



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:02 PM GMT

PDB ID : 3SN6
Title : Crystal structure of the beta2 adrenergic receptor-Gs protein complex
Authors : Rasmussen, S.G.F.; DeVree, B.T.; Zou, Y.; Kruse, A.C.; Chung, K.Y.; Kobilka, T.S.; Thian, F.S.; Chae, P.S.; Pardon, E.; Calinski, D.; Mathiesen, J.M.; Shah, S.T.A.; Lyons, J.A.; Caffrey, M.; Gellman, S.H.; Steyaert, J.; Skiniotis, G.; Weis, W.I.; Sunahara, R.K.; Kobilka, B.K.
Deposited on : 2011-06-28
Resolution : 3.20 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

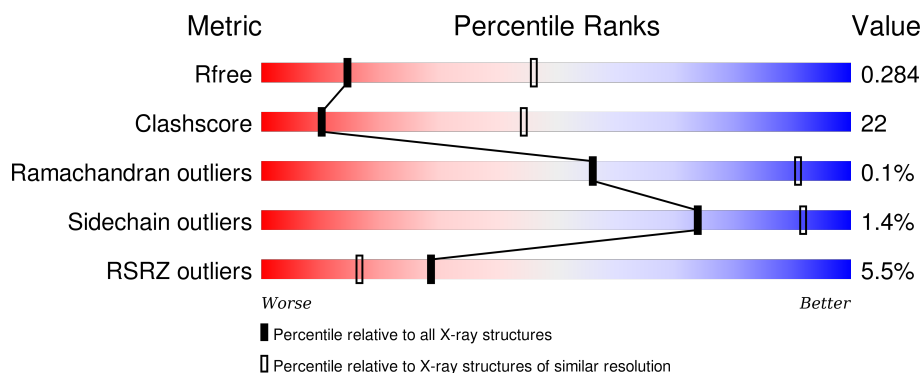
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	380	<div> <div>6%</div> <div>58% 33% 8%</div> </div>
2	B	351	<div> <div>%</div> <div>57% 39% .</div> </div>
3	G	68	<div> <div>%</div> <div>68% 18% 15%</div> </div>
4	R	514	<div> <div>9%</div> <div>49% 36% . 14%</div> </div>
5	N	138	<div> <div></div> <div>51% 41% . 7%</div> </div>

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
6	P0G	R	366	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10274 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 Guanine nucleotide-binding protein G(s) subunit alpha isoforms short.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	0	0
			2814	1786	494	522	12			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	86	SER	GLY	ENGINEERED MUTATION	UNP P04896

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	340	Total	C	N	O	S	0	0	0
			2592	1600	463	508	21			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	MET	-	EXPRESSION TAG	UNP P54311
B	-9	HIS	-	EXPRESSION TAG	UNP P54311
B	-8	HIS	-	EXPRESSION TAG	UNP P54311
B	-7	HIS	-	EXPRESSION TAG	UNP P54311
B	-6	HIS	-	EXPRESSION TAG	UNP P54311
B	-5	HIS	-	EXPRESSION TAG	UNP P54311
B	-4	HIS	-	EXPRESSION TAG	UNP P54311
B	-3	GLY	-	EXPRESSION TAG	UNP P54311
B	-2	SER	-	EXPRESSION TAG	UNP P54311
B	-1	LEU	-	EXPRESSION TAG	UNP P54311
B	0	LEU	-	EXPRESSION TAG	UNP P54311
B	1	GLN	-	EXPRESSION TAG	UNP P54311

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	58	Total	C	N	O	S	0	0	0
			438	274	76	85	3			

- Molecule 4 is a protein called Lysozyme, Beta-2 adrenergic receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	R	443	Total	C	N	O	S	0	0	0
			3433	2234	572	605	22			

There are 20 discrepancies between the modelled and reference sequences:

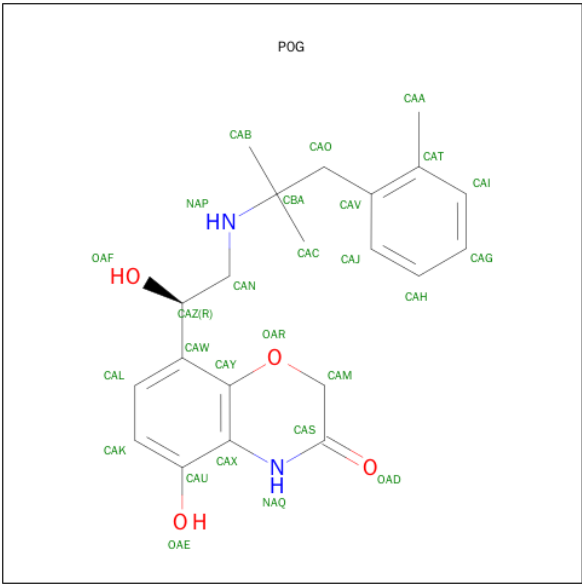
Chain	Residue	Modelled	Actual	Comment	Reference
R	987	ASP	-	EXPRESSION TAG	UNP P07550
R	988	TYR	-	EXPRESSION TAG	UNP P07550
R	989	LYS	-	EXPRESSION TAG	UNP P07550
R	990	ASP	-	EXPRESSION TAG	UNP P07550
R	991	ASP	-	EXPRESSION TAG	UNP P07550
R	992	ASP	-	EXPRESSION TAG	UNP P07550
R	993	ASP	-	EXPRESSION TAG	UNP P07550
R	994	ALA	-	EXPRESSION TAG	UNP P07550
R	995	GLU	-	EXPRESSION TAG	UNP P07550
R	996	ASN	-	EXPRESSION TAG	UNP P07550
R	997	LEU	-	EXPRESSION TAG	UNP P07550
R	998	TYR	-	EXPRESSION TAG	UNP P07550
R	999	PHE	-	EXPRESSION TAG	UNP P07550
R	1000	GLN	-	EXPRESSION TAG	UNP P07550
R	1001	GLY	-	EXPRESSION TAG	UNP P07550
R	96	THR	MET	ENGINEERED MUTATION	UNP P07550
R	98	THR	MET	ENGINEERED MUTATION	UNP P07550
R	187	GLU	ASN	ENGINEERED MUTATION	UNP P07550
R	1054	THR	CYS	ENGINEERED MUTATION	UNP P00720
R	1097	ALA	CYS	ENGINEERED MUTATION	UNP P00720

- Molecule 5 is a protein called Camelid antibody VHH fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	N	128	Total	C	N	O	S	0	0	0
			970	604	170	190	6			

- Molecule 6 is 8-[(1R)-2-{{1,1-DIMETHYL-2-(2-METHYLPHENYL)ETHYL}AMINO}-1-

HYDROXYETHYL]-5-HYDROXY-2H-1,4-BENZOXAZIN-3(4H)-ONE (three-letter code: P0G) (formula: C₂₁H₂₆N₂O₄).

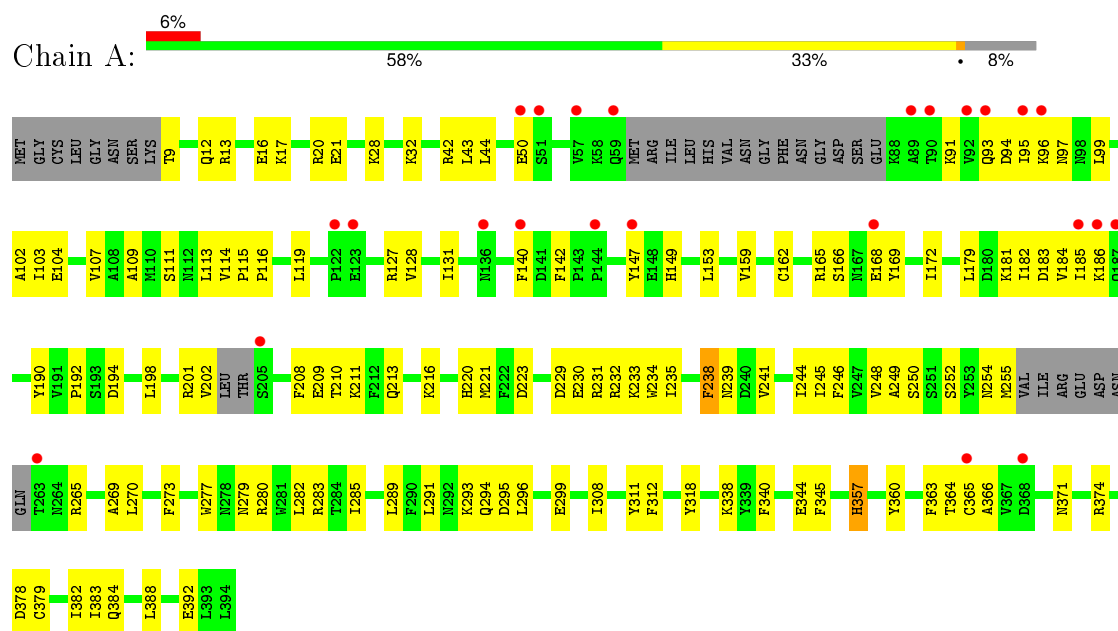


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	R	1	27	21	2	4	0	0

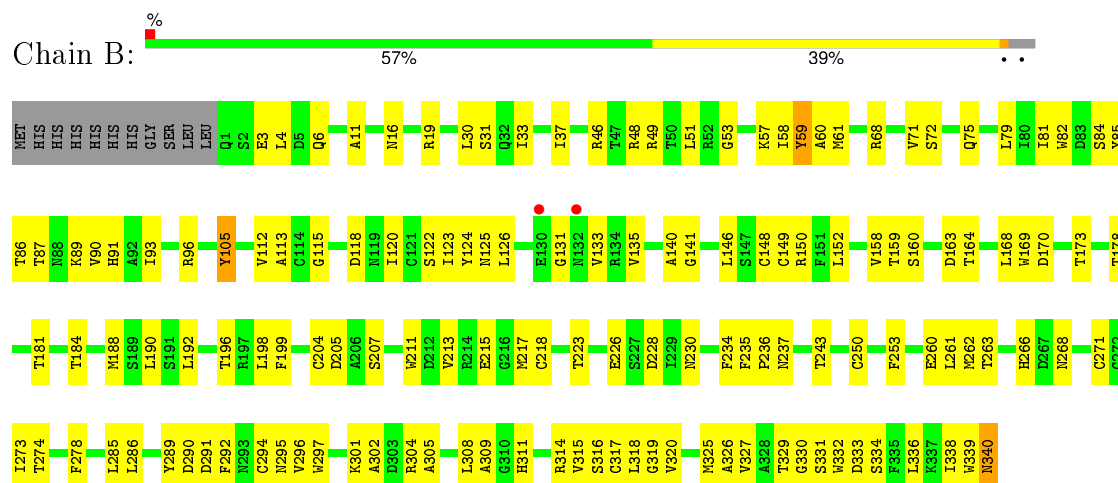
3 Residue-property plots

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.

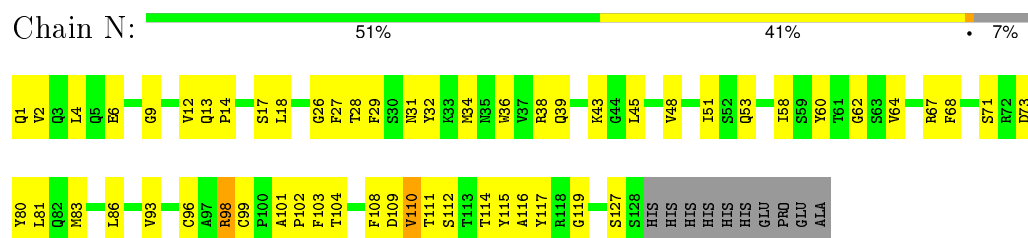
- Molecule 1: Guanine nucleotide-binding protein G(s) subunit alpha isoforms short



- Molecule 2: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1



- Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	119.34Å 64.56Å 131.24Å 90.00° 91.67° 90.00°	Depositor
Resolution (Å)	40.67 – 3.20 40.68 – 3.20	Depositor EDS
% Data completeness (in resolution range)	92.8 (40.67-3.20) 92.9 (40.68-3.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 3.18Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.226 , 0.277 0.236 , 0.284	Depositor DCC
R_{free} test set	1558 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	68.1	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 61.2	EDS
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 31099 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	10274	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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: P0G

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	0.23	0/2870	0.40	0/3879
2	B	0.23	0/2639	0.46	0/3580
3	G	0.22	0/444	0.38	0/601
4	R	0.23	0/3502	0.40	0/4762
5	N	0.28	0/990	0.53	1/1341 (0.1%)
All	All	0.23	0/10445	0.43	1/14163 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	N	99	CYS	CA-CB-SG	-5.45	104.19	114.00

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	2814	0	2719	110	0
2	B	2592	0	2485	134	0
3	G	438	0	443	11	0
4	R	3433	0	3429	143	0
5	N	970	0	930	59	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	R	27	0	25	7	0
All	All	10274	0	10031	440	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 22.

All (440) 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)
4:R:132:TYR:HB2	4:R:218:VAL:HG13	1.43	0.99
4:R:137:SER:HB3	4:R:140:LYS:HD3	1.43	0.97
1:A:103:ILE:HG22	1:A:179:LEU:HD21	1.51	0.93
1:A:114:VAL:HB	1:A:115:PRO:HD3	1.48	0.92
4:R:293:ASN:HD21	6:R:366:P0G:HAMA	1.39	0.88
4:R:72:ILE:HD11	4:R:326:TYR:HB3	1.57	0.83
1:A:294:GLN:HB2	1:A:366:ALA:HB2	1.61	0.83
2:B:262:MET:SD	2:B:302:ALA:HB2	2.18	0.82
1:A:202:VAL:CG2	1:A:211:LYS:HD2	2.09	0.81
5:N:1:GLN:HG2	5:N:2:VAL:H	1.47	0.81
5:N:71:SER:O	5:N:80:TYR:HB2	1.81	0.80
1:A:378:ASP:O	1:A:382:ILE:HG13	1.80	0.80
4:R:1146:ALA:O	4:R:1150:ILE:HG13	1.83	0.78
2:B:37:ILE:HD11	3:G:38:MET:SD	2.23	0.78
2:B:53:GLY:N	2:B:82:TRP:HH2	1.81	0.77
1:A:388:LEU:HD21	4:R:226:ALA:HA	1.67	0.77
2:B:57:LYS:HE2	2:B:75:GLN:HG3	1.67	0.76
2:B:149:CYS:O	2:B:150:ARG:HD3	1.86	0.75
2:B:4:LEU:HD13	3:G:9:ILE:HG22	1.68	0.75
2:B:286:LEU:HD22	2:B:296:VAL:HG22	1.68	0.75
1:A:99:LEU:HD11	1:A:182:ILE:HG12	1.68	0.74
1:A:202:VAL:HG23	1:A:211:LYS:HD2	1.70	0.74
1:A:202:VAL:HG22	1:A:202:VAL:O	1.88	0.73
2:B:68:ARG:HG3	2:B:85:TYR:CD2	2.25	0.72
4:R:74:SER:HB2	4:R:154:ILE:HD12	1.72	0.72
1:A:107:VAL:HG11	1:A:153:LEU:HD13	1.72	0.72
4:R:117:VAL:HG21	6:R:366:P0G:HAL	1.72	0.71
4:R:149:LYS:O	4:R:153:ILE:HG12	1.91	0.71
1:A:168:GLU:HB2	2:B:131:GLY:HA3	1.72	0.70
5:N:32:TYR:OH	5:N:102:PRO:HG3	1.91	0.70
5:N:14:PRO:HD3	5:N:127:SER:O	1.90	0.70
4:R:288:PRO:HB2	4:R:311:LEU:HD22	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:304:ARG:HG3	2:B:304:ARG:O	1.92	0.69
2:B:120:ILE:HG12	2:B:140:ALA:HB2	1.74	0.69
3:G:16:VAL:O	3:G:20:LYS:HB2	1.92	0.69
1:A:179:LEU:HD23	1:A:182:ILE:HD11	1.73	0.69
2:B:68:ARG:HG3	2:B:85:TYR:HD2	1.58	0.69
1:A:166:SER:HA	1:A:169:TYR:CE1	2.28	0.69
2:B:49:ARG:HB2	2:B:338:ILE:HB	1.74	0.68
2:B:168:LEU:HD22	2:B:213:VAL:HG13	1.74	0.68
4:R:1025:TYR:HE2	4:R:1039:LEU:HD12	1.58	0.68
1:A:114:VAL:O	1:A:116:PRO:HD3	1.93	0.67
2:B:163:ASP:O	2:B:164:THR:HB	1.94	0.67
4:R:1126:TRP:HB3	4:R:1154:ARG:HA	1.76	0.66
1:A:229:ASP:CG	5:N:111:THR:HG23	2.16	0.66
5:N:1:GLN:O	5:N:26:GLY:HA3	1.95	0.66
1:A:208:PHE:HB3	1:A:223:ASP:HB3	1.78	0.65
4:R:287:LEU:HB3	4:R:288:PRO:HD3	1.78	0.65
4:R:210:VAL:HB	4:R:211:PRO:HD3	1.78	0.65
5:N:110:VAL:O	5:N:110:VAL:HG12	1.96	0.65
5:N:9:GLY:HA2	5:N:18:LEU:HD21	1.78	0.65
1:A:202:VAL:HB	1:A:211:LYS:CG	2.26	0.65
2:B:146:LEU:HD11	2:B:159:THR:HB	1.78	0.65
2:B:51:LEU:HB3	2:B:82:TRP:CZ3	2.31	0.65
1:A:9:THR:HG22	1:A:13:ARG:HH21	1.62	0.64
2:B:3:GLU:O	2:B:6:GLN:HG2	1.97	0.64
2:B:318:LEU:HG	2:B:329:THR:HG22	1.79	0.64
5:N:12:VAL:HG11	5:N:86:LEU:HD22	1.79	0.64
2:B:141:GLY:N	2:B:169:TRP:HH2	1.96	0.64
1:A:311:TYR:O	5:N:62:GLY:HA3	1.97	0.64
2:B:311:HIS:HB3	2:B:333:ASP:OD2	1.98	0.63
2:B:90:VAL:HG12	2:B:91:HIS:ND1	2.13	0.63
2:B:198:LEU:HB2	2:B:211:TRP:O	1.98	0.63
4:R:86:VAL:HG22	4:R:112:ILE:HG23	1.81	0.62
5:N:112:SER:O	5:N:114:THR:HG22	1.99	0.62
3:G:19:LEU:O	3:G:23:ALA:HB2	1.98	0.62
2:B:266:HIS:CD2	2:B:268:ASN:HB2	2.33	0.62
2:B:124:TYR:CD1	2:B:135:VAL:HA	2.34	0.62
4:R:1059:THR:HB	4:R:1062:GLU:HG3	1.80	0.62
2:B:81:ILE:HB	2:B:91:HIS:HB2	1.81	0.62
2:B:126:LEU:HD23	2:B:133:VAL:HG11	1.82	0.62
2:B:115:GLY:HA3	2:B:146:LEU:HD23	1.82	0.61
2:B:160:SER:HB3	2:B:190:LEU:HD23	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:109:TRP:O	4:R:112:ILE:HG22	2.00	0.61
2:B:311:HIS:H	2:B:311:HIS:CD2	2.16	0.61
1:A:279:ASN:O	1:A:283:ARG:HG3	2.00	0.61
4:R:69:ASN:O	4:R:72:ILE:HG22	2.02	0.60
1:A:239:ASN:O	1:A:285:ILE:HD11	2.02	0.60
1:A:246:PHE:HE1	1:A:273:PHE:HB2	1.65	0.60
2:B:93:ILE:HG12	2:B:133:VAL:HG21	1.83	0.60
4:R:1014:ARG:HG3	4:R:1018:TYR:CD2	2.37	0.59
4:R:132:TYR:HB2	4:R:218:VAL:CG1	2.26	0.59
4:R:285:CYS:HA	4:R:314:ILE:HG22	1.84	0.59
2:B:301:LYS:O	2:B:302:ALA:HB3	2.03	0.59
4:R:1025:TYR:CE2	4:R:1039:LEU:HD12	2.38	0.59
1:A:249:ALA:HB1	1:A:293:LYS:HD2	1.85	0.58
1:A:202:VAL:HG21	1:A:211:LYS:HD2	1.85	0.58
1:A:190:TYR:CE2	1:A:192:PRO:HG3	2.38	0.58
1:A:109:ALA:O	1:A:113:LEU:HD13	2.02	0.58
4:R:136:THR:HG23	4:R:221:ARG:HH21	1.68	0.58
1:A:364:THR:HG22	1:A:365:CYS:N	2.19	0.58
2:B:294:CYS:HB3	2:B:308:LEU:HB2	1.85	0.58
1:A:291:LEU:HD12	1:A:363:PHE:CE1	2.39	0.58
5:N:51:ILE:HG13	5:N:58:ILE:HG12	1.86	0.58
4:R:289:PHE:HB2	4:R:311:LEU:HB3	1.85	0.57
1:A:93:GLN:HA	1:A:96:LYS:HE2	1.86	0.57
2:B:309:ALA:O	2:B:339:TRP:HH2	1.88	0.57
1:A:338:LYS:HG2	1:A:363:PHE:CE1	2.39	0.57
1:A:9:THR:HG22	1:A:13:ARG:NH2	2.19	0.57
1:A:293:LYS:HD3	1:A:296:LEU:HD23	1.86	0.57
4:R:167:LEU:HB2	4:R:168:PRO:HD3	1.86	0.57
5:N:60:TYR:HB3	5:N:64:VAL:HG23	1.86	0.57
4:R:54:VAL:HG22	4:R:336:PHE:HZ	1.69	0.57
4:R:83:GLY:HA2	4:R:87:VAL:HB	1.87	0.57
4:R:199:TYR:O	4:R:203:SER:CB	2.53	0.57
2:B:168:LEU:HB3	2:B:178:THR:CG2	2.35	0.56
1:A:254:ASN:O	1:A:255:MET:HB3	2.04	0.56
2:B:71:VAL:HG22	2:B:81:ILE:HG12	1.87	0.56
2:B:235:PHE:CD1	2:B:236:PRO:HD2	2.40	0.56
1:A:183:ASP:HA	1:A:186:LYS:HG3	1.86	0.56
4:R:292:VAL:HG13	4:R:303:ILE:HD13	1.86	0.56
1:A:103:ILE:HG13	1:A:104:GLU:N	2.20	0.56
2:B:30:LEU:HD12	2:B:261:LEU:HD13	1.87	0.56
5:N:109:ASP:OD1	5:N:110:VAL:HG23	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:121:ILE:HG23	4:R:282:PHE:CE1	2.41	0.56
4:R:1117:SER:HA	4:R:1120:MET:HE3	1.86	0.56
2:B:152:LEU:HD22	2:B:196:THR:HB	1.87	0.56
1:A:280:ARG:HA	1:A:283:ARG:NE	2.20	0.56
1:A:248:VAL:HG11	1:A:269:ALA:HB1	1.88	0.56
5:N:32:TYR:CZ	5:N:102:PRO:HG3	2.41	0.56
4:R:1033:LEU:HD13	4:R:1046:LEU:HB2	1.88	0.56
4:R:199:TYR:O	4:R:203:SER:HB2	2.06	0.55
2:B:207:SER:CB	2:B:223:THR:HG22	2.37	0.55
2:B:158:VAL:HG11	2:B:192:LEU:HD21	1.89	0.55
4:R:1005:GLU:O	4:R:1009:ILE:HG13	2.06	0.55
1:A:213:GLN:NE2	1:A:216:LYS:HA	2.20	0.55
4:R:239:ARG:N	4:R:239:ARG:HD2	2.22	0.55
5:N:31:ASN:OD1	5:N:103:PHE:CE1	2.59	0.55
4:R:275:LEU:HA	4:R:278:ILE:HG22	1.87	0.55
1:A:102:ALA:HB2	1:A:172:ILE:HD11	1.89	0.55
1:A:44:LEU:HD22	1:A:238:PHE:CD1	2.42	0.54
5:N:73:ASP:HB3	5:N:78:THR:HG23	1.88	0.54
2:B:211:TRP:CZ3	2:B:218:CYS:HB2	2.42	0.54
4:R:1145:ARG:O	4:R:1149:VAL:HG23	2.07	0.54
2:B:318:LEU:HD23	2:B:319:GLY:N	2.23	0.54
3:G:6:THR:O	3:G:9:ILE:HG13	2.08	0.54
1:A:229:ASP:OD2	5:N:111:THR:HG23	2.07	0.54
1:A:127:ARG:HG2	1:A:149:HIS:CD2	2.43	0.54
1:A:294:GLN:CB	1:A:366:ALA:HB2	2.36	0.54
4:R:184:CYS:SG	4:R:190:CYS:O	2.66	0.54
1:A:103:ILE:CG2	1:A:179:LEU:HD21	2.33	0.54
1:A:291:LEU:HD12	1:A:363:PHE:HE1	1.72	0.54
1:A:371:ASN:HA	1:A:374:ARG:HG3	1.89	0.54
1:A:119:LEU:HD12	1:A:119:LEU:H	1.73	0.53
1:A:165:ARG:O	1:A:168:GLU:HG2	2.07	0.53
1:A:338:LYS:HG2	1:A:363:PHE:CZ	2.43	0.53
4:R:84:LEU:O	4:R:88:PRO:HG2	2.08	0.53
2:B:278:PHE:CE1	2:B:285:LEU:HD13	2.43	0.53
4:R:309:ILE:HG12	6:R:366:P0G:HAH	1.89	0.53
4:R:325:ILE:HG22	4:R:325:ILE:O	2.09	0.53
1:A:202:VAL:HB	1:A:211:LYS:HG3	1.89	0.53
4:R:163:LEU:HG	4:R:167:LEU:HD12	1.89	0.53
1:A:115:PRO:HD2	1:A:165:ARG:HH22	1.73	0.53
2:B:148:CYS:SG	2:B:190:LEU:HG	2.49	0.53
4:R:1002:ASN:CG	4:R:1003:ILE:H	2.12	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:71:VAL:HG23	2:B:105:TYR:CD2	2.44	0.53
1:A:318:TYR:CE2	1:A:340:PHE:HB2	2.44	0.53
2:B:314:ARG:HD3	2:B:332:TRP:CE2	2.44	0.53
4:R:1034:THR:HG22	4:R:1042:ALA:HB2	1.89	0.53
4:R:51:ASN:OD1	4:R:79:ASP:HB2	2.08	0.52
4:R:174:TYR:HA	4:R:196:ASN:HD21	1.75	0.52
1:A:111:SER:HA	1:A:116:PRO:HB3	1.92	0.52
2:B:289:TYR:HB2	2:B:291:ASP:OD2	2.09	0.52
2:B:60:ALA:HA	2:B:317:CYS:HB3	1.92	0.52
1:A:244:ILE:HG13	1:A:285:ILE:HG21	1.92	0.52
4:R:291:ILE:O	4:R:295:VAL:HG23	2.09	0.52
2:B:226:GLU:O	5:N:98:ARG:NH1	2.37	0.52
5:N:102:PRO:HB2	5:N:103:PHE:CD2	2.44	0.52
4:R:1024:TYR:CE2	4:R:1035:LYS:HD2	2.45	0.52
5:N:1:GLN:HG2	5:N:2:VAL:N	2.21	0.52
2:B:168:LEU:HB3	2:B:178:THR:HG23	1.91	0.52
4:R:136:THR:C	4:R:138:PRO:HD3	2.30	0.52
2:B:148:CYS:HB3	2:B:160:SER:OG	2.09	0.52
4:R:218:VAL:HG12	4:R:219:TYR:N	2.25	0.51
4:R:1091:LEU:HD11	4:R:1121:LEU:HD13	1.91	0.51
5:N:17:SER:HA	5:N:86:LEU:HD13	1.91	0.51
2:B:31:SER:HB2	2:B:262:MET:CE	2.40	0.51
2:B:274:THR:HG22	2:B:314:ARG:HH21	1.74	0.51
4:R:1120:MET:HG2	4:R:1125:ARG:HD2	1.93	0.51
4:R:212:LEU:O	4:R:216:VAL:HG23	2.09	0.51
4:R:228:ARG:O	4:R:232:LYS:HG3	2.09	0.51
1:A:209:GLU:HG3	1:A:220:HIS:NE2	2.25	0.51
1:A:113:LEU:HG	1:A:168:GLU:OE2	2.11	0.51
4:R:296:HIS:HB2	4:R:303:ILE:HD12	1.93	0.51
2:B:72:SER:O	2:B:79:LEU:HD12	2.11	0.51
4:R:1026:THR:HG23	4:R:1031:HIS:O	2.09	0.51
2:B:235:PHE:HD2	2:B:237:ASN:OD1	1.92	0.51
4:R:277:ILE:HD12	4:R:278:ILE:N	2.26	0.51
2:B:61:MET:HG3	2:B:317:CYS:SG	2.50	0.51
1:A:12:GLN:O	1:A:16:GLU:HG2	2.10	0.51
4:R:1108:GLU:HG3	4:R:1109:THR:N	2.26	0.51
2:B:113:ALA:HB2	2:B:123:ILE:HD13	1.93	0.51
4:R:142:GLN:NE2	4:R:142:GLN:H	2.09	0.50
1:A:364:THR:O	1:A:365:CYS:HB3	2.11	0.50
4:R:1074:ALA:O	4:R:1078:ILE:HG13	2.11	0.50
2:B:126:LEU:HA	2:B:133:VAL:HG12	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:322:ASN:HB2	4:R:323:PRO:HD3	1.92	0.50
5:N:1:GLN:CG	5:N:2:VAL:H	2.22	0.50
4:R:1124:LYS:HA	4:R:1126:TRP:CH2	2.47	0.50
1:A:17:LYS:O	1:A:21:GLU:HG2	2.11	0.50
1:A:293:LYS:H	1:A:365:CYS:HB3	1.77	0.49
1:A:202:VAL:HG21	1:A:211:LYS:CD	2.41	0.49
4:R:1094:VAL:HG11	4:R:1156:GLY:O	2.11	0.49
2:B:141:GLY:N	2:B:169:TRP:CH2	2.79	0.49
1:A:231:ARG:HG2	1:A:234:TRP:CZ2	2.47	0.49
4:R:329:SER:HB3	4:R:332:PHE:HD1	1.77	0.49
1:A:91:LYS:O	1:A:95:ILE:HG12	2.12	0.49
4:R:110:THR:O	4:R:113:ASP:HB3	2.13	0.49
4:R:293:ASN:ND2	6:R:366:P0G:HAMA	2.20	0.49
2:B:51:LEU:HB3	2:B:82:TRP:CE3	2.48	0.49
1:A:277:TRP:CE2	1:A:357:HIS:CE1	3.01	0.49
1:A:202:VAL:CG2	1:A:211:LYS:CD	2.87	0.48
5:N:67:ARG:O	5:N:68:PHE:HD1	1.96	0.48
2:B:271:CYS:HB2	2:B:290:ASP:HB2	1.93	0.48
5:N:45:LEU:H	5:N:45:LEU:HD12	1.79	0.48
5:N:6:GLU:OE2	5:N:119:GLY:HA3	2.14	0.48
5:N:1:GLN:O	5:N:26:GLY:CA	2.62	0.48
4:R:327:CYS:O	4:R:332:PHE:HB3	2.12	0.48
2:B:57:LYS:CE	2:B:75:GLN:HG3	2.41	0.48
5:N:39:GLN:HG3	5:N:43:LYS:O	2.14	0.48
4:R:121:ILE:HG23	4:R:282:PHE:HE1	1.79	0.48
2:B:289:TYR:HE1	2:B:295:ASN:HB2	1.79	0.48
4:R:197:GLN:HG3	4:R:297:VAL:CG1	2.44	0.48
1:A:289:LEU:HD21	1:A:291:LEU:HD21	1.96	0.48
4:R:169:ILE:HA	4:R:174:TYR:CE2	2.48	0.48
1:A:379:CYS:O	1:A:383:ILE:HG22	2.13	0.48
4:R:147:LYS:HD3	4:R:151:ARG:NH2	2.29	0.48
2:B:53:GLY:N	2:B:82:TRP:CH2	2.72	0.48
2:B:93:ILE:HG21	2:B:124:TYR:CD2	2.49	0.48
2:B:292:PHE:CD1	2:B:292:PHE:N	2.82	0.48
5:N:36:TRP:NE1	5:N:81:LEU:HB2	2.29	0.48
5:N:28:THR:OG1	5:N:31:ASN:ND2	2.47	0.47
2:B:207:SER:HB2	2:B:223:THR:HG22	1.96	0.47
1:A:119:LEU:N	1:A:119:LEU:HD12	2.29	0.47
4:R:1096:ARG:O	4:R:1100:ILE:HG13	2.14	0.47
4:R:1125:ARG:HG3	4:R:1125:ARG:O	2.14	0.47
2:B:289:TYR:CE1	2:B:295:ASN:HB2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:153:ILE:O	4:R:157:VAL:HG23	2.14	0.47
2:B:192:LEU:HD23	2:B:199:PHE:HB3	1.96	0.47
4:R:100:THR:HG22	4:R:100:THR:O	2.14	0.47
1:A:114:VAL:CB	1:A:115:PRO:HD3	2.28	0.47
4:R:79:ASP:O	4:R:82:MET:HB3	2.13	0.47
2:B:205:ASP:HA	5:N:116:ALA:CB	2.44	0.47
1:A:194:ASP:O	1:A:198:LEU:HD13	2.14	0.47
2:B:315:VAL:O	2:B:315:VAL:HG23	2.15	0.47
4:R:136:THR:O	4:R:138:PRO:HD3	2.15	0.47
1:A:99:LEU:HD12	1:A:185:ILE:HD12	1.96	0.47
2:B:164:THR:HG22	2:B:164:THR:O	2.14	0.47
1:A:50:GLU:O	1:A:293:LYS:HE3	2.14	0.47
1:A:202:VAL:HB	1:A:211:LYS:HG2	1.97	0.47
1:A:241:VAL:O	1:A:285:ILE:HD12	2.15	0.47
4:R:204:SER:HB3	4:R:209:TYR:HE2	1.79	0.47
2:B:243:THR:O	2:B:250:CYS:HA	2.14	0.47
4:R:132:TYR:O	4:R:136:THR:HB	2.15	0.47
5:N:115:TYR:HE2	5:N:117:TYR:O	1.98	0.47
2:B:309:ALA:O	2:B:339:TRP:CH2	2.68	0.47
1:A:119:LEU:HD21	1:A:128:VAL:HG21	1.97	0.47
2:B:291:ASP:O	2:B:292:PHE:HB2	2.15	0.47
4:R:1046:LEU:HD11	4:R:1058:ILE:HG21	1.97	0.47
5:N:13:GLN:HG3	5:N:14:PRO:HD2	1.98	0.46
2:B:84:SER:HB2	3:G:61:PHE:HE2	1.80	0.46
4:R:289:PHE:HA	4:R:311:LEU:HD13	1.96	0.46
2:B:205:ASP:HA	5:N:116:ALA:HB2	1.96	0.46
4:R:31:VAL:O	4:R:31:VAL:HG12	2.15	0.46
1:A:392:GLU:HB2	4:R:274:THR:HG21	1.97	0.46
4:R:140:LYS:N	4:R:140:LYS:HD2	2.29	0.46
4:R:1119:ARG:O	4:R:1122:GLN:HG2	2.16	0.46
4:R:205:ILE:HA	4:R:209:TYR:HB2	1.96	0.46
4:R:313:TRP:O	4:R:316:TYR:HB2	2.15	0.46
2:B:325:MET:O	2:B:340:ASN:ND2	2.47	0.46
1:A:42:ARG:C	1:A:43:LEU:HD12	2.36	0.46
2:B:58:ILE:HG13	2:B:334:SER:HA	1.98	0.46
2:B:96:ARG:HA	2:B:96:ARG:HD3	1.78	0.46
2:B:181:THR:O	2:B:211:TRP:HH2	1.98	0.46
2:B:235:PHE:CG	2:B:236:PRO:HD2	2.51	0.46
4:R:1016:LYS:HG3	4:R:1057:VAL:HG22	1.98	0.46
5:N:13:GLN:CG	5:N:14:PRO:HD2	2.46	0.46
2:B:340:ASN:HD22	2:B:340:ASN:HA	1.56	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:122:SER:HB3	2:B:124:TYR:CE1	2.51	0.46
4:R:1098:ALA:O	4:R:1102:MET:HG3	2.16	0.46
4:R:1127:ASP:O	4:R:1131:VAL:HG23	2.16	0.46
2:B:53:GLY:HA3	2:B:89:LYS:HE3	1.98	0.45
1:A:384:GLN:O	1:A:388:LEU:HB2	2.17	0.45
4:R:237:GLU:HG2	4:R:239:ARG:H	1.81	0.45
2:B:16:ASN:HA	2:B:19:ARG:NH1	2.32	0.45
4:R:1046:LEU:O	4:R:1050:ILE:HG12	2.16	0.45
2:B:297:TRP:HA	2:B:304:ARG:HA	1.98	0.45
4:R:1031:HIS:CE1	4:R:1066:LEU:HD22	2.51	0.45
1:A:166:SER:HA	1:A:169:TYR:CZ	2.50	0.45
5:N:109:ASP:CG	5:N:110:VAL:HG23	2.37	0.45
5:N:45:LEU:N	5:N:45:LEU:HD12	2.32	0.45
5:N:48:VAL:HG13	5:N:64:VAL:HG11	1.97	0.45
5:N:76:LYS:HB3	5:N:78:THR:HG22	1.98	0.45
2:B:57:LYS:HB2	2:B:332:TRP:HA	1.99	0.45
5:N:101:ALA:HA	5:N:102:PRO:HD3	1.85	0.45
2:B:30:LEU:HD23	2:B:30:LEU:O	2.17	0.45
1:A:140:PHE:CE2	1:A:142:PHE:HA	2.51	0.45
2:B:86:THR:O	2:B:87:THR:HG22	2.17	0.45
2:B:215:GLU:HB3	2:B:217:MET:HG2	1.98	0.45
4:R:110:THR:HG23	6:R:366:P0G:HAB	1.98	0.45
4:R:169:ILE:HA	4:R:174:TYR:CD2	2.51	0.45
5:N:67:ARG:C	5:N:68:PHE:HD1	2.21	0.45
4:R:124:LEU:HD22	4:R:326:TYR:OH	2.17	0.45
2:B:59:TYR:O	2:B:60:ALA:HB2	2.17	0.45
4:R:68:THR:HG23	4:R:130:ASP:OD2	2.17	0.45
2:B:85:TYR:HE1	3:G:60:PRO:HB2	1.81	0.44
2:B:48:ARG:HE	2:B:340:ASN:HB2	1.82	0.44
4:R:1111:VAL:O	4:R:1114:PHE:HB2	2.17	0.44
4:R:125:CYS:HB2	4:R:211:PRO:HB3	1.97	0.44
4:R:1033:LEU:CD1	4:R:1046:LEU:HB2	2.47	0.44
2:B:326:ALA:HA	2:B:340:ASN:ND2	2.33	0.44
1:A:43:LEU:HD23	1:A:245:ILE:HD11	1.98	0.44
4:R:1027:ILE:HG12	4:R:1028:GLY:N	2.32	0.44
4:R:109:TRP:HZ3	6:R:366:P0G:HAAA	1.83	0.44
4:R:126:VAL:HG21	4:R:157:VAL:HG21	1.98	0.44
2:B:320:VAL:HG22	2:B:327:VAL:HG22	2.00	0.44
1:A:270:LEU:HD23	1:A:345:PHE:CE2	2.52	0.44
1:A:364:THR:CG2	1:A:365:CYS:N	2.81	0.44
4:R:169:ILE:HG23	4:R:174:TYR:CE2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:THR:O	1:A:221:MET:N	2.48	0.44
5:N:38:ARG:HA	5:N:93:VAL:O	2.17	0.44
2:B:230:ASN:HB2	2:B:273:ILE:O	2.18	0.44
1:A:308:ILE:HG22	1:A:312:PHE:HB2	1.99	0.44
2:B:262:MET:HG3	2:B:263:THR:N	2.32	0.44
2:B:53:GLY:H	2:B:82:TRP:HH2	1.60	0.44
3:G:9:ILE:HD12	3:G:9:ILE:C	2.38	0.44
2:B:207:SER:HB3	2:B:223:THR:HG22	2.00	0.44
4:R:1004:PHE:O	4:R:1008:ARG:HB2	2.17	0.44
2:B:316:SER:H	2:B:331:SER:HA	1.81	0.44
4:R:174:TYR:HA	4:R:196:ASN:ND2	2.32	0.44
2:B:184:THR:O	2:B:184:THR:HG22	2.17	0.44
2:B:46:ARG:HD2	2:B:48:ARG:NH2	2.33	0.44
2:B:163:ASP:O	2:B:164:THR:CB	2.64	0.44
4:R:44:VAL:HG22	4:R:87:VAL:HG12	1.99	0.44
1:A:127:ARG:O	1:A:131:ILE:HG12	2.17	0.44
1:A:232:ARG:HB3	5:N:108:PHE:HB2	1.99	0.44
1:A:159:VAL:O	1:A:162:CYS:HB3	2.18	0.44
1:A:360:TYR:CE2	1:A:382:ILE:HG12	2.52	0.44
2:B:149:CYS:CB	2:B:159:THR:HG22	2.47	0.44
5:N:111:THR:CG2	5:N:114:THR:HG23	2.47	0.44
5:N:6:GLU:OE1	5:N:119:GLY:HA3	2.17	0.44
3:G:28:ILE:HD12	3:G:32:LYS:HD3	2.00	0.44
5:N:115:TYR:CE2	5:N:117:TYR:O	2.71	0.43
1:A:238:PHE:HB3	1:A:241:VAL:HB	1.99	0.43
5:N:34:MET:SD	5:N:98:ARG:HB3	2.58	0.43
5:N:43:LYS:HA	5:N:43:LYS:HD2	1.74	0.43
4:R:288:PRO:O	4:R:311:LEU:HD13	2.17	0.43
4:R:121:ILE:HD12	4:R:282:PHE:CE1	2.53	0.43
4:R:225:GLU:OE1	4:R:225:GLU:HA	2.18	0.43
4:R:40:MET:HE2	4:R:94:ILE:HD12	2.00	0.43
1:A:94:ASP:O	1:A:97:ASN:HB2	2.19	0.43
2:B:296:VAL:O	2:B:305:ALA:N	2.51	0.43
2:B:291:ASP:O	2:B:292:PHE:CB	2.67	0.43
2:B:253:PHE:CE1	2:B:260:GLU:HB3	2.53	0.43
2:B:331:SER:OG	2:B:332:TRP:N	2.52	0.43
2:B:150:ARG:HG2	2:B:190:LEU:HD11	2.00	0.43
1:A:295:ASP:O	1:A:299:GLU:HG3	2.18	0.43
2:B:314:ARG:HD3	2:B:332:TRP:CZ2	2.54	0.43
5:N:102:PRO:HB2	5:N:103:PHE:CE2	2.54	0.43
3:G:12:ALA:O	3:G:16:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:62:GLU:H	4:R:62:GLU:CD	2.22	0.43
2:B:72:SER:HB3	2:B:336:LEU:HD11	2.01	0.43
4:R:155:LEU:O	4:R:159:ILE:HG13	2.19	0.43
4:R:38:ILE:O	4:R:42:LEU:HG	2.19	0.43
4:R:303:ILE:HG22	4:R:304:ARG:O	2.19	0.43
1:A:235:ILE:HG12	1:A:282:LEU:HD11	2.01	0.43
4:R:94:ILE:HD11	4:R:313:TRP:NE1	2.33	0.43
1:A:179:LEU:O	1:A:182:ILE:HG13	2.19	0.42
6:R:366:P0G:HABB	6:R:366:P0G:HANA	1.85	0.42
4:R:1059:THR:HG22	4:R:1060:LYS:N	2.34	0.42
4:R:53:LEU:O	4:R:57:ALA:HB2	2.19	0.42
2:B:170:ASP:HB3	2:B:173:THR:OG1	2.19	0.42
4:R:90:GLY:O	4:R:94:ILE:HG13	2.18	0.42
4:R:1072:ASP:O	4:R:1076:ARG:HG3	2.19	0.42
2:B:118:ASP:OD2	2:B:120:ILE:HB	2.19	0.42
4:R:91:ALA:HB2	4:R:313:TRP:HH2	1.84	0.42
4:R:30:GLU:HB2	4:R:32:TRP:H	1.84	0.42
5:N:109:ASP:HB3	5:N:115:TYR:CZ	2.55	0.42
2:B:124:TYR:HD1	2:B:135:VAL:HA	1.80	0.42
4:R:1066:LEU:HA	4:R:1066:LEU:HD23	1.93	0.42
1:A:96:LYS:HD2	1:A:147:TYR:OH	2.19	0.42
1:A:201:ARG:O	1:A:202:VAL:C	2.58	0.42
2:B:149:CYS:HB2	2:B:159:THR:HG22	2.00	0.42
1:A:230:GLU:HG2	5:N:111:THR:OG1	2.19	0.42
1:A:250:SER:OG	1:A:291:LEU:HB3	2.19	0.42
1:A:340:PHE:O	1:A:344:GLU:HG2	2.19	0.42
5:N:29:PHE:CZ	5:N:34:MET:HG3	2.54	0.42
4:R:52:VAL:HG23	4:R:80:LEU:HD11	2.01	0.42
5:N:109:ASP:CB	5:N:115:TYR:CZ	3.03	0.42
5:N:83:MET:HB3	5:N:86:LEU:HD11	2.02	0.42
1:A:231:ARG:O	1:A:235:ILE:HB	2.20	0.42
4:R:307:VAL:O	4:R:311:LEU:HG	2.19	0.42
2:B:112:VAL:HG23	2:B:126:LEU:HD11	2.02	0.42
1:A:233:LYS:HG3	2:B:188:MET:SD	2.59	0.42
1:A:115:PRO:O	1:A:116:PRO:C	2.58	0.41
1:A:28:LYS:HE2	1:A:32:LYS:HE3	2.01	0.41
4:R:71:PHE:HA	4:R:154:ILE:HD11	2.01	0.41
4:R:71:PHE:HD1	4:R:154:ILE:HD11	1.84	0.41
2:B:118:ASP:O	2:B:120:ILE:HG13	2.20	0.41
4:R:1071:VAL:O	4:R:1075:VAL:HG23	2.20	0.41
1:A:238:PHE:CD1	1:A:238:PHE:N	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:317:CYS:SG	2:B:330:GLY:HA3	2.60	0.41
2:B:204:CYS:HA	2:B:228:ASP:HB3	2.03	0.41
1:A:252:SER:HB2	1:A:265:ARG:HB2	2.03	0.41
1:A:229:ASP:OD1	5:N:111:THR:HG23	2.19	0.41
2:B:125:ASN:O	2:B:133:VAL:HG12	2.20	0.41
1:A:246:PHE:CE1	1:A:273:PHE:HB2	2.51	0.41
1:A:17:LYS:O	1:A:20:ARG:HB3	2.21	0.41
4:R:197:GLN:O	4:R:201:ILE:HG13	2.21	0.41
1:A:181:LYS:HG2	1:A:184:VAL:HB	2.03	0.41
4:R:71:PHE:CD1	4:R:154:ILE:HD11	2.56	0.41
4:R:86:VAL:CG2	4:R:112:ILE:HG23	2.48	0.41
4:R:1093:ALA:O	4:R:1096:ARG:HB2	2.20	0.41
4:R:1107:GLY:O	4:R:1111:VAL:HG23	2.19	0.41
4:R:1036:SER:HA	4:R:1037:PRO:HD3	1.91	0.41
5:N:83:MET:CB	5:N:86:LEU:HD11	2.51	0.41
5:N:4:LEU:HD21	5:N:27:PHE:CE2	2.56	0.41
4:R:305:LYS:HA	4:R:305:LYS:HD3	1.85	0.41
2:B:314:ARG:HD3	2:B:332:TRP:CD2	2.56	0.41
4:R:210:VAL:O	4:R:214:ILE:HG13	2.21	0.41
4:R:87:VAL:HB	4:R:88:PRO:HD3	2.02	0.41
2:B:271:CYS:SG	2:B:289:TYR:HB3	2.61	0.41
2:B:326:ALA:HA	2:B:340:ASN:HD21	1.84	0.41
1:A:99:LEU:HD23	1:A:99:LEU:C	2.42	0.40
2:B:71:VAL:HG23	2:B:105:TYR:CE2	2.56	0.40
2:B:192:LEU:HD23	2:B:199:PHE:CB	2.51	0.40
4:R:1010:ASP:HB3	4:R:1145:ARG:NE	2.37	0.40
4:R:1106:MET:HE1	4:R:1138:TRP:CD1	2.56	0.40
2:B:308:LEU:C	2:B:309:ALA:O	2.57	0.40
2:B:339:TRP:CD1	2:B:339:TRP:N	2.89	0.40
4:R:295:VAL:HG12	4:R:303:ILE:HD11	2.03	0.40
4:R:90:GLY:O	4:R:93:HIS:HB3	2.22	0.40
2:B:48:ARG:CG	2:B:340:ASN:HB2	2.51	0.40
5:N:53:GLN:NE2	5:N:104:THR:H	2.19	0.40
4:R:1085:LYS:N	4:R:1086:PRO:HD2	2.36	0.40
2:B:11:ALA:HB2	3:G:16:VAL:HG22	2.03	0.40
2:B:181:THR:O	2:B:211:TRP:CH2	2.74	0.40
5:N:109:ASP:O	5:N:110:VAL:HB	2.21	0.40
2:B:48:ARG:HG3	2:B:340:ASN:HB2	2.03	0.40
4:R:1077:GLY:HA2	4:R:1080:ARG:HG2	2.02	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	341/380 (90%)	328 (96%)	13 (4%)	0	100	100
2	B	338/351 (96%)	323 (96%)	15 (4%)	0	100	100
3	G	56/68 (82%)	54 (96%)	2 (4%)	0	100	100
4	R	435/514 (85%)	407 (94%)	28 (6%)	0	100	100
5	N	126/138 (91%)	117 (93%)	8 (6%)	1 (1%)	24	69
All	All	1296/1451 (89%)	1229 (95%)	66 (5%)	1 (0%)	56	91

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	N	110	VAL

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	296/342 (86%)	294 (99%)	2 (1%)	88	97
2	B	278/293 (95%)	273 (98%)	5 (2%)	66	89
3	G	46/56 (82%)	46 (100%)	0	100	100
4	R	358/441 (81%)	352 (98%)	6 (2%)	68	90
5	N	104/115 (90%)	102 (98%)	2 (2%)	65	89
All	All	1082/1247 (87%)	1067 (99%)	15 (1%)	74	92

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	238	PHE
1	A	357	HIS
2	B	33	ILE
2	B	59	TYR
2	B	105	TYR
2	B	234	PHE
2	B	340	ASN
4	R	1104	PHE
4	R	106	CYS
4	R	142	GLN
4	R	312	ASN
4	R	332	PHE
4	R	336	PHE
5	N	96	CYS
5	N	98	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	98	ASN
1	A	292	ASN
1	A	357	HIS
2	B	9	GLN
2	B	220	GLN
2	B	230	ASN
2	B	239	ASN
2	B	268	ASN
2	B	293	ASN
2	B	340	ASN
3	G	5	ASN
4	R	1132	ASN
4	R	142	GLN
4	R	170	GLN
4	R	224	GLN
4	R	293	ASN
5	N	1	GLN
5	N	13	GLN
5	N	31	ASN
5	N	77	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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$
6	P0G	R	366	-	28,29,29	1.27	3 (10%)	36,42,42	1.48	4 (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
6	P0G	R	366	-	-	0/15/24/24	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	R	366	P0G	CAW-CAZ	-4.66	1.48	1.52
6	R	366	P0G	CBA-NAP	-3.44	1.45	1.49
6	R	366	P0G	CAX-NAQ	-2.23	1.35	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	R	366	P0G	OAR-CAY-CAX	-3.21	117.44	121.06
6	R	366	P0G	CAX-NAQ-CAS	-2.88	120.34	124.57
6	R	366	P0G	OAR-CAY-CAW	3.41	121.75	116.61
6	R	366	P0G	CAN-NAP-CBA	5.34	122.01	116.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	R	366	P0G	7	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	349/380 (91%)	0.35	24 (6%) 20 11	32, 93, 156, 191	0
2	B	340/351 (96%)	-0.12	2 (0%) 90 84	29, 59, 101, 162	0
3	G	58/68 (85%)	-0.18	1 (1%) 73 60	36, 80, 132, 146	0
4	R	443/514 (86%)	0.52	46 (10%) 8 5	56, 122, 175, 214	0
5	N	128/138 (92%)	-0.20	0 100 100	31, 59, 93, 115	0
All	All	1318/1451 (90%)	0.21	73 (5%) 29 16	29, 89, 158, 214	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	R	192	ASP	5.5
1	A	147	TYR	4.8
1	A	136	ASN	4.2
4	R	144	LEU	4.1
4	R	294	ILE	4.1
2	B	130	GLU	4.0
1	A	185	ILE	3.9
4	R	201	ILE	3.9
4	R	197	GLN	3.9
4	R	105	TRP	3.8
4	R	1094	VAL	3.8
4	R	1017	ILE	3.8
4	R	1014	ARG	3.8
1	A	90	THR	3.7
4	R	1037	PRO	3.7
4	R	181	ALA	3.7
3	G	5	ASN	3.6
4	R	298	ILE	3.5
1	A	89	ALA	3.4
4	R	34	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
4	R	183	ASN	3.2
4	R	193	PHE	3.2
1	A	92	VAL	3.2
1	A	93	GLN	3.2
4	R	1033	LEU	3.2
1	A	57	VAL	3.1
2	B	132	ASN	3.1
4	R	184	CYS	3.1
4	R	32	TRP	3.1
1	A	51	SER	3.0
4	R	301	ASN	3.0
1	A	96	LYS	3.0
4	R	1088	TYR	3.0
4	R	57	ALA	3.0
4	R	275	LEU	2.9
1	A	122	PRO	2.9
4	R	202	ALA	2.9
1	A	123	GLU	2.9
1	A	205	SER	2.8
4	R	1025	TYR	2.7
1	A	144	PRO	2.7
1	A	168	GLU	2.6
4	R	274	THR	2.6
1	A	187	GLN	2.6
4	R	91	ALA	2.5
4	R	191	CYS	2.5
4	R	92	ALA	2.4
4	R	1012	GLY	2.4
4	R	39	VAL	2.4
1	A	50	GLU	2.3
4	R	1159	ASP	2.3
1	A	365	CYS	2.3
1	A	368	ASP	2.2
4	R	341	CYS	2.2
1	A	59	GLN	2.2
4	R	1097	ALA	2.2
4	R	1018	TYR	2.2
4	R	309	ILE	2.2
1	A	140	PHE	2.1
4	R	1036	SER	2.1
4	R	172	HIS	2.1
4	R	300	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	95	ILE	2.1
4	R	33	VAL	2.1
4	R	1055	ASN	2.1
4	R	95	LEU	2.1
4	R	185	TYR	2.1
4	R	1053	ASN	2.1
1	A	186	LYS	2.1
4	R	1028	GLY	2.1
4	R	1041	ALA	2.1
4	R	1007	LEU	2.0
1	A	263	THR	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(Å ²)	Q<0.9
6	P0G	R	366	27/27	0.60	0.57	1.88	152,166,192,195	0

6.5 Other polymers [i](#)

There are no such residues in this entry.