	2.36	0.40
1:A:37:ILE:CG2	1:A:38:CYS:N	2.84	0.40
2:B:431:LYS:HG2	2:B:432:GLU:N	2.36	0.40
2:B:363:ASN:C	2:B:363:ASN:OD1	2.60	0.40
1:A:184:MET:HE3	1:A:184:MET:HA	2.02	0.40
1:A:99:GLY:HA3	1:A:382:ILE:HG23	2.02	0.40
2:B:341:ILE:HD12	2:B:341:ILE:N	2.36	0.40
1:A:295:LEU:HD22	1:A:300:GLU:OE2	2.22	0.40
1:A:2:ILE:HG22	1:A:3:SER:H	1.85	0.40
1:A:295:LEU:HD23	1:A:299:ALA:HB3	2.04	0.40
1:A:300:GLU:HA	1:A:300:GLU:OE2	2.20	0.40
2:B:180:ILE:O	2:B:180:ILE:HG22	2.21	0.40
2:B:383:TRP:O	2:B:384:GLY:C	2.60	0.40
1:A:214:LEU:HA	1:A:214:LEU:HD23	1.78	0.40
1:A:254:VAL:HG22	1:A:293:ILE:CD1	2.50	0.40
1:A:118:VAL:O	1:A:148:VAL:CG2	2.69	0.40
2:B:345:PRO:O	2:B:346:PHE:CB	2.70	0.40
1:A:356:ARG:NH1	1:A:367:GLN:HB3	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	539/543 (99%)	470 (87%)	49 (9%)	20 (4%)	4	17
2	B	402/440 (91%)	349 (87%)	47 (12%)	6 (2%)	13	42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	941/983 (96%)	819 (87%)	96 (10%)	26 (3%)	6	24

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	ILE
1	A	242	GLN
2	B	162	SER
1	A	88	TRP
1	A	91	GLN
1	A	92	LEU
1	A	137	ASN
1	A	230	MET
1	A	358	ARG
2	B	361	HIS
1	A	18	GLY
1	A	68	SER
1	A	250	ASP
1	A	360	ALA
2	B	267	ALA
2	B	238	LYS
1	A	4	PRO
1	A	69	THR
1	A	85	GLN
1	A	286	THR
1	A	310	LEU
1	A	345	PRO
1	A	236	PRO
1	A	195	ILE
2	B	167	ILE
2	B	236	PRO

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	485/485 (100%)	437 (90%)	48 (10%)	10	29
2	B	370/400 (92%)	338 (91%)	32 (9%)	13	36
All	All	855/885 (97%)	775 (91%)	80 (9%)	11	32

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	PRO
1	A	6	GLU
1	A	11	LYS
1	A	22	LYS
1	A	37	ILE
1	A	40	GLU
1	A	42	GLU
1	A	88	TRP
1	A	89	GLU
1	A	92	LEU
1	A	105	SER
1	A	122	GLU
1	A	161	GLN
1	A	182	GLN
1	A	184	MET
1	A	189	VAL
1	A	194	GLU
1	A	205	LEU
1	A	215	THR
1	A	216	THR
1	A	228	LEU
1	A	230	MET
1	A	238	LYS
1	A	250	ASP
1	A	253	THR
1	A	295	LEU
1	A	296	THR
1	A	303	LEU
1	A	312	GLU
1	A	336	GLN
1	A	340	GLN
1	A	358	ARG
1	A	361	HIS
1	A	362	THR
1	A	369	THR

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Mol	Chain	Res	Type
1	A	373	GLN
1	A	386	THR
1	A	388	LYS
1	A	424	LYS
1	A	449	GLU
1	A	475	GLN
1	A	480	GLN
1	A	493	VAL
1	A	496	VAL
1	A	514	GLU
1	A	517	LEU
1	A	527	LYS
1	A	540	LYS
2	B	6	GLU
2	B	24	TRP
2	B	55	PRO
2	B	63	ILE
2	B	65	LYS
2	B	69	THR
2	B	72	ARG
2	B	87	PHE
2	B	100	LEU
2	B	103	ASN
2	B	123	ASP
2	B	134	SER
2	B	161	GLN
2	B	162	SER
2	B	163	SER
2	B	164	MET
2	B	186	ASP
2	B	194	GLU
2	B	236	PRO
2	B	237	ASP
2	B	242	GLN
2	B	281	LYS
2	B	283	LEU
2	B	287	LYS
2	B	289	LEU
2	B	291	GLU
2	B	358	ARG
2	B	362	THR
2	B	363	ASN

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Mol	Chain	Res	Type
2	B	368	LEU
2	B	418	ASN
2	B	424	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	103	ASN
1	A	137	ASN
1	A	151	GLN
1	A	161	GLN
1	A	197	GLN
1	A	278	GLN
1	A	332	GLN
1	A	336	GLN
1	A	367	GLN
1	A	475	GLN
1	A	480	GLN
2	B	57	ASN
2	B	85	GLN
2	B	147	ASN
2	B	151	GLN
2	B	242	GLN
2	B	278	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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
1	CSD	A	280	1	3,7,8	0.78	0	3,8,10	1.98	2 (66%)

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
1	CSD	A	280	1	-	1/2/6/8	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	CSD	O-C-CA	-2.24	119.66	125.49
1	A	280	CSD	OD1-SG-CB	2.58	109.69	105.40

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	280	CSD	CA-CB-SG-OD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	NVP	A	999	-	18,23,23	1.35	4 (22%)	18,34,34	1.19	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
3	NVP	A	999	-	-	0/0/6/6	0/2/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	999	NVP	C4-N3	2.09	1.36	1.32
3	A	999	NVP	C13-N14	2.20	1.36	1.32
3	A	999	NVP	C12-C11	2.34	1.42	1.36
3	A	999	NVP	C10-C15	3.19	1.43	1.40

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	999	NVP	CC-CA-N1	-2.27	115.91	118.25
3	A	999	NVP	CB-CA-N1	-2.05	116.14	118.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	999	NVP	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	541/543 (99%)	-0.09	26 (4%) 34 28	20, 63, 127, 150	0
2	B	408/440 (92%)	-0.12	16 (3%) 43 36	23, 62, 114, 149	0
All	All	949/983 (96%)	-0.10	42 (4%) 38 32	20, 63, 122, 150	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	435	VAL	8.0
2	B	436	GLY	6.7
2	B	434	ILE	6.1
2	B	437	ALA	5.6
1	A	67	ASP	5.5
1	A	68	SER	5.3
1	A	69	THR	5.0
1	A	66	LYS	4.2
1	A	29	GLU	3.7
1	A	139	THR	3.6
2	B	438	GLU	3.6
2	B	67	ASP	3.5
1	A	71	TRP	3.4
2	B	71	TRP	3.3
1	A	52	PRO	3.3
1	A	91	GLN	3.2
2	B	213	GLY	3.1
1	A	28	GLU	3.0
1	A	138	GLU	3.0
1	A	542	ILE	3.0
2	B	197	GLN	3.0
1	A	62	ALA	2.9
1	A	135	ILE	2.8
2	B	432	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	14	PRO	2.7
1	A	63	ILE	2.5
2	B	242	GLN	2.5
2	B	284	ARG	2.4
1	A	140	PRO	2.4
1	A	473	THR	2.3
1	A	51	GLY	2.3
2	B	362	THR	2.3
1	A	90	VAL	2.3
1	A	245	VAL	2.2
2	B	85	GLN	2.2
1	A	448	ARG	2.2
1	A	32	LYS	2.2
1	A	357	MET	2.1
1	A	21	VAL	2.1
2	B	6	GLU	2.1
1	A	136	ASN	2.1
2	B	317	VAL	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CSD	A	280	8/9	0.96	0.13	-	36,64,80,84	0

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
3	NVP	A	999	20/20	0.96	0.17	0.45	34,54,73,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.