



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:12 PM UTC

PDB ID : 3SN6 / pdb_00003sn6
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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

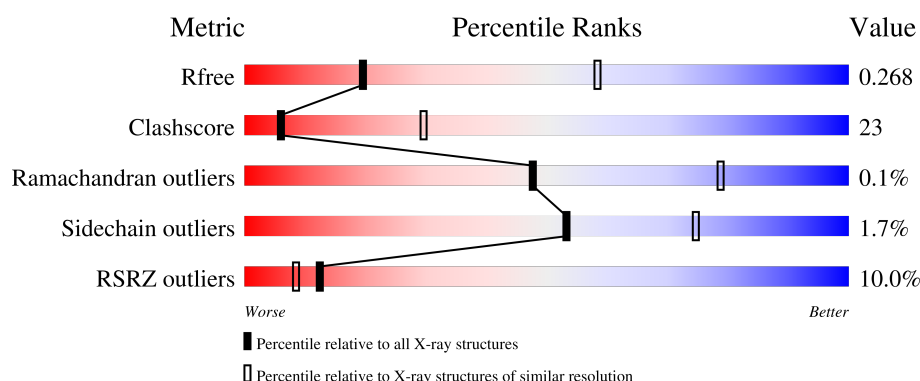
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	380	10% 58% 33% 8%
2	B	351	5% 56% 40% •
3	G	68	6% 65% 21% 15%
4	R	514	13% 47% 38% • 14%
5	N	138	4% 48% 44% • 7%

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 Endolysin,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			

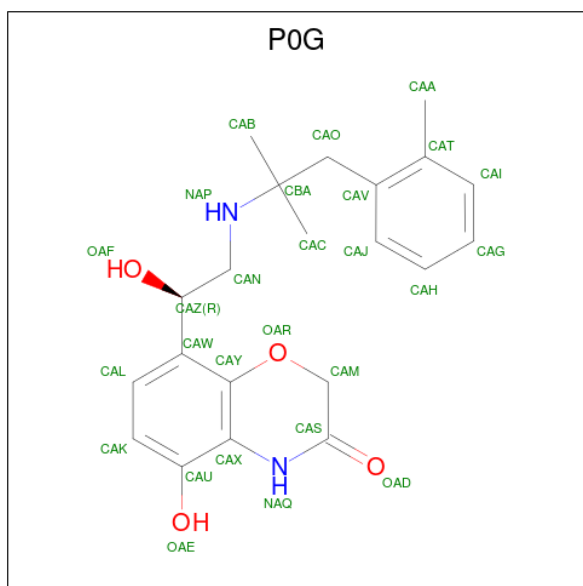
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	987	ASP	-	expression tag	UNP P00720
R	988	TYR	-	expression tag	UNP P00720
R	989	LYS	-	expression tag	UNP P00720
R	990	ASP	-	expression tag	UNP P00720
R	991	ASP	-	expression tag	UNP P00720
R	992	ASP	-	expression tag	UNP P00720
R	993	ASP	-	expression tag	UNP P00720
R	994	ALA	-	expression tag	UNP P00720
R	995	GLU	-	expression tag	UNP P00720
R	996	ASN	-	expression tag	UNP P00720
R	997	LEU	-	expression tag	UNP P00720
R	998	TYR	-	expression tag	UNP P00720
R	999	PHE	-	expression tag	UNP P00720
R	1000	GLN	-	expression tag	UNP P00720
R	1001	GLY	-	expression tag	UNP P00720
R	1012	GLY	ARG	conflict	UNP P00720
R	1054	THR	CYS	engineered mutation	UNP P00720
R	1097	ALA	CYS	engineered mutation	UNP P00720
R	1137	ARG	ILE	conflict	UNP P00720
R	?	ALA	-	linker	UNP P00720
R	?	ALA	-	linker	UNP P00720
R	96	THR	MET	engineered mutation	UNP P07550
R	98	THR	MET	engineered mutation	UNP P07550
R	187	GLU	ASN	engineered mutation	UNP P07550

- 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 (CCD ID: P0G) (formula: $C_{21}H_{26}N_2O_4$).

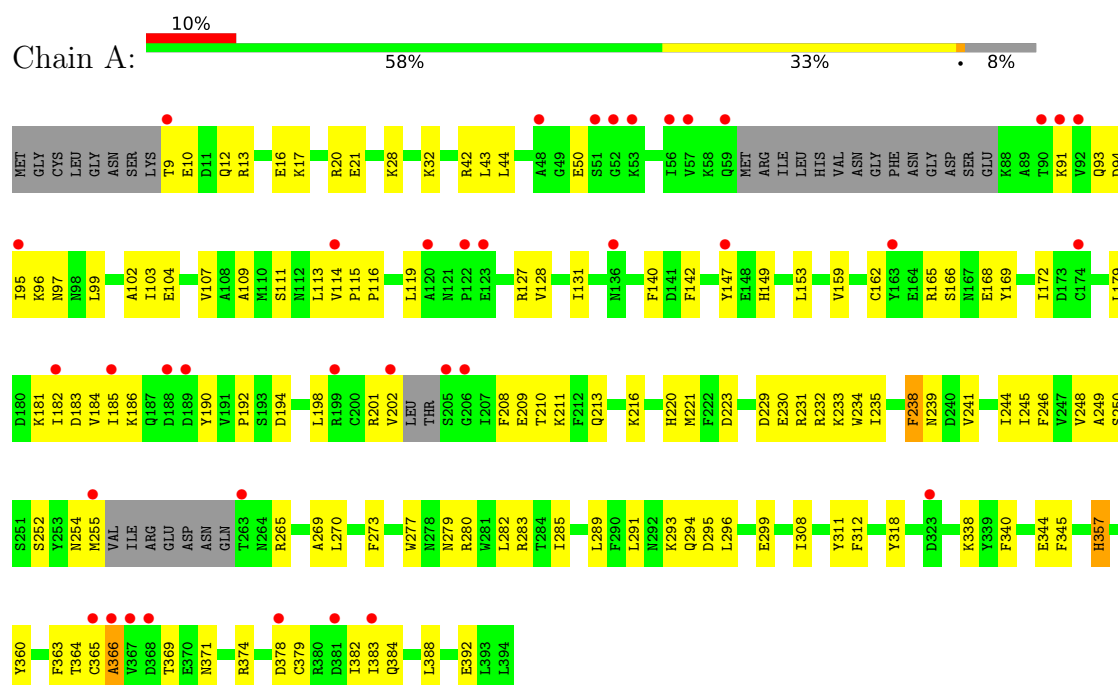


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	R	1	Total	C	N	O	0	0
			27	21	2	4		

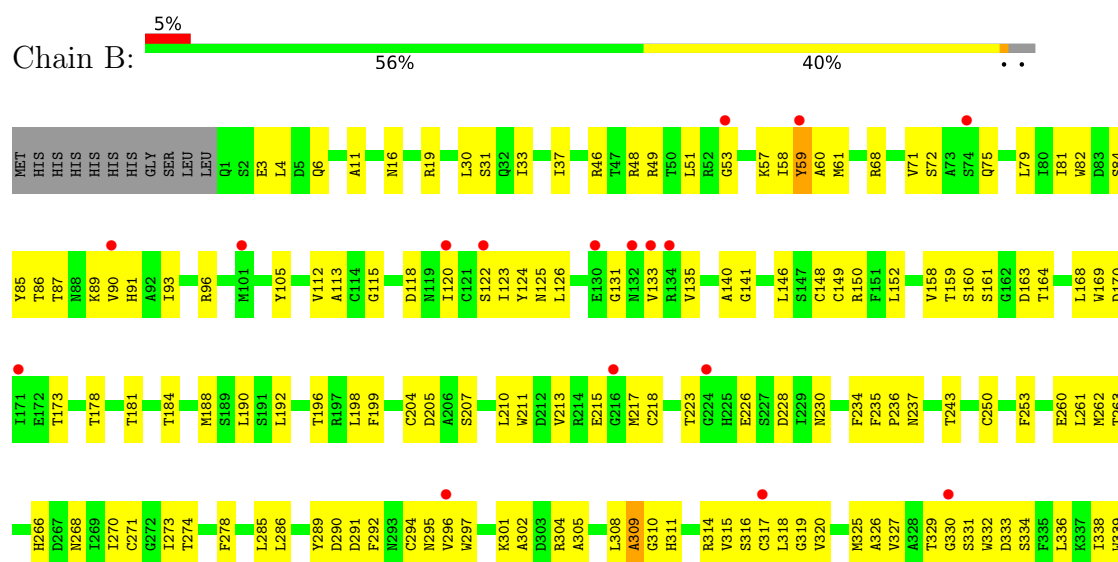
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 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





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 40.67 – 3.20	Depositor EDS
% Data completeness (in resolution range)	92.8 (40.67-3.20) 92.8 (40.67-3.20)	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 0.223 , 0.268	Depositor DCC
R_{free} test set	1558 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	68.1	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 74.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-l	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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.28	0/2870	0.73	3/3879 (0.1%)
2	B	0.29	0/2639	0.77	2/3580 (0.1%)
3	G	0.32	0/444	0.75	0/601
4	R	0.29	0/3502	0.73	2/4762 (0.0%)
5	N	0.34	0/990	0.84	1/1341 (0.1%)
All	All	0.29	0/10445	0.75	8/14163 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	309	ALA	N-CA-C	-8.97	99.07	110.19
2	B	310	GLY	N-CA-C	8.50	121.17	110.29
1	A	366	ALA	N-CA-C	-6.87	103.72	111.07
4	R	106	CYS	N-CA-C	-6.66	104.11	111.36
1	A	369	THR	CB-CA-C	-6.16	107.72	116.34
5	N	113	THR	CB-CA-C	-6.10	108.53	115.79
4	R	31	VAL	N-CA-C	-5.99	106.91	112.83
1	A	10	GLU	CB-CA-C	-5.01	110.77	116.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2814	0	2719	111	0
2	B	2592	0	2485	138	0
3	G	438	0	443	12	0
4	R	3433	0	3429	149	0
5	N	970	0	930	65	0
6	R	27	0	25	8	0
All	All	10274	0	10031	460	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 23.

All (460) 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:1601: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:O	1:A:202:VAL:HG22	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:1601:P0G:HAL	1.72	0.71
4:R:149:LYS:O	4:R:153:ILE:HG12	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ASP:CG	5:N:111:THR:HG23	2.16	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
2:B:120:ILE:HG12	2:B:140:ALA:HB2	1.74	0.69
2:B:304:ARG:O	2:B:304:ARG:HG3	1.92	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
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:9:GLY:HA2	5:N:18:LEU:HD21	1.78	0.65
5:N:110:VAL:HG12	5:N:110:VAL:O	1.96	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
1:A:311:TYR:O	5:N:62:GLY:HA3	1.97	0.64
2:B:141:GLY:N	2:B:169:TRP:HH2	1.96	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:124:TYR:CD1	2:B:135:VAL:HA	2.34	0.62
2:B:266:HIS:CD2	2:B:268:ASN:HB2	2.33	0.62
4:R:1059:THR:HB	4:R:1062:GLU:HG3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:279:ASN:O	1:A:283:ARG:HG3	2.00	0.61
2:B:311:HIS:CD2	2:B:311:HIS:H	2.16	0.61
4:R:109:TRP:O	4:R:112:ILE:HG22	2.00	0.61
1:A:239:ASN:O	1:A:285:ILE:HD11	2.02	0.60
4:R:69:ASN:O	4:R:72:ILE:HG22	2.02	0.60
1:A:246:PHE:HE1	1:A:273:PHE:HB2	1.65	0.60
4:R:36:MET:HB3	4:R:40:MET:HE3	1.84	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
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:190:TYR:CE2	1:A:192:PRO:HG3	2.38	0.58
1:A:202:VAL:HG21	1:A:211:LYS:HD2	1.85	0.58
1:A:109:ALA:O	1:A:113:LEU:HD13	2.02	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
4:R:136:THR:HG23	4:R:221:ARG:HH21	1.68	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:1002:ASN:CG	4:R:1003:ILE:H	2.12	0.57
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
2:B:301:LYS:O	2:B:302:ALA:HB3	2.03	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
1:A:254:ASN:O	1:A:255:MET:HB3	2.04	0.56
2:B:168:LEU:HB3	2:B:178:THR:CG2	2.35	0.56
1:A:183:ASP:HA	1:A:186:LYS:HG3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
4:R:136:THR:C	4:R:138:PRO:HD3	2.30	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
4:R:121:ILE:HG23	4:R:282:PHE:CE1	2.41	0.56
5:N:109:ASP:OD1	5:N:110:VAL:HG23	2.05	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
4:R:1033:LEU:HD13	4:R:1046:LEU:HB2	1.88	0.56
5:N:32:TYR:CZ	5:N:102:PRO:HG3	2.41	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:HD2	4:R:239:ARG:N	2.22	0.55
5:N:31:ASN:OD1	5:N:103:PHE:CE1	2.59	0.55
1:A:102:ALA:HB2	1:A:172:ILE:HD11	1.89	0.55
4:R:275:LEU:HA	4:R:278:ILE:HG22	1.87	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
1:A:127:ARG:HG2	1:A:149:HIS:CD2	2.43	0.54
1:A:229:ASP:OD2	5:N:111:THR:HG23	2.07	0.54
3:G:6:THR:O	3:G:9:ILE:HG13	2.08	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
4:R:215:MET:HB3	4:R:279:MET:HE1	1.89	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
2:B:278:PHE:CE1	2:B:285:LEU:HD13	2.43	0.53
4:R:84:LEU:O	4:R:88:PRO:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:309:ILE:HG12	6:R:1601: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
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:60:ALA:HA	2:B:317:CYS:HB3	1.92	0.52
2:B:289:TYR:HB2	2:B:291:ASP:OD2	2.09	0.52
1:A:244:ILE:HG13	1:A:285:ILE:HG21	1.92	0.52
2:B:226:GLU:O	5:N:98:ARG:NH1	2.37	0.52
4:R:291:ILE:O	4:R:295:VAL:HG23	2.09	0.52
4:R:1024:TYR:CE2	4:R:1035:LYS:HD2	2.45	0.52
5:N:34:MET:HE3	5:N:79:LEU:HD22	1.90	0.52
5:N:102:PRO:HB2	5:N:103:PHE:CD2	2.44	0.52
2:B:168:LEU:HB3	2:B:178:THR:HG23	1.91	0.52
5:N:1:GLN:HG2	5:N:2:VAL:N	2.21	0.52
2:B:148:CYS:HB3	2:B:160:SER:OG	2.09	0.52
4:R:1091:LEU:HD11	4:R:1121:LEU:HD13	1.91	0.51
4:R:218:VAL:HG12	4:R:219:TYR:N	2.25	0.51
5:N:17:SER:HA	5:N:86:LEU:HD13	1.91	0.51
1:A:209:GLU:HG3	1:A:220:HIS:NE2	2.25	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:113:LEU:HG	1:A:168:GLU:OE2	2.11	0.51
2:B:72:SER:O	2:B:79:LEU:HD12	2.11	0.51
4:R:296:HIS:HB2	4:R:303:ILE:HD12	1.93	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ARG:C	1:A:43:LEU:HD12	2.36	0.51
2:B:113:ALA:HB2	2:B:123:ILE:HD13	1.93	0.51
4:R:1108:GLU:HG3	4:R:1109:THR:N	2.26	0.51
4:R:142:GLN:H	4:R:142:GLN:NE2	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
4:R:322:ASN:HB2	4:R:323:PRO:HD3	1.92	0.50
4:R:1124:LYS:HA	4:R:1126:TRP:CH2	2.47	0.50
5:N:1:GLN:CG	5:N:2:VAL:H	2.22	0.50
5:N:109:ASP:CG	5:N:110:VAL:HG23	2.37	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
5:N:67:ARG:C	5:N:68:PHE:HD1	2.21	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
1:A:231:ARG:HG2	1:A:234:TRP:CZ2	2.47	0.49
2:B:31:SER:HB2	2:B:262:MET:HE1	1.93	0.49
2:B:141:GLY:N	2:B:169:TRP:CH2	2.79	0.49
1:A:91:LYS:O	1:A:95:ILE:HG12	2.12	0.49
4:R:329:SER:HB3	4:R:332:PHE:HD1	1.77	0.49
4:R:110:THR:O	4:R:113:ASP:HB3	2.13	0.49
4:R:293:ASN:ND2	6:R:1601:P0G:HAMA	2.20	0.49
2:B:51:LEU:HB3	2:B:82:TRP:CE3	2.48	0.49
3:G:9:ILE:C	3:G:9:ILE:HD12	2.38	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:6:GLU:OE2	5:N:119:GLY:HA3	2.14	0.48
5:N:45:LEU:HD12	5:N:45:LEU:H	1.79	0.48
4:R:327:CYS:O	4:R:332:PHE:HB3	2.12	0.48
5:N:1:GLN:O	5:N:26:GLY:CA	2.62	0.48
5:N:18:LEU:HB3	5:N:83:MET:HE3	1.95	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
2:B:289:TYR:HE1	2:B:295:ASN:HB2	1.79	0.48
4:R:121:ILE:HG23	4:R:282:PHE:HE1	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
1:A:379:CYS:O	1:A:383:ILE:HG22	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:147:LYS:HD3	4:R:151:ARG:NH2	2.29	0.48
4:R:169:ILE:HA	4:R:174:TYR:CE2	2.48	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
1:A:119:LEU:HD12	1:A:119:LEU:N	2.29	0.47
2:B:207:SER:HB2	2:B:223:THR:HG22	1.96	0.47
4:R:1096:ARG:O	4:R:1100:ILE:HG13	2.14	0.47
4:R:62:GLU:CD	4:R:62:GLU:H	2.22	0.47
5:N:28:THR:OG1	5:N:31:ASN:ND2	2.47	0.47
2:B:289:TYR:CE1	2:B:295:ASN:HB2	2.49	0.47
4:R:1125:ARG:O	4:R:1125:ARG:HG3	2.14	0.47
2:B:192:LEU:HD23	2:B:199:PHE:HB3	1.96	0.47
4:R:153:ILE:O	4:R:157:VAL:HG23	2.14	0.47
4:R:1006:MET:HE1	4:R:1098:ALA:HA	1.97	0.47
4:R:1117:SER:HA	4:R:1120:MET:HE3	1.96	0.47
4:R:100:THR:O	4:R:100:THR:HG22	2.14	0.47
1:A:114:VAL:CB	1:A:115:PRO:HD3	2.28	0.47
2:B:205:ASP:HA	5:N:116:ALA:CB	2.44	0.47
4:R:79:ASP:O	4:R:82:MET:HB3	2.13	0.47
1:A:194:ASP:O	1:A:198:LEU:HD13	2.14	0.47
1:A:50:GLU:O	1:A:293:LYS:HE3	2.14	0.47
1:A:99:LEU:HD12	1:A:185:ILE:HD12	1.96	0.47
2:B:164:THR:O	2:B:164:THR:HG22	2.14	0.47
4:R:136:THR:O	4:R:138:PRO:HD3	2.15	0.47
1:A:201:ARG:O	1:A:202:VAL:C	2.58	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
2:B:243:THR:O	2:B:250:CYS:HA	2.14	0.47
4:R:1120:MET:HE1	4:R:1132:ASN:CG	2.39	0.47
4:R:204:SER:HB3	4:R:209:TYR:HE2	1.79	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
2:B:309:ALA:O	2:B:339:TRP:CH2	2.68	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
4:R:1046:LEU:HD11	4:R:1058:ILE:HG21	1.97	0.47
2:B:84:SER:HB2	3:G:61:PHE:HE2	1.80	0.46
5:N:13:GLN:HG3	5:N:14:PRO:HD2	1.98	0.46
1:A:392:GLU:HB2	4:R:274:THR:HG21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:205:ASP:HA	5:N:116:ALA:HB2	1.96	0.46
4:R:31:VAL:HG12	4:R:31:VAL:O	2.15	0.46
4:R:289:PHE:HA	4:R:311:LEU:HD13	1.96	0.46
4:R:1119:ARG:O	4:R:1122:GLN:HG2	2.16	0.46
4:R:140:LYS:HD2	4:R:140:LYS:N	2.29	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
2:B:58:ILE:HG13	2:B:334:SER:HA	1.98	0.46
2:B:96:ARG:HD3	2:B:96:ARG:HA	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
4:R:1106:MET:HB3	4:R:1106:MET:HE2	1.78	0.46
1:A:115:PRO:O	1:A:116:PRO:C	2.58	0.46
2:B:340:ASN:HD22	2:B:340:ASN:HA	1.56	0.46
3:G:35:ALA:HA	3:G:38:MET:HE2	1.98	0.46
5:N:13:GLN:CG	5:N:14:PRO:HD2	2.46	0.46
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
1:A:384:GLN:O	1:A:388:LEU:HB2	2.17	0.45
2:B:16:ASN:HA	2:B:19:ARG:NH1	2.32	0.45
2:B:53:GLY:HA3	2:B:89:LYS:HE3	1.98	0.45
4:R:237:GLU:HG2	4:R:239:ARG:H	1.81	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:45:LEU:HD12	5:N:45:LEU:N	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
1:A:140:PHE:CE2	1:A:142:PHE:HA	2.51	0.45
2:B:30:LEU:O	2:B:30:LEU:HD23	2.17	0.45
2:B:57:LYS:HB2	2:B:332:TRP:HA	1.99	0.45
2:B:86:THR:O	2:B:87:THR:HG22	2.17	0.45
5:N:101:ALA:HA	5:N:102:PRO:HD3	1.85	0.45
2:B:215:GLU:HB3	2:B:217:MET:HG2	1.98	0.45
2:B:315:VAL:HG23	2:B:315:VAL:O	2.15	0.45
2:B:308:LEU:C	2:B:309:ALA:O	2.57	0.45
4:R:110:THR:HG23	6:R:1601:P0G:HAB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:169:ILE:HA	4:R:174:TYR:CD2	2.51	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
4:R:124:LEU:HD22	4:R:326:TYR:OH	2.17	0.45
2:B:48:ARG:HE	2:B:340:ASN:HB2	1.82	0.44
2:B:85:TYR:HE1	3:G:60:PRO:HB2	1.81	0.44
4:R:1106:MET:HE1	4:R:1138:TRP:HB2	1.99	0.44
4:R:1111:VAL:O	4:R:1114:PHE:HB2	2.17	0.44
1:A:43:LEU:HD23	1:A:245:ILE:HD11	1.98	0.44
2:B:326:ALA:HA	2:B:340:ASN:ND2	2.33	0.44
4:R:1027:ILE:HG12	4:R:1028:GLY:N	2.32	0.44
4:R:1033:LEU:CD1	4:R:1046:LEU:HB2	2.47	0.44
4:R:125:CYS:HB2	4:R:211:PRO:HB3	1.97	0.44
1:A:99:LEU:HD23	1:A:99:LEU:C	2.42	0.44
1:A:270:LEU:HD23	1:A:345:PHE:CE2	2.52	0.44
2:B:320:VAL:HG22	2:B:327:VAL:HG22	2.00	0.44
4:R:126:VAL:HG21	4:R:157:VAL:HG21	1.98	0.44
1:A:210:THR:O	1:A:221:MET:N	2.48	0.44
1:A:308:ILE:HG22	1:A:312:PHE:HB2	1.99	0.44
1:A:364:THR:CG2	1:A:365:CYS:N	2.81	0.44
2:B:230:ASN:HB2	2:B:273:ILE:O	2.18	0.44
4:R:169:ILE:HG23	4:R:174:TYR:CE2	2.51	0.44
5:N:38:ARG:HA	5:N:93:VAL:O	2.17	0.44
2:B:53:GLY:H	2:B:82:TRP:HH2	1.60	0.44
2:B:207:SER:HB3	2:B:223:THR:HG22	2.00	0.44
2:B:262:MET:HG3	2:B:263:THR:N	2.32	0.44
4:R:1004:PHE:O	4:R:1008:ARG:HB2	2.17	0.44
2:B:184:THR:HG22	2:B:184:THR:O	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:46:ARG:HD2	2:B:48:ARG:NH2	2.33	0.44
1:A:127:ARG:O	1:A:131:ILE:HG12	2.17	0.44
1:A:159:VAL:O	1:A:162:CYS:HB3	2.18	0.44
1:A:232:ARG:HB3	5:N:108:PHE:HB2	1.99	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: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
3:G:28:ILE:HD12	3:G:32:LYS:HD3	2.00	0.44
5:N:6:GLU:OE1	5:N:119:GLY:HA3	2.17	0.44
5:N:111:THR:CG2	5:N:114:THR:HG23	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
5:N:115:TYR:CE2	5:N:117:TYR:O	2.71	0.43
4:R:121:ILE:HD12	4:R:282:PHE:CE1	2.53	0.43
4:R:288:PRO:O	4:R:311:LEU:HD13	2.17	0.43
4:R:225:GLU:OE1	4:R:225:GLU:HA	2.18	0.43
5:N:83:MET:HE1	5:N:124:VAL:HG21	2.00	0.43
1:A:94:ASP:O	1:A:97:ASN:HB2	2.19	0.43
4:R:165:SER:C	4:R:168:PRO:HD2	2.43	0.43
2:B:253:PHE:CE1	2:B:260:GLU:HB3	2.53	0.43
2:B:291:ASP:O	2:B:292:PHE:CB	2.67	0.43
2:B:296:VAL:O	2:B:305:ALA:N	2.51	0.43
4:R:103:ASN:CG	4:R:188:GLU:HA	2.43	0.43
2:B:331:SER:OG	2:B:332:TRP:N	2.52	0.43
1:A:295:ASP:O	1:A:299:GLU:HG3	2.18	0.43
2:B:150:ARG:HG2	2:B:190:LEU:HD11	2.00	0.43
5:N:6:GLU:CD	5:N:119:GLY:HA3	2.44	0.43
2:B:314:ARG:HD3	2:B:332:TRP:CZ2	2.54	0.43
3:G:12:ALA:O	3:G:16:VAL:HG23	2.18	0.43
5:N:102:PRO:HB2	5:N:103:PHE:CE2	2.54	0.43
2:B:72:SER:HB3	2:B:336:LEU:HD11	2.01	0.43
4:R:38:ILE:O	4:R:42:LEU:HG	2.19	0.43
4:R:155:LEU:O	4:R:159:ILE:HG13	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
4:R:303:ILE:HG22	4:R:304:ARG:O	2.19	0.43
1:A:179:LEU:O	1:A:182:ILE:HG13	2.19	0.42
2:B:170:ASP:HB3	2:B:173:THR:OG1	2.19	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
6:R:1601:P0G:HABB	6:R:1601:P0G:HANA	1.85	0.42
4:R:1072:ASP:O	4:R:1076:ARG:HG3	2.19	0.42
4:R:90:GLY:O	4:R:94:ILE:HG13	2.18	0.42
6:R:1601:P0G:HABA	6:R:1601:P0G:HAAA	2.00	0.42
2:B:118:ASP:OD2	2:B:120:ILE:HB	2.19	0.42
4:R:30:GLU:HB2	4:R:32:TRP:H	1.84	0.42
4:R:91:ALA:HB2	4:R:313:TRP:HH2	1.84	0.42
2:B:124:TYR:HD1	2:B:135:VAL:HA	1.80	0.42
4:R:1066:LEU:HD23	4:R:1066:LEU:HA	1.93	0.42
5:N:13:GLN:CD	5:N:14:PRO:HD2	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:109:ASP:HB3	5:N:115:TYR:CZ	2.55	0.42
1:A:96:LYS:HD2	1:A:147:TYR:OH	2.19	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
2:B:149:CYS:HB2	2:B:159:THR:HG22	2.00	0.42
4:R:52:VAL:HG23	4:R:80:LEU:HD11	2.01	0.42
5:N:29:PHE:CZ	5:N:34:MET:HG3	2.54	0.42
5:N:83:MET:HB3	5:N:86:LEU:HD11	2.02	0.42
5:N:109:ASP:CB	5:N:115:TYR:CZ	3.03	0.42
1:A:231:ARG:O	1:A:235:ILE:HB	2.20	0.42
1:A:233:LYS:HG3	2:B:188:MET:SD	2.59	0.42
2:B:112:VAL:HG23	2:B:126:LEU:HD11	2.02	0.42
2:B:270:ILE:HD12	2:B:270:ILE:C	2.45	0.42
4:R:307:VAL:O	4:R:311:LEU:HG	2.19	0.42
5:N:38:ARG:HE	5:N:46:GLU:CD	2.27	0.42
1:A:28:LYS:HE2	1:A:32:LYS:HE3	2.01	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
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
1:A:238:PHE:CD1	1:A:238:PHE:N	2.89	0.41
1:A:252:SER:HB2	1:A:265:ARG:HB2	2.03	0.41
2:B:204:CYS:HA	2:B:228:ASP:HB3	2.03	0.41
2:B:317:CYS:SG	2:B:330:GLY:HA3	2.60	0.41
1:A:17:LYS:O	1:A:20:ARG:HB3	2.21	0.41
1:A:181:LYS:HG2	1:A:184:VAL:HB	2.03	0.41
1:A:229:ASP:OD1	5:N:111:THR:HG23	2.19	0.41
1:A:246:PHE:CE1	1:A:273:PHE:HB2	2.51	0.41
2:B:125:ASN:O	2:B:133:VAL:HG12	2.20	0.41
4:R:1002:ASN:CG	4:R:1003:ILE:N	2.78	0.41
4:R:197:GLN:O	4:R:201:ILE:HG13	2.21	0.41
1:A:294:GLN:CD	1:A:363:PHE:HB3	2.45	0.41
2:B:190:LEU:HD12	2:B:190:LEU:C	2.46	0.41
4:R:71:PHE:CD1	4:R:154:ILE:HD11	2.56	0.41
4:R:1036:SER:HA	4:R:1037:PRO:HD3	1.91	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:86:VAL:CG2	4:R:112:ILE:HG23	2.48	0.41
4:R:305:LYS:HD3	4:R:305:LYS:HA	1.85	0.41
5:N:4:LEU:HD21	5:N:27:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:83:MET:CB	5:N:86:LEU:HD11	2.51	0.41
2:B:271:CYS:SG	2:B:289:TYR:HB3	2.61	0.41
2:B:314:ARG:HD3	2:B:332:TRP:CD2	2.56	0.41
4:R:87:VAL:HB	4:R:88:PRO:HD3	2.02	0.41
4:R:210:VAL:O	4:R:214:ILE:HG13	2.21	0.41
2:B:326:ALA:HA	2:B:340:ASN:HD21	1.84	0.41
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
2:B:48:ARG:CG	2:B:340:ASN:HB2	2.51	0.40
2:B:161:SER:C	2:B:163:ASP:H	2.28	0.40
2:B:339:TRP:CD1	2:B:339:TRP:N	2.89	0.40
4:R:1085:LYS:N	4:R:1086:PRO:HD2	2.36	0.40
4:R:90:GLY:O	4:R:93:HIS:HB3	2.22	0.40
4:R:295:VAL:HG12	4:R:303:ILE:HD11	2.03	0.40
5:N:53:GLN:NE2	5:N:104:THR:H	2.19	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
2:B:48:ARG:HG3	2:B:340:ASN:HB2	2.03	0.40
6:R:1601:P0G:HAOA	6:R:1601:P0G:HAAB	1.89	0.40
5:N:109:ASP:O	5:N:110:VAL:HB	2.21	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 [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	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

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	R	435/514 (85%)	407 (94%)	28 (6%)	0	100	100
5	N	126/138 (91%)	117 (93%)	8 (6%)	1 (1%)	16	50
All	All	1296/1451 (89%)	1229 (95%)	66 (5%)	1 (0%)	48	79

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%)	76	83
2	B	278/293 (95%)	273 (98%)	5 (2%)	51	74
3	G	46/56 (82%)	45 (98%)	1 (2%)	45	71
4	R	358/441 (81%)	350 (98%)	8 (2%)	45	71
5	N	104/115 (90%)	102 (98%)	2 (2%)	50	73
All	All	1082/1247 (87%)	1064 (98%)	18 (2%)	53	75

All (18) 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	210	LEU
2	B	234	PHE
2	B	340	ASN
3	G	30	VAL
4	R	1016	LYS
4	R	1065	LYS

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Mol	Chain	Res	Type
4	R	1104	PHE
4	R	72	ILE
4	R	142	GLN
4	R	218	VAL
4	R	292	VAL
4	R	312	ASN
5	N	3	GLN
5	N	96	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	41	HIS
1	A	97	ASN
1	A	98	ASN
1	A	292	ASN
1	A	357	HIS
2	B	9	GLN
2	B	230	ASN
2	B	239	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	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 [i](#)

There are no oligosaccharides 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
6	P0G	R	1601	-	27,29,29	1.25	3 (11%)	34,42,42	1.41	3 (8%)

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	1601	-	-	3/15/24/24	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	R	1601	P0G	CAW-CAZ	-4.30	1.48	1.52
6	R	1601	P0G	CBA-NAP	-3.62	1.45	1.49
6	R	1601	P0G	CAX-NAQ	-2.10	1.35	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	R	1601	P0G	CAN-NAP-CBA	5.27	122.01	116.35
6	R	1601	P0G	OAR-CAY-CAW	3.57	121.75	116.76
6	R	1601	P0G	OAR-CAY-CAX	-2.98	117.44	121.16

There are no chirality outliers.

All (3) torsion outliers are listed below:

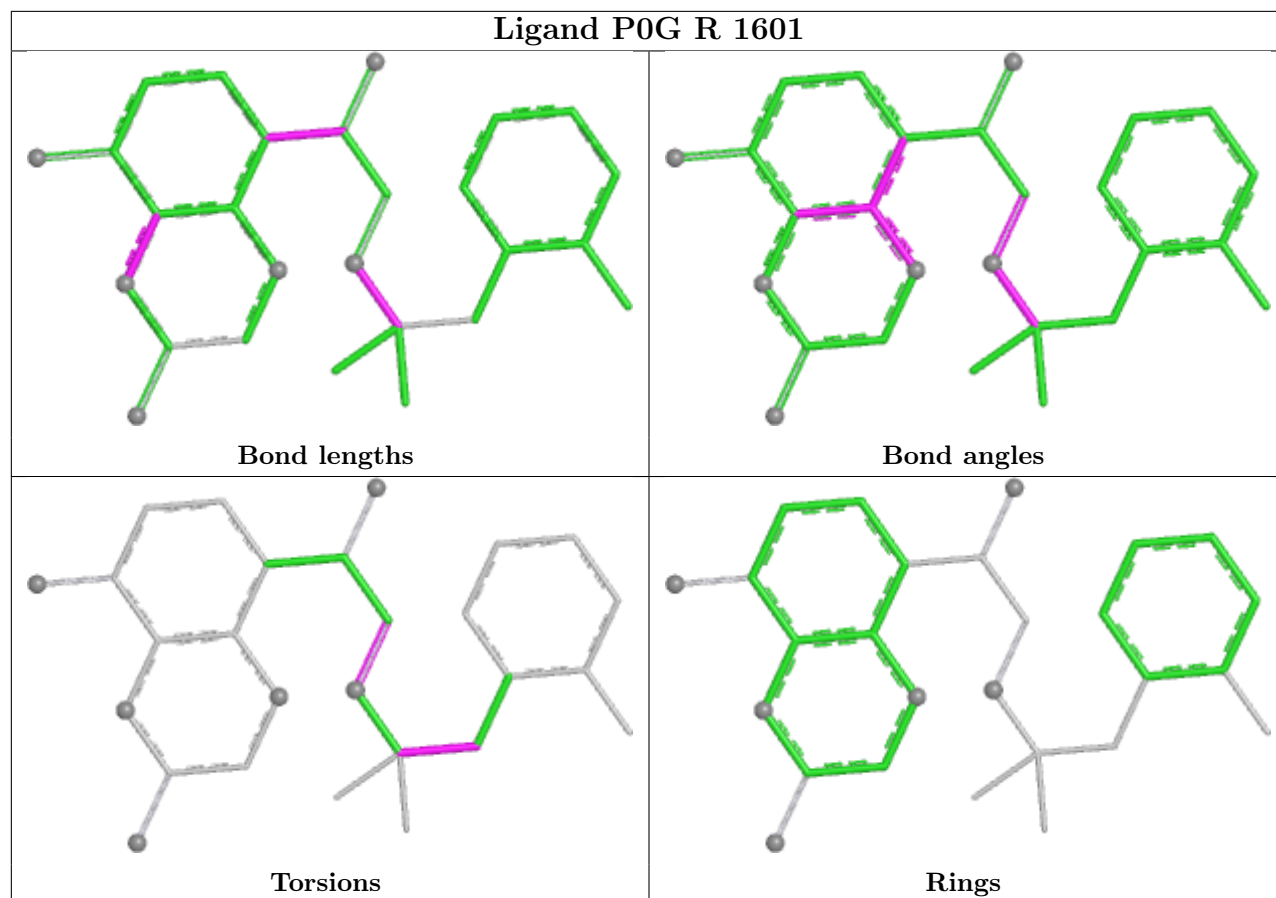
Mol	Chain	Res	Type	Atoms
6	R	1601	P0G	CAV-CAO-CBA-NAP
6	R	1601	P0G	CAZ-CAN-NAP-CBA
6	R	1601	P0G	CAV-CAO-CBA-CAB

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	R	1601	P0G	8	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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%)	0.88	38 (10%) 10 7	32, 93, 156, 191	0
2	B	340/351 (96%)	0.45	17 (5%) 34 22	29, 59, 101, 162	0
3	G	58/68 (85%)	0.69	4 (6%) 23 14	36, 80, 132, 146	0
4	R	443/514 (86%)	1.11	67 (15%) 5 4	56, 122, 175, 214	0
5	N	128/138 (92%)	0.53	6 (4%) 36 23	31, 59, 93, 115	0
All	All	1318/1451 (90%)	0.80	132 (10%) 12 8	29, 89, 158, 214	0

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	130	GLU	5.4
4	R	192	ASP	5.0
1	A	185	ILE	4.3
4	R	1094	VAL	4.1
4	R	274	THR	4.1
5	N	8	GLY	4.1
1	A	205	SER	3.7
1	A	366	ALA	3.6
1	A	56	ILE	3.6
4	R	145	LEU	3.6
1	A	206	GLY	3.5
1	A	367	VAL	3.5
4	R	309	ILE	3.5
4	R	105	TRP	3.5
4	R	1088	TYR	3.5
1	A	202	VAL	3.4
4	R	1017	ILE	3.4
4	R	1007	LEU	3.3
4	R	275	LEU	3.3
3	G	56	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
4	R	202	ALA	3.3
1	A	122	PRO	3.2
2	B	53	GLY	3.2
4	R	195	THR	3.2
1	A	136	ASN	3.1
5	N	17	SER	3.1
2	B	330	GLY	3.1
2	B	74	SER	3.1
4	R	144	LEU	3.1
1	A	199	ARG	3.1
4	R	316	TYR	3.0
1	A	59	GLN	3.0
1	A	365	CYS	3.0
1	A	51	SER	3.0
2	B	132	ASN	3.0
1	A	188	ASP	2.9
4	R	201	ILE	2.9
2	B	134	ARG	2.9
4	R	266	LEU	2.9
1	A	323	ASP	2.9
5	N	109	ASP	2.8
1	A	368	ASP	2.8
4	R	294	ILE	2.8
1	A	255	MET	2.8
5	N	128	SER	2.8
4	R	1155	THR	2.8
1	A	91	LYS	2.7
4	R	182	ILE	2.6
1	A	123	GLU	2.6
1	A	57	VAL	2.6
2	B	296	VAL	2.6
4	R	194	PHE	2.6
4	R	1097	ALA	2.6
4	R	181	ALA	2.6
4	R	1159	ASP	2.6
3	G	50	LEU	2.5
4	R	160	VAL	2.5
4	R	298	ILE	2.5
4	R	139	PHE	2.5
1	A	53	LYS	2.5
4	R	67	VAL	2.5
4	R	108	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
4	R	82	MET	2.4
5	N	110	VAL	2.4
4	R	268	GLU	2.4
4	R	100	THR	2.4
4	R	331	ASP	2.4
4	R	209	TYR	2.4
4	R	167	LEU	2.4
1	A	182	ILE	2.4
4	R	271	ALA	2.4
2	B	101	MET	2.4
4	R	297	VAL	2.4
1	A	147	TYR	2.4
4	R	278	ILE	2.3
4	R	318	ASN	2.3
1	A	9	THR	2.3
1	A	120	ALA	2.3
1	A	263	THR	2.3
4	R	92	ALA	2.3
4	R	104	PHE	2.3
3	G	42	GLU	2.3
1	A	114	VAL	2.3
2	B	133	VAL	2.3
4	R	172	HIS	2.3
1	A	189	ASP	2.3
1	A	381	ASP	2.3
1	A	52	GLY	2.3
1	A	90	THR	2.3
2	B	317	CYS	2.3
4	R	153	ILE	2.2
2	B	224	GLY	2.2
1	A	163	TYR	2.2
1	A	174	CYS	2.2
4	R	30	GLU	2.2
1	A	95	ILE	2.2
2	B	120	ILE	2.2
2	B	171	ILE	2.2
4	R	117	VAL	2.2
2	B	122	SER	2.2
1	A	48	ALA	2.2
4	R	330	PRO	2.2
1	A	378	ASP	2.2
4	R	147	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
4	R	110	THR	2.2
4	R	296	HIS	2.2
4	R	211	PRO	2.2
4	R	235	LYS	2.2
2	B	90	VAL	2.1
1	A	383	ILE	2.1
2	B	59	TYR	2.1
2	B	216	GLY	2.1
4	R	1023	GLY	2.1
4	R	267	LYS	2.1
4	R	1130	ALA	2.1
5	N	57	SER	2.1
1	A	92	VAL	2.1
3	G	54	VAL	2.1
4	R	307	VAL	2.1
4	R	109	TRP	2.1
4	R	1076	ARG	2.1
4	R	193	PHE	2.1
4	R	319	SER	2.1
4	R	206	VAL	2.1
4	R	1089	ASP	2.1
4	R	1042	ALA	2.1
4	R	91	ALA	2.0
4	R	98	THR	2.0
4	R	1003	ILE	2.0
4	R	1050	ILE	2.0
4	R	184	CYS	2.0
4	R	191	CYS	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 oligosaccharides in this entry.

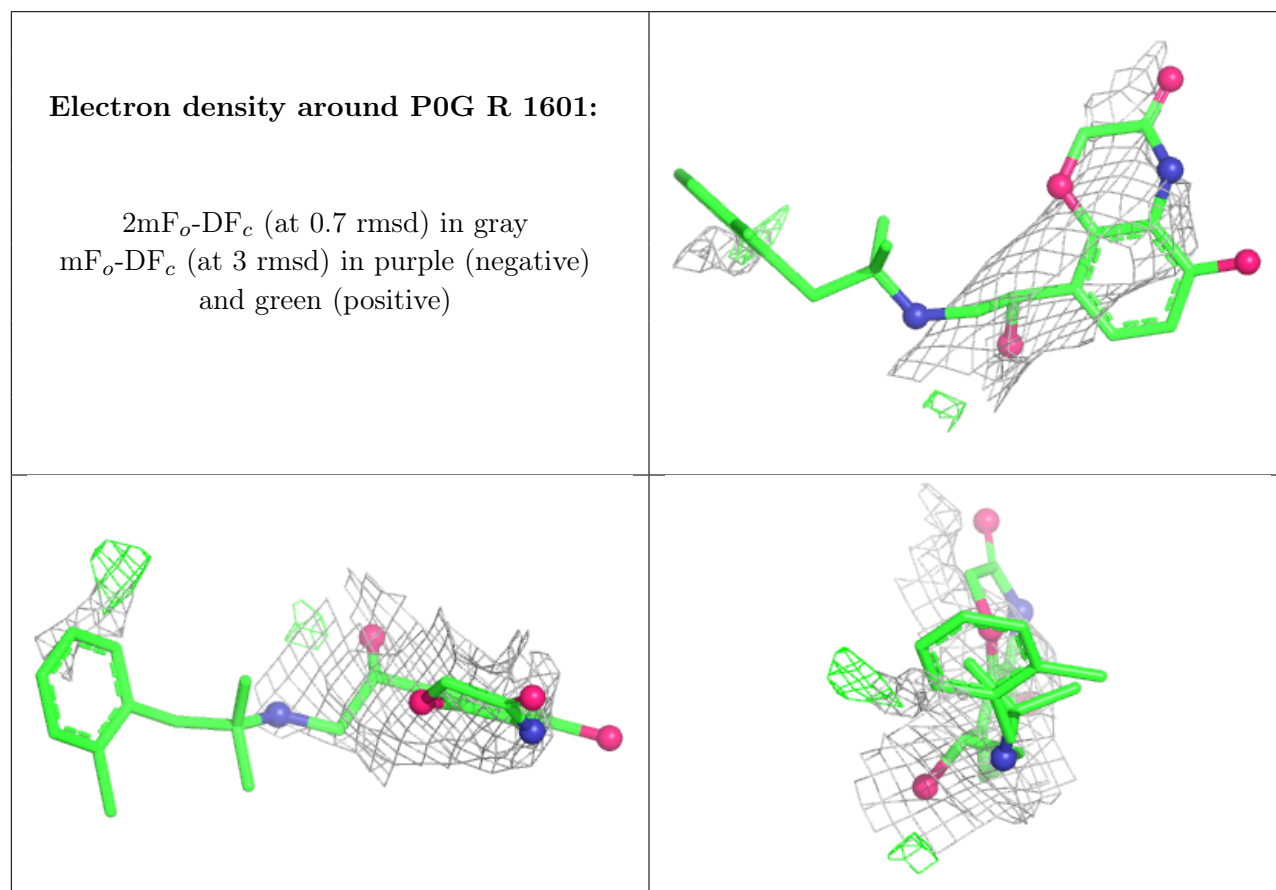
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	P0G	R	1601	27/27	0.72	0.30	152,166,192,195	0

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



6.5 Other polymers [i](#)

There are no such residues in this entry.