



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 11:13 AM EDT

PDB ID : 7VXG / pdb_00007vxg
Title : Non-receptor Protein Tyrosine Phosphatase SHP2 in Complex with Allosteric Inhibitor TK-453
Authors : Li, T.H.; Guo, H.T.; Ji, X.Y.
Deposited on : 2021-11-12
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	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

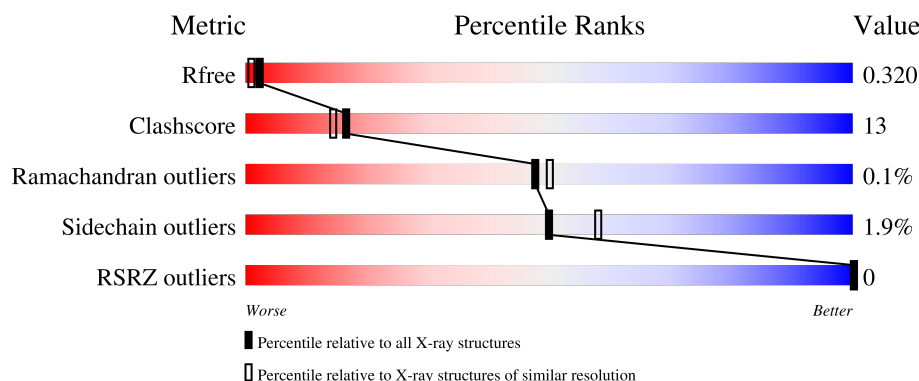
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

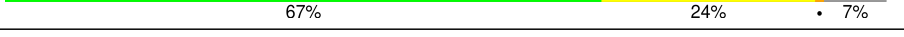
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	526	
1	B	526	
1	C	526	
1	D	526	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 29981 atoms, of which 14395 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.

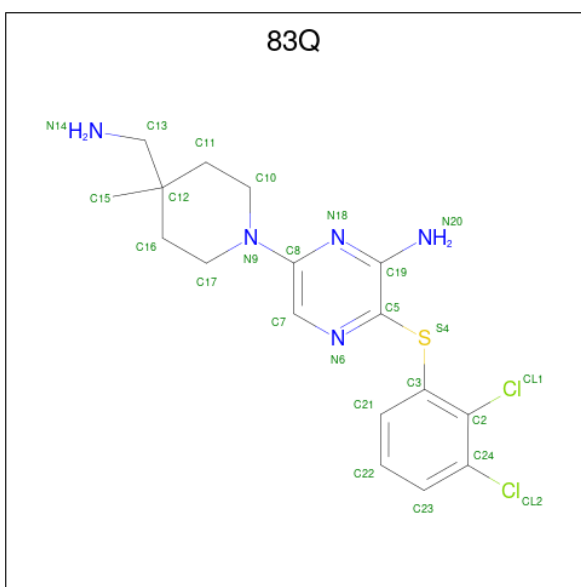
- Molecule 1 is a protein called Tyrosine-protein phosphatase non-receptor type 11.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	477	Total	C	H	N	O	S	0	0	0
			7338	2371	3581	653	715	18			
1	B	480	Total	C	H	N	O	S	0	0	0
			7409	2387	3624	666	715	17			
1	C	481	Total	C	H	N	O	S	0	0	0
			7213	2352	3478	649	718	16			
1	D	487	Total	C	H	N	O	S	0	0	0
			7457	2410	3628	674	728	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q06124
B	0	SER	-	expression tag	UNP Q06124
C	0	SER	-	expression tag	UNP Q06124
D	0	SER	-	expression tag	UNP Q06124

- Molecule 2 is 6-[4-(aminomethyl)-4-methyl-piperidin-1-yl]-3-[2,3-bis(chloranyl)phenyl]sulfanylyl-pyrazin-2-amine (CCD ID: 83Q) (formula: C₁₇H₂₁Cl₂N₅S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 46	C 17	Cl 2	H 21	N 5	S 1	0	0
2	B	1	Total 46	C 17	Cl 2	H 21	N 5	S 1	0	0
2	C	1	Total 46	C 17	Cl 2	H 21	N 5	S 1	0	0
2	D	1	Total 46	C 17	Cl 2	H 21	N 5	S 1	0	0

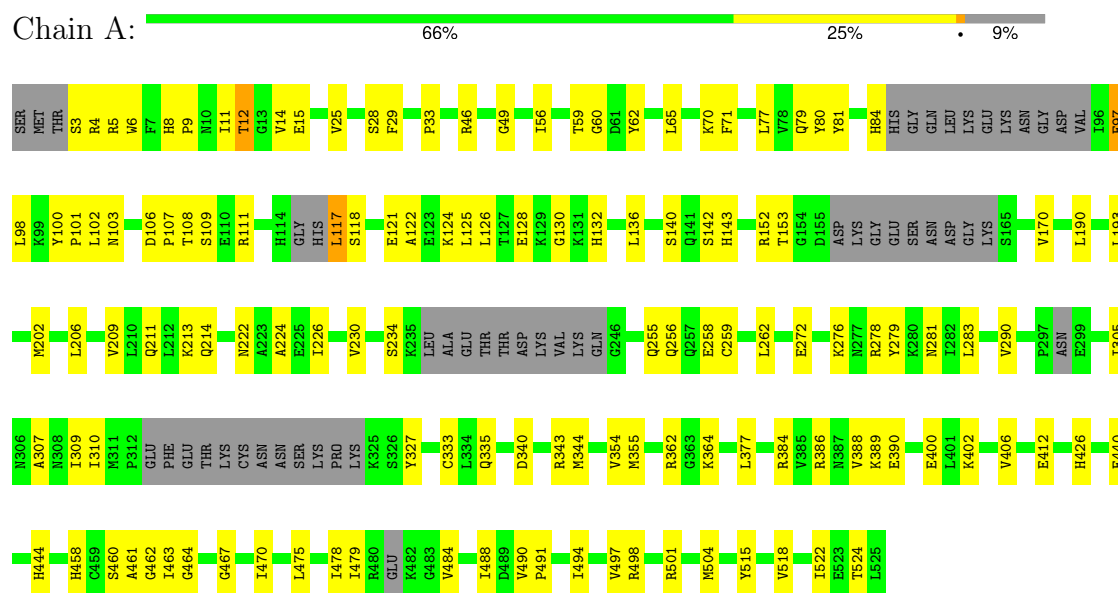
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	112	Total	O	0	0
			112	112		
3	B	87	Total	O	0	0
			87	87		
3	C	85	Total	O	0	0
			85	85		
3	D	96	Total	O	0	0
			96	96		

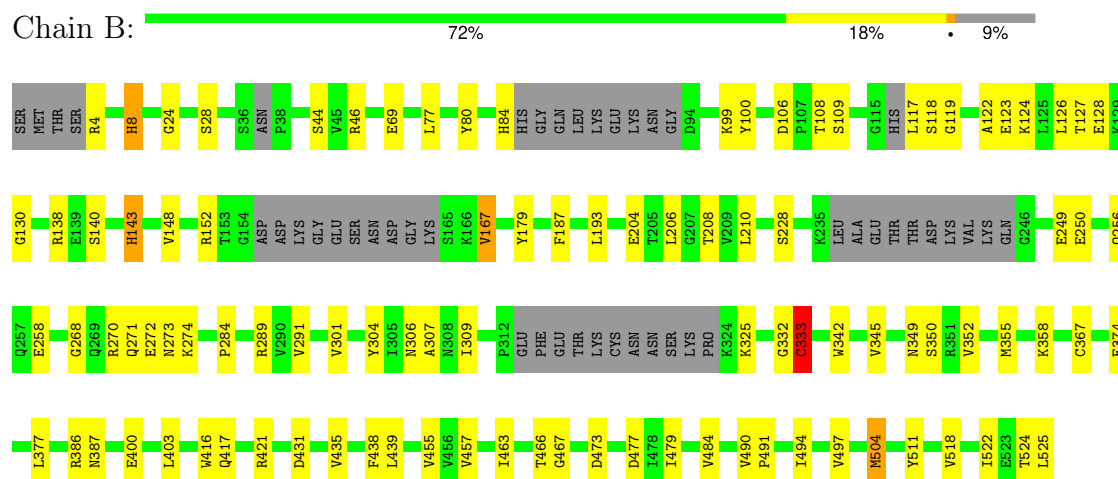
3 Residue-property plots

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: Tyrosine-protein phosphatase non-receptor type 11

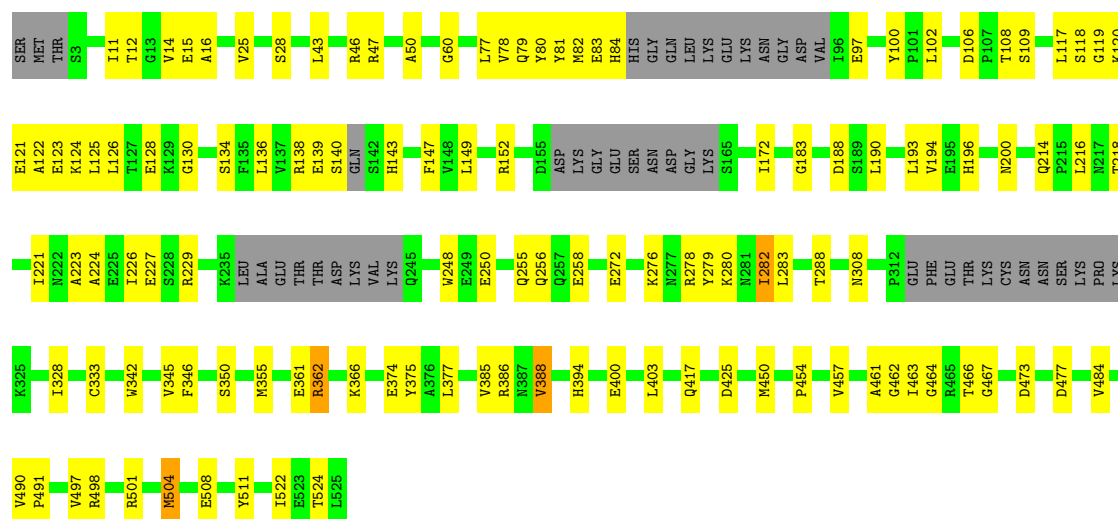


- Molecule 1: Tyrosine-protein phosphatase non-receptor type 11



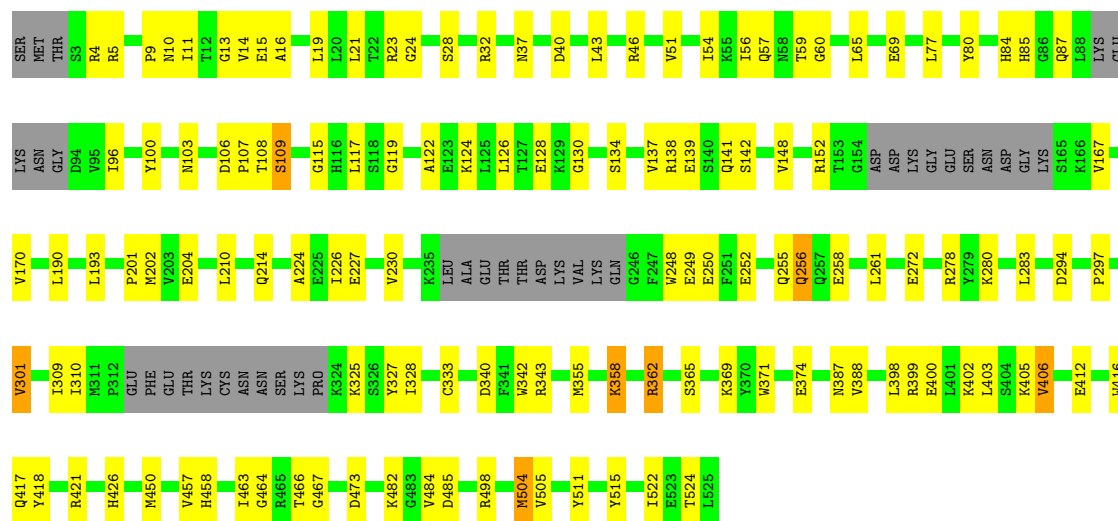
- Molecule 1: Tyrosine-protein phosphatase non-receptor type 11

Chain C:  68% 22% 9%



• Molecule 1: Tyrosine-protein phosphatase non-receptor type 11

Chain D:  67% 24% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.40Å 55.60Å 218.84Å 88.66° 102.01° 96.19°	Depositor
Resolution (Å)	38.14 – 2.10 38.14 – 2.10	Depositor EDS
% Data completeness (in resolution range)	83.7 (38.14-2.10) 83.6 (38.14-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.270 , 0.325 0.271 , 0.320	Depositor DCC
R_{free} test set	2000 reflections (1.64%)	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.430 for -h,-k,2*h+l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	29981	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.54% 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: 83Q

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	0.64	1/3835 (0.0%)	0.73	0/5195
1	B	0.65	0/3864	0.78	4/5232 (0.1%)
1	C	0.64	0/3815	0.76	2/5184 (0.0%)
1	D	0.66	1/3911 (0.0%)	0.76	2/5302 (0.0%)
All	All	0.65	2/15425 (0.0%)	0.76	8/20913 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	344	MET	SD-CE	-6.45	1.63	1.79
1	D	297	PRO	N-CD	5.49	1.55	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	504	MET	CA-C-N	-6.53	114.05	123.06
1	C	504	MET	C-N-CA	-6.53	114.05	123.06
1	B	333	CYS	N-CA-CB	6.40	121.31	110.49
1	D	504	MET	CA-C-N	-5.65	115.74	123.14
1	D	504	MET	C-N-CA	-5.65	115.74	123.14
1	B	307	ALA	N-CA-C	5.44	117.55	110.43
1	B	504	MET	CA-C-N	-5.24	116.28	123.14
1	B	504	MET	C-N-CA	-5.24	116.28	123.14

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	3757	3581	3577	101	0
1	B	3785	3624	3620	78	0
1	C	3735	3478	3486	99	0
1	D	3829	3628	3636	104	0
2	A	25	21	0	2	0
2	B	25	21	0	1	0
2	C	25	21	0	1	0
2	D	25	21	0	1	0
3	A	112	0	0	3	0
3	B	87	0	0	4	0
3	C	85	0	0	1	0
3	D	96	0	0	2	0
All	All	15586	14395	14319	381	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 13.

All (381) 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:D:4:ARG:NH1	1:D:258:GLU:OE2	1.92	1.01
1:B:4:ARG:NH1	1:B:258:GLU:OE2	1.96	0.98
1:B:249:GLU:OE2	2:B:601:83Q:N14	1.96	0.98
1:C:109:SER:OG	1:C:256:GLN:NE2	2.02	0.93
1:D:130:GLY:O	1:D:152:ARG:NH1	2.06	0.88
1:A:12:THR:OG1	1:A:15:GLU:OE1	1.89	0.88
1:D:362:ARG:HG2	1:D:362:ARG:HH11	1.35	0.87
1:C:83:GLU:OE1	1:C:83:GLU:N	2.08	0.85
1:D:84:HIS:CE1	1:D:85:HIS:HD1	1.96	0.83
1:D:362:ARG:HH11	1:D:362:ARG:CG	1.92	0.82
1:B:123:GLU:O	1:B:127:THR:OG1	1.98	0.82
1:B:352:VAL:HG21	1:B:455:VAL:HG22	1.60	0.81
1:C:342:TRP:CE3	1:C:403:LEU:HD13	2.16	0.80
1:B:352:VAL:CG2	1:B:455:VAL:HG22	2.13	0.79
1:B:518:VAL:O	1:B:522:ILE:HD12	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:250:GLU:OE2	1:D:511:TYR:OH	2.02	0.78
1:A:15:GLU:OE1	1:A:15:GLU:N	2.19	0.76
1:B:24:GLY:O	1:B:46:ARG:NH1	2.19	0.75
1:B:117:LEU:HD22	1:B:122:ALA:HB2	1.69	0.73
1:C:12:THR:OG1	1:C:15:GLU:HG3	1.90	0.72
1:D:402:LYS:HZ1	1:D:412:GLU:CD	1.97	0.72
1:B:140:SER:OG	1:B:143:HIS:O	2.07	0.71
1:B:325:LYS:NZ	1:B:473:ASP:OD2	2.19	0.71
1:B:355:MET:HE3	1:B:417:GLN:HE21	1.55	0.70
1:B:69:GLU:HA	1:B:69:GLU:OE1	1.90	0.69
1:B:479:ILE:HD13	1:B:522:ILE:HD11	1.76	0.69
1:C:82:MET:C	1:C:83:GLU:OE1	2.37	0.68
1:A:140:SER:OG	1:A:143:HIS:O	2.11	0.68
1:C:126:LEU:HD23	1:C:216:LEU:HD13	1.74	0.68
1:D:402:LYS:NZ	1:D:402:LYS:HB3	2.09	0.67
1:D:56:ILE:HG12	1:D:65:LEU:HD12	1.77	0.67
1:C:140:SER:OG	1:C:143:HIS:O	2.09	0.66
1:B:352:VAL:HG23	1:B:455:VAL:HG13	1.76	0.66
1:C:224:ALA:HB2	1:C:484:VAL:CG2	2.26	0.66
1:D:190:LEU:HD12	1:D:190:LEU:H	1.61	0.66
1:D:278:ARG:NH2	1:D:333:CYS:O	2.28	0.66
1:A:14:VAL:HG23	1:A:15:GLU:OE1	1.96	0.66
1:A:278:ARG:NH2	1:A:333:CYS:O	2.27	0.66
1:B:332:GLY:O	3:B:701:HOH:O	2.15	0.65
1:C:108:THR:O	2:C:601:83Q:N14	2.29	0.65
1:D:204:GLU:HG3	1:D:210:LEU:HD21	1.77	0.65
1:C:218:THR:O	1:C:229:ARG:NH2	2.28	0.65
1:B:250:GLU:OE2	1:B:511:TYR:OH	2.14	0.65
1:B:272:GLU:HB2	1:B:301:VAL:HG21	1.78	0.65
1:B:332:GLY:O	1:B:367:CYS:SG	2.47	0.65
1:D:524:THR:O	1:D:524:THR:HG22	1.97	0.64
1:D:294:ASP:HB2	1:D:343:ARG:HE	1.63	0.64
1:C:385:VAL:HG22	1:C:403:LEU:HD12	1.80	0.63
1:C:224:ALA:HB2	1:C:484:VAL:HG22	1.80	0.63
1:A:107:PRO:HG3	1:A:190:LEU:HD12	1.80	0.63
1:C:226:ILE:HG21	1:C:522:ILE:HD12	1.81	0.62
1:B:463:ILE:O	1:B:504:MET:O	2.18	0.62
1:D:388:VAL:HG22	1:D:400:GLU:O	1.99	0.62
1:D:342:TRP:CE3	1:D:403:LEU:HD13	2.35	0.62
1:A:117:LEU:CD1	1:A:122:ALA:HB2	2.30	0.62
1:C:362:ARG:HB3	1:C:362:ARG:HH11	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:VAL:HG21	1:A:515:TYR:CE1	2.34	0.62
1:B:289:ARG:HH11	1:B:291:VAL:HG22	1.63	0.62
1:A:462:GLY:O	1:A:501:ARG:NH1	2.32	0.62
1:A:130:GLY:O	1:A:152:ARG:NH1	2.33	0.62
1:D:224:ALA:HB2	1:D:484:VAL:HG22	1.81	0.62
1:C:272:GLU:O	1:C:276:LYS:HE3	1.99	0.61
1:D:84:HIS:CE1	1:D:85:HIS:ND1	2.66	0.61
1:A:388:VAL:HG22	1:A:400:GLU:O	2.00	0.61
1:A:389:LYS:HD3	1:A:390:GLU:N	2.15	0.61
1:B:204:GLU:HG3	1:B:210:LEU:HD21	1.81	0.61
1:C:282:ILE:C	1:C:283:LEU:HD12	2.26	0.61
1:D:340:ASP:OD1	1:D:343:ARG:NH1	2.33	0.61
1:A:463:ILE:HG23	1:A:501:ARG:HH12	1.66	0.60
1:D:388:VAL:HG23	1:D:400:GLU:HB3	1.84	0.60
1:D:24:GLY:O	1:D:46:ARG:NH1	2.35	0.60
1:B:342:TRP:CE3	1:B:403:LEU:HD13	2.37	0.59
1:A:272:GLU:O	1:A:276:LYS:HE3	2.02	0.59
1:A:467:GLY:HA3	1:A:504:MET:O	2.03	0.59
1:A:340:ASP:OD1	1:A:343:ARG:NH1	2.35	0.59
1:D:467:GLY:HA3	1:D:504:MET:O	2.02	0.59
1:A:5:ARG:HB3	1:A:103:ASN:OD1	2.03	0.58
1:D:11:ILE:HD12	1:D:16:ALA:HB2	1.85	0.58
1:C:190:LEU:HD12	1:C:190:LEU:H	1.68	0.58
1:D:4:ARG:NH1	1:D:258:GLU:CD	2.62	0.58
1:D:325:LYS:NZ	1:D:473:ASP:OD2	2.29	0.58
1:A:206:LEU:HB2	3:A:780:HOH:O	2.03	0.58
1:C:126:LEU:O	1:C:152:ARG:NH1	2.34	0.58
1:A:170:VAL:HG11	1:A:202:MET:HE1	1.85	0.57
1:D:77:LEU:HD23	1:D:77:LEU:C	2.29	0.57
1:C:280:LYS:O	1:C:283:LEU:HD11	2.03	0.57
1:D:498:ARG:NH2	1:D:505:VAL:O	2.30	0.57
1:A:440:GLU:O	1:A:444:HIS:HD2	1.88	0.57
1:B:193:LEU:C	1:B:193:LEU:HD23	2.29	0.57
1:A:117:LEU:HD11	1:A:122:ALA:HB2	1.85	0.57
1:C:130:GLY:O	1:C:152:ARG:NH1	2.38	0.57
1:C:193:LEU:HD23	1:C:193:LEU:C	2.30	0.57
1:C:497:VAL:HG12	1:C:504:MET:HG3	1.87	0.57
1:C:106:ASP:OD1	1:C:108:THR:HG23	2.05	0.56
1:C:463:ILE:HG23	1:C:501:ARG:NH1	2.20	0.56
1:D:362:ARG:HG2	1:D:362:ARG:NH1	2.11	0.56
1:D:463:ILE:O	1:D:504:MET:O	2.23	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:524:THR:O	1:D:524:THR:CG2	2.53	0.56
1:B:289:ARG:NH1	1:B:291:VAL:HG22	2.20	0.56
1:D:80:TYR:O	1:D:84:HIS:HB3	2.06	0.56
1:C:467:GLY:HA3	1:C:504:MET:O	2.05	0.56
1:A:193:LEU:C	1:A:193:LEU:HD23	2.31	0.56
1:A:377:LEU:HD13	1:A:386:ARG:HB2	1.88	0.56
1:D:402:LYS:NZ	1:D:412:GLU:OE2	2.36	0.56
1:B:358:LYS:CE	1:B:421:ARG:O	2.54	0.56
1:C:388:VAL:HG22	1:C:400:GLU:O	2.06	0.56
1:A:498:ARG:HB3	3:A:733:HOH:O	2.05	0.56
1:B:4:ARG:NH1	1:B:258:GLU:CD	2.64	0.56
1:B:301:VAL:HG23	1:B:301:VAL:O	2.06	0.55
1:B:332:GLY:O	1:B:333:CYS:HB2	2.05	0.55
1:A:126:LEU:O	1:A:152:ARG:NH1	2.38	0.55
1:C:462:GLY:O	1:C:501:ARG:NH1	2.39	0.55
1:D:230:VAL:HG22	1:D:515:TYR:HD2	1.71	0.55
1:D:362:ARG:CG	1:D:362:ARG:NH1	2.63	0.55
1:A:108:THR:O	2:A:601:83Q:N14	2.39	0.55
1:C:362:ARG:HD3	1:C:425:ASP:HB3	1.89	0.55
1:A:71:PHE:CZ	1:A:81:TYR:HE1	2.24	0.55
1:B:467:GLY:HA3	1:B:504:MET:O	2.07	0.55
1:C:328:ILE:HG13	1:C:454:PRO:HB2	1.89	0.55
1:A:524:THR:HG22	1:A:524:THR:O	2.06	0.54
1:B:457:VAL:HG12	1:B:466:THR:HG23	1.90	0.54
1:D:115:GLY:HA2	1:D:139:GLU:H	1.71	0.54
1:A:56:ILE:HD11	1:A:65:LEU:HD13	1.88	0.54
1:A:362:ARG:NH1	1:A:362:ARG:HG2	2.22	0.54
1:C:43:LEU:HD23	1:C:43:LEU:O	2.07	0.54
1:C:278:ARG:NH2	1:C:333:CYS:O	2.41	0.54
1:C:121:GLU:O	1:C:125:LEU:HG	2.07	0.54
1:A:362:ARG:HG2	1:A:362:ARG:HH11	1.73	0.54
1:C:124:LYS:O	1:C:128:GLU:N	2.36	0.54
1:A:111:ARG:HG2	2:A:601:83Q:N6	2.23	0.54
1:C:84:HIS:HD1	1:C:84:HIS:N	2.06	0.54
1:C:139:GLU:HG3	1:C:147:PHE:CE1	2.43	0.54
1:C:117:LEU:HD22	1:C:122:ALA:HB2	1.90	0.53
1:C:388:VAL:CG2	1:C:400:GLU:O	2.57	0.53
1:A:488:ILE:CD1	1:A:518:VAL:HG11	2.39	0.52
1:C:388:VAL:HG23	1:C:400:GLU:HB3	1.91	0.52
1:A:463:ILE:HG23	1:A:501:ARG:NH1	2.25	0.52
1:B:117:LEU:CD2	1:B:122:ALA:HB2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:VAL:HG22	1:A:515:TYR:HD2	1.74	0.52
1:D:193:LEU:C	1:D:193:LEU:HD23	2.35	0.52
1:B:325:LYS:HE3	1:B:477:ASP:OD2	2.10	0.52
1:B:284:PRO:HB3	1:B:306:ASN:O	2.09	0.52
1:B:271:GLN:HA	1:B:274:LYS:HE3	1.91	0.52
1:C:377:LEU:HD13	1:C:386:ARG:HB2	1.91	0.52
1:C:102:LEU:C	1:C:102:LEU:HD23	2.35	0.52
1:D:325:LYS:HZ2	1:D:473:ASP:CG	2.17	0.52
1:C:134:SER:HA	1:C:214:GLN:O	2.09	0.52
1:A:463:ILE:CG2	1:A:501:ARG:NH1	2.73	0.52
1:A:490:VAL:HB	1:A:491:PRO:CD	2.40	0.52
1:B:28:SER:HA	1:B:100:TYR:O	2.10	0.52
1:B:358:LYS:HE2	1:B:421:ARG:O	2.11	0.51
1:A:124:LYS:O	1:A:128:GLU:N	2.39	0.51
1:B:400:GLU:HG3	1:B:416:TRP:CZ3	2.46	0.51
1:D:272:GLU:HB2	1:D:301:VAL:HG21	1.93	0.51
1:A:389:LYS:HD3	1:A:390:GLU:C	2.35	0.51
1:B:179:TYR:O	1:B:187:PHE:HD1	1.93	0.51
1:C:79:GLN:O	1:C:83:GLU:OE1	2.29	0.51
1:D:84:HIS:HE1	1:D:85:HIS:CE1	2.28	0.51
1:A:109:SER:HB3	1:A:256:GLN:NE2	2.25	0.51
1:A:354:VAL:O	1:A:354:VAL:HG12	2.10	0.51
1:A:490:VAL:HB	1:A:491:PRO:HD3	1.91	0.51
1:B:126:LEU:HD13	1:B:167:VAL:HG22	1.91	0.51
1:C:288:THR:OG1	1:C:308:ASN:ND2	2.43	0.51
1:C:250:GLU:OE2	1:C:511:TYR:OH	2.28	0.51
1:D:255:GLN:OE1	1:D:498:ARG:NH1	2.38	0.51
1:B:8:HIS:N	1:B:8:HIS:CD2	2.80	0.50
1:A:255:GLN:OE1	1:A:498:ARG:NH1	2.39	0.50
1:D:398:LEU:HD13	1:D:418:TYR:CE2	2.46	0.50
1:A:121:GLU:O	1:A:125:LEU:HG	2.11	0.50
1:D:280:LYS:O	1:D:283:LEU:HD21	2.10	0.50
1:B:479:ILE:CD1	1:B:522:ILE:HD11	2.40	0.50
1:A:475:LEU:O	1:A:479:ILE:HG13	2.12	0.50
1:A:494:ILE:O	1:A:497:VAL:HG22	2.12	0.50
1:D:402:LYS:NZ	1:D:412:GLU:CD	2.69	0.50
1:A:388:VAL:HG23	1:A:400:GLU:HB3	1.94	0.49
1:D:80:TYR:O	1:D:84:HIS:CB	2.60	0.49
1:C:221:ILE:CG2	1:C:229:ARG:NH2	2.75	0.49
1:A:80:TYR:HD1	1:A:81:TYR:CD1	2.30	0.49
1:B:349:ASN:O	1:B:349:ASN:ND2	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:362:ARG:HB3	1:C:362:ARG:NH1	2.27	0.49
1:C:221:ILE:HG22	1:C:229:ARG:HH21	1.76	0.49
1:D:106:ASP:OD2	1:D:108:THR:HG23	2.12	0.49
1:D:355:MET:HE1	1:D:371:TRP:CZ3	2.47	0.49
1:A:97:GLU:OE1	1:A:97:GLU:N	2.43	0.49
1:C:77:LEU:C	1:C:77:LEU:HD23	2.38	0.49
1:D:9:PRO:HD2	1:D:10:ASN:H	1.78	0.49
1:D:5:ARG:HD2	1:D:103:ASN:HB3	1.95	0.49
1:D:109:SER:HB2	1:D:256:GLN:NE2	2.28	0.49
1:C:80:TYR:HD2	1:C:81:TYR:CD2	2.31	0.48
1:A:479:ILE:HD13	1:A:522:ILE:HD11	1.95	0.48
1:B:377:LEU:HD12	1:B:386:ARG:HB2	1.95	0.48
1:B:524:THR:O	1:B:525:LEU:HG	2.14	0.48
1:C:377:LEU:CD1	1:C:386:ARG:HB2	2.43	0.48
1:D:399:ARG:HD2	1:D:417:GLN:OE1	2.14	0.48
1:B:106:ASP:OD1	1:B:108:THR:OG1	2.22	0.48
1:B:490:VAL:HB	1:B:491:PRO:HD3	1.95	0.48
1:C:77:LEU:HD23	1:C:78:VAL:N	2.29	0.48
1:D:457:VAL:HG12	1:D:466:THR:HG23	1.94	0.48
1:C:190:LEU:H	1:C:190:LEU:CD1	2.26	0.48
1:A:226:ILE:O	1:A:230:VAL:HG23	2.14	0.48
1:B:477:ASP:HB3	3:B:771:HOH:O	2.13	0.48
1:B:518:VAL:O	1:B:522:ILE:CD1	2.59	0.48
1:D:258:GLU:HA	1:D:261:LEU:HD12	1.95	0.48
1:C:120:LYS:O	1:C:123:GLU:N	2.45	0.47
1:D:21:LEU:HD21	1:D:51:VAL:HG23	1.96	0.47
1:D:201:PRO:HD2	3:D:788:HOH:O	2.13	0.47
1:A:209:VAL:HG23	1:A:211:GLN:HG3	1.95	0.47
1:C:462:GLY:C	1:C:501:ARG:NH1	2.72	0.47
1:D:170:VAL:HG11	1:D:202:MET:HE1	1.95	0.47
1:D:28:SER:HA	1:D:100:TYR:O	2.14	0.47
1:D:124:LYS:O	1:D:128:GLU:N	2.35	0.47
1:A:402:LYS:HD2	1:A:412:GLU:OE2	2.15	0.47
1:B:119:GLY:HA2	1:B:138:ARG:CZ	2.44	0.47
1:D:107:PRO:HG3	1:D:190:LEU:HD13	1.97	0.47
1:B:270:ARG:NE	3:B:712:HOH:O	2.47	0.47
1:B:457:VAL:CG1	1:B:466:THR:HG23	2.44	0.47
1:D:325:LYS:HE2	3:D:750:HOH:O	2.15	0.47
1:A:460:SER:OG	1:A:461:ALA:N	2.47	0.47
1:A:4:ARG:CZ	1:A:258:GLU:OE2	2.62	0.47
1:A:384:ARG:NH2	1:A:406:VAL:HG13	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:GLU:CG	1:B:210:LEU:HD21	2.45	0.47
1:A:132:HIS:ND1	1:A:153:THR:HA	2.30	0.47
1:B:490:VAL:O	1:B:494:ILE:HG13	2.15	0.47
1:C:223:ALA:O	1:C:522:ILE:HD13	2.15	0.47
1:A:46:ARG:HE	1:A:49:GLY:HA2	1.79	0.46
1:D:10:ASN:ND2	1:D:248:TRP:HH2	2.13	0.46
1:D:54:ILE:HG21	1:D:65:LEU:HD11	1.97	0.46
1:A:102:LEU:C	1:A:102:LEU:HD23	2.39	0.46
1:A:402:LYS:HB2	1:A:402:LYS:NZ	2.31	0.46
1:A:136:LEU:HD12	1:A:136:LEU:C	2.41	0.46
1:A:484:VAL:HG22	1:A:484:VAL:O	2.15	0.46
1:C:457:VAL:HG12	1:C:466:THR:HG23	1.96	0.46
1:D:249:GLU:OE2	2:D:601:83Q:N14	2.49	0.46
1:D:402:LYS:NZ	1:D:412:GLU:CG	2.79	0.46
1:B:273:ASN:O	1:B:274:LYS:C	2.58	0.46
1:B:435:VAL:O	1:B:438:PHE:HB3	2.15	0.46
1:A:290:VAL:HG12	1:A:307:ALA:HB3	1.97	0.46
1:B:117:LEU:HD23	1:B:118:SER:O	2.14	0.46
1:C:524:THR:HG22	1:C:524:THR:O	2.16	0.46
1:D:138:ARG:HG3	1:D:148:VAL:HB	1.98	0.46
1:D:402:LYS:HB3	1:D:402:LYS:HZ2	1.81	0.45
1:A:402:LYS:HB2	1:A:402:LYS:HZ3	1.81	0.45
1:A:470:ILE:HD12	1:A:504:MET:HE1	1.98	0.45
1:B:268:GLY:HA2	1:B:304:TYR:CE2	2.51	0.45
1:C:190:LEU:HD12	1:C:190:LEU:N	2.30	0.45
1:D:405:LYS:O	1:D:406:VAL:C	2.58	0.45
1:A:281:ASN:C	1:A:283:LEU:HD12	2.41	0.45
1:C:355:MET:HE3	1:C:417:GLN:NE2	2.31	0.45
1:C:473:ASP:O	1:C:477:ASP:HB2	2.17	0.45
1:D:87:GLN:CB	1:D:96:ILE:H	2.30	0.45
1:A:402:LYS:HD3	1:A:412:GLU:CD	2.41	0.45
1:C:47:ARG:O	1:C:50:ALA:N	2.46	0.45
1:A:29:PHE:CD2	1:A:98:LEU:HD23	2.51	0.45
1:A:222:ASN:OD1	1:A:224:ALA:N	2.48	0.45
1:B:109:SER:OG	1:B:256:GLN:NE2	2.49	0.45
1:C:248:TRP:CE3	1:C:508:GLU:OE1	2.70	0.45
1:A:279:TYR:CE2	1:A:461:ALA:HB2	2.52	0.45
1:B:80:TYR:O	1:B:84:HIS:HB3	2.17	0.45
1:C:80:TYR:CD1	1:C:84:HIS:CE1	3.05	0.45
1:A:4:ARG:HB2	1:A:259:CYS:HB3	1.98	0.45
1:C:226:ILE:HG23	1:C:227:GLU:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:LYS:O	1:B:128:GLU:N	2.38	0.45
1:B:355:MET:HE3	1:B:417:GLN:NE2	2.27	0.45
1:C:490:VAL:HB	1:C:491:PRO:CD	2.47	0.45
1:C:490:VAL:HB	1:C:491:PRO:HD3	1.98	0.45
1:D:40:ASP:OD2	1:D:57:GLN:HG2	2.16	0.44
1:A:80:TYR:O	1:A:84:HIS:CD2	2.71	0.44
1:A:8:HIS:HB2	1:A:11:ILE:CD1	2.47	0.44
1:C:149:LEU:HB2	1:C:172:ILE:HD11	1.98	0.44
1:A:117:LEU:HD13	1:A:118:SER:O	2.18	0.44
1:A:389:LYS:HD3	1:A:390:GLU:O	2.18	0.44
1:C:60:GLY:O	1:C:464:GLY:HA3	2.17	0.44
1:C:117:LEU:CD2	1:C:118:SER:O	2.66	0.44
1:D:301:VAL:CG2	1:D:301:VAL:O	2.65	0.44
1:D:310:ILE:HB	1:D:327:TYR:HB2	2.00	0.44
1:A:283:LEU:HD12	1:A:283:LEU:N	2.32	0.44
1:D:134:SER:HA	1:D:214:GLN:O	2.18	0.44
1:A:213:LYS:C	1:A:214:GLN:HG3	2.42	0.44
1:D:450:MET:HE2	1:D:450:MET:HA	2.00	0.44
1:B:77:LEU:HD23	1:B:77:LEU:C	2.43	0.43
1:D:13:GLY:HA2	1:D:32:ARG:CZ	2.48	0.43
1:D:362:ARG:HH21	1:D:426:HIS:CE1	2.36	0.43
1:A:490:VAL:O	1:A:491:PRO:C	2.60	0.43
1:C:84:HIS:N	1:C:84:HIS:ND1	2.65	0.43
1:C:109:SER:HG	1:C:256:GLN:NE2	2.12	0.43
1:D:69:GLU:OE2	1:D:69:GLU:HA	2.17	0.43
1:D:369:LYS:HE2	1:D:371:TRP:NE1	2.32	0.43
1:A:117:LEU:HD13	1:A:117:LEU:C	2.43	0.43
1:A:28:SER:HA	1:A:100:TYR:O	2.19	0.43
1:C:342:TRP:CE3	1:C:403:LEU:CD1	2.95	0.43
1:C:28:SER:HA	1:C:100:TYR:O	2.18	0.43
1:C:255:GLN:O	1:C:258:GLU:HG2	2.18	0.43
1:C:136:LEU:HD12	1:C:136:LEU:C	2.43	0.43
1:C:450:MET:HE2	1:C:450:MET:HA	2.01	0.43
1:D:457:VAL:CG1	1:D:466:THR:HG23	2.49	0.43
1:A:59:THR:HG21	1:A:426:HIS:CE1	2.54	0.43
1:C:183:GLY:N	3:C:710:HOH:O	2.51	0.43
1:A:6:TRP:HB3	1:A:101:PRO:HB3	2.01	0.43
1:C:361:GLU:HB2	1:C:366:LYS:HE3	2.00	0.43
1:C:374:GLU:HG2	1:C:375:TYR:CD2	2.54	0.43
1:A:9:PRO:O	1:A:33:PRO:HG3	2.19	0.43
1:C:80:TYR:O	1:C:84:HIS:ND1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:346:PHE:HB2	1:C:403:LEU:HD21	2.01	0.43
1:A:8:HIS:HB2	1:A:11:ILE:HD11	1.99	0.42
1:A:79:GLN:CD	1:A:262:LEU:HD22	2.44	0.42
1:A:310:ILE:HB	1:A:327:TYR:HB2	2.01	0.42
1:D:19:LEU:HD21	1:D:23:ARG:NH2	2.33	0.42
1:D:60:GLY:O	1:D:464:GLY:HA3	2.19	0.42
1:C:97:GLU:HA	1:C:97:GLU:OE1	2.19	0.42
1:C:463:ILE:CG2	1:C:501:ARG:NH1	2.83	0.42
1:C:501:ARG:HB3	1:C:504:MET:HG2	2.00	0.42
1:D:374:GLU:HA	1:D:387:ASN:CG	2.44	0.42
1:B:439:LEU:HD23	1:B:439:LEU:HA	1.92	0.42
1:B:138:ARG:O	1:B:148:VAL:N	2.45	0.42
1:B:497:VAL:HG12	1:B:504:MET:HG3	2.01	0.42
1:B:130:GLY:O	1:B:152:ARG:NH1	2.53	0.42
1:C:196:HIS:O	1:C:200:ASN:HB2	2.20	0.42
1:C:279:TYR:CE2	1:C:461:ALA:HB2	2.54	0.42
1:C:345:VAL:HG13	1:C:350:SER:HB2	2.01	0.42
1:D:43:LEU:HB3	1:D:54:ILE:HB	2.02	0.42
1:D:119:GLY:HA2	1:D:138:ARG:CZ	2.49	0.42
1:A:355:MET:HB2	1:A:458:HIS:NE2	2.35	0.42
1:A:46:ARG:HH21	1:A:49:GLY:HA2	1.84	0.42
1:A:71:PHE:CZ	1:A:81:TYR:CE1	3.07	0.42
1:C:117:LEU:HD23	1:C:118:SER:N	2.34	0.42
1:D:141:GLN:OE1	1:D:141:GLN:N	2.52	0.42
1:A:290:VAL:HG13	1:A:305:ILE:CG2	2.50	0.42
1:A:402:LYS:CD	1:A:412:GLU:OE2	2.67	0.42
1:C:394:HIS:CD2	1:D:365:SER:HB2	2.55	0.42
1:D:40:ASP:CG	1:D:57:GLN:HG2	2.44	0.42
1:A:71:PHE:CD1	1:A:77:LEU:HD12	2.55	0.42
1:A:475:LEU:HA	1:A:478:ILE:HD12	2.02	0.42
1:C:119:GLY:HA2	1:C:138:ARG:CZ	2.50	0.42
1:C:221:ILE:HG22	1:C:229:ARG:NH2	2.34	0.42
1:A:106:ASP:CG	1:A:108:THR:HG1	2.28	0.41
1:D:355:MET:HE2	1:D:458:HIS:NE2	2.35	0.41
1:D:482:LYS:HE3	1:D:485:ASP:O	2.20	0.41
1:A:25:VAL:C	1:A:46:ARG:HD3	2.45	0.41
1:A:498:ARG:NH2	3:A:720:HOH:O	2.53	0.41
1:B:138:ARG:HG3	1:B:148:VAL:HB	2.02	0.41
1:B:271:GLN:HA	1:B:274:LYS:CE	2.50	0.41
1:D:126:LEU:O	1:D:152:ARG:NH1	2.50	0.41
1:D:252:GLU:OE1	1:D:252:GLU:HA	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:GLY:O	1:A:464:GLY:HA3	2.20	0.41
1:B:387:ASN:OD1	1:B:387:ASN:C	2.63	0.41
1:C:25:VAL:O	1:C:46:ARG:HD3	2.19	0.41
1:D:226:ILE:HG21	1:D:522:ILE:HD13	2.02	0.41
1:D:226:ILE:HG23	1:D:227:GLU:N	2.36	0.41
1:C:43:LEU:HD23	1:C:43:LEU:C	2.46	0.41
1:D:309:ILE:HD13	1:D:328:ILE:HG12	2.01	0.41
1:D:137:VAL:HA	1:D:148:VAL:O	2.21	0.41
1:C:255:GLN:OE1	1:C:498:ARG:NH1	2.49	0.41
1:C:11:ILE:HD12	1:C:16:ALA:HB2	2.01	0.41
1:C:14:VAL:O	1:C:15:GLU:C	2.62	0.41
1:D:84:HIS:CE1	1:D:85:HIS:CE1	3.07	0.41
1:B:117:LEU:CD2	1:B:118:SER:O	2.69	0.41
1:B:270:ARG:HG2	3:B:712:HOH:O	2.21	0.41
1:C:463:ILE:O	1:C:504:MET:O	2.39	0.41
1:D:37:ASN:HB2	1:D:40:ASP:HB2	2.03	0.41
1:D:358:LYS:CE	1:D:421:ARG:O	2.69	0.41
1:D:400:GLU:HG3	1:D:416:TRP:CZ3	2.56	0.41
1:B:374:GLU:HA	1:B:387:ASN:CG	2.46	0.41
1:D:59:THR:HG21	1:D:426:HIS:NE2	2.36	0.41
1:D:355:MET:HG3	1:D:458:HIS:CE1	2.56	0.41
1:B:345:VAL:HG13	1:B:350:SER:HB2	2.02	0.40
1:B:352:VAL:CG2	1:B:455:VAL:HG13	2.45	0.40
1:C:190:LEU:O	1:C:194:VAL:HG23	2.21	0.40
1:D:14:VAL:O	1:D:15:GLU:C	2.64	0.40
1:A:3:SER:C	1:A:5:ARG:H	2.28	0.40
1:D:398:LEU:HD13	1:D:418:TYR:CZ	2.57	0.40
1:A:109:SER:HB3	1:A:256:GLN:HE22	1.86	0.40
1:A:364:LYS:HE2	1:A:364:LYS:HA	2.03	0.40
1:B:106:ASP:OD1	1:B:108:THR:HG23	2.21	0.40
1:C:221:ILE:CG2	1:C:229:ARG:CZ	2.99	0.40
1:D:56:ILE:HG12	1:D:65:LEU:CD1	2.48	0.40
1:B:117:LEU:HD23	1:B:118:SER:N	2.35	0.40
1:B:272:GLU:HB2	1:B:301:VAL:CG2	2.48	0.40
1:D:117:LEU:HD23	1:D:122:ALA:HB2	2.03	0.40
1:D:402:LYS:HB3	1:D:402:LYS:HZ3	1.85	0.40
1:A:62:TYR:CE1	1:A:70:LYS:HD3	2.57	0.40
1:B:99:LYS:HE3	1:B:100:TYR:CE2	2.57	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	461/526 (88%)	445 (96%)	16 (4%)	0	100	100
1	B	466/526 (89%)	452 (97%)	13 (3%)	1 (0%)	43	44
1	C	469/526 (89%)	448 (96%)	21 (4%)	0	100	100
1	D	477/526 (91%)	458 (96%)	19 (4%)	0	100	100
All	All	1873/2104 (89%)	1803 (96%)	69 (4%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	333	CYS

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	401/468 (86%)	394 (98%)	7 (2%)	53	62
1	B	403/468 (86%)	392 (97%)	11 (3%)	39	45
1	C	391/468 (84%)	387 (99%)	4 (1%)	68	76
1	D	406/468 (87%)	398 (98%)	8 (2%)	48	56
All	All	1601/1872 (86%)	1571 (98%)	30 (2%)	50	58

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	THR
1	A	97	GLU
1	A	117	LEU
1	A	142	SER
1	A	234	SER
1	A	309	ILE
1	A	335	GLN
1	B	8	HIS
1	B	44	SER
1	B	143	HIS
1	B	167	VAL
1	B	206	LEU
1	B	208	THR
1	B	228	SER
1	B	309	ILE
1	B	333	CYS
1	B	431	ASP
1	B	484	VAL
1	C	188	ASP
1	C	282	ILE
1	C	362	ARG
1	C	388	VAL
1	D	109	SER
1	D	142	SER
1	D	167	VAL
1	D	256	GLN
1	D	301	VAL
1	D	358	LYS
1	D	362	ARG
1	D	406	VAL

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	169	HIS
1	A	256	GLN
1	A	287	HIS
1	A	426	HIS
1	B	10	ASN
1	B	256	GLN
1	B	408	GLN
1	B	417	GLN
1	B	520	HIS

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Mol	Chain	Res	Type
1	C	196	HIS
1	C	256	GLN
1	C	308	ASN
1	C	335	GLN
1	C	426	HIS
1	D	10	ASN
1	D	257	GLN
1	D	269	GLN
1	D	281	ASN
1	D	293	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	83Q	D	601	-	25,27,27	2.44	10 (40%)	31,39,39	1.65	8 (25%)
2	83Q	B	601	-	25,27,27	2.81	13 (52%)	31,39,39	1.92	8 (25%)
2	83Q	A	601	-	25,27,27	1.84	4 (16%)	31,39,39	1.90	8 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	83Q	C	601	-	25,27,27	2.46	8 (32%)	31,39,39	1.73	9 (29%)

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
2	83Q	D	601	-	-	1/11/23/23	0/3/3/3
2	83Q	B	601	-	-	4/11/23/23	0/3/3/3
2	83Q	A	601	-	-	1/11/23/23	0/3/3/3
2	83Q	C	601	-	-	1/11/23/23	0/3/3/3

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	83Q	C2-CL1	-6.83	1.58	1.72
2	B	601	83Q	C10-N9	6.75	1.58	1.46
2	B	601	83Q	C2-CL1	-6.57	1.59	1.72
2	C	601	83Q	C24-CL2	-6.23	1.58	1.73
2	C	601	83Q	C10-N9	5.60	1.56	1.46
2	A	601	83Q	C10-N9	5.56	1.56	1.46
2	C	601	83Q	C24-C2	-4.52	1.32	1.39
2	D	601	83Q	C11-C12	-4.30	1.44	1.53
2	B	601	83Q	C7-C8	3.95	1.48	1.40
2	B	601	83Q	C11-C12	-3.71	1.46	1.53
2	B	601	83Q	C17-N9	-3.40	1.40	1.46
2	D	601	83Q	C10-N9	3.32	1.52	1.46
2	B	601	83Q	C24-CL2	-3.27	1.65	1.73
2	B	601	83Q	C8-N18	3.25	1.39	1.34
2	D	601	83Q	C24-CL2	-3.19	1.66	1.73
2	C	601	83Q	C8-N18	3.07	1.38	1.34
2	D	601	83Q	C17-N9	3.03	1.51	1.46
2	C	601	83Q	C7-C8	2.91	1.46	1.40
2	B	601	83Q	C24-C2	-2.83	1.35	1.39
2	D	601	83Q	C16-C12	-2.80	1.48	1.53
2	D	601	83Q	C19-N18	2.69	1.38	1.35
2	D	601	83Q	C5-N6	2.67	1.37	1.34
2	B	601	83Q	C23-C24	-2.60	1.33	1.38
2	A	601	83Q	C24-C2	-2.59	1.35	1.39
2	C	601	83Q	C2-C3	-2.56	1.33	1.39
2	C	601	83Q	C5-N6	2.50	1.37	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	83Q	C11-C12	-2.41	1.48	1.53
2	B	601	83Q	C5-N6	2.41	1.37	1.34
2	A	601	83Q	C24-CL2	-2.40	1.68	1.73
2	B	601	83Q	C19-N18	2.36	1.38	1.35
2	B	601	83Q	C16-C12	-2.32	1.49	1.53
2	D	601	83Q	C8-N18	2.17	1.37	1.34
2	A	601	83Q	C11-C12	-2.11	1.49	1.53
2	B	601	83Q	C2-C3	-2.10	1.34	1.39
2	D	601	83Q	C24-C2	-2.10	1.36	1.39

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	83Q	C24-C2-CL1	-5.24	113.38	120.02
2	B	601	83Q	C24-C2-CL1	-4.34	114.52	120.02
2	D	601	83Q	C24-C2-CL1	-4.20	114.69	120.02
2	B	601	83Q	C3-S4-C5	4.19	114.88	102.93
2	A	601	83Q	C3-S4-C5	3.91	114.08	102.93
2	C	601	83Q	C7-C8-N18	-3.69	116.56	121.22
2	B	601	83Q	N18-C8-N9	3.61	121.11	116.56
2	D	601	83Q	C3-S4-C5	3.55	113.05	102.93
2	C	601	83Q	C24-C2-CL1	-3.42	115.69	120.02
2	C	601	83Q	C3-S4-C5	3.35	112.49	102.93
2	A	601	83Q	C7-C8-N18	-3.30	117.05	121.22
2	C	601	83Q	C15-C12-C13	-3.28	105.28	108.22
2	A	601	83Q	N18-C8-N9	3.16	120.55	116.56
2	B	601	83Q	C11-C10-N9	3.03	116.02	110.53
2	A	601	83Q	C16-C17-N9	2.90	115.79	110.53
2	C	601	83Q	C11-C10-N9	2.85	115.69	110.53
2	B	601	83Q	C16-C17-N9	2.82	115.64	110.53
2	D	601	83Q	C7-C8-N18	-2.81	117.67	121.22
2	B	601	83Q	C2-C3-S4	2.74	126.23	120.27
2	B	601	83Q	C7-C8-N18	-2.69	117.82	121.22
2	A	601	83Q	C11-C10-N9	2.46	115.00	110.53
2	C	601	83Q	N18-C8-N9	2.46	119.66	116.56
2	A	601	83Q	C8-C7-N6	2.32	124.31	121.31
2	D	601	83Q	C22-C23-C24	-2.31	115.95	119.42
2	D	601	83Q	C11-C10-N9	2.26	114.63	110.53
2	D	601	83Q	C21-C3-C2	-2.26	114.71	117.67
2	D	601	83Q	C2-C3-S4	2.23	125.13	120.27
2	C	601	83Q	C23-C24-C2	2.15	123.11	120.52
2	C	601	83Q	C22-C23-C24	-2.11	116.25	119.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	83Q	C8-C7-N6	2.10	124.03	121.31
2	A	601	83Q	C23-C24-C2	2.08	123.02	120.52
2	B	601	83Q	C15-C12-C11	2.03	112.84	109.75
2	D	601	83Q	C8-C7-N6	2.02	123.92	121.31

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	83Q	N6-C5-S4-C3
2	B	601	83Q	C15-C12-C13-N14
2	B	601	83Q	C11-C12-C13-N14
2	B	601	83Q	C16-C12-C13-N14
2	B	601	83Q	N6-C5-S4-C3
2	C	601	83Q	N6-C5-S4-C3
2	D	601	83Q	N6-C5-S4-C3

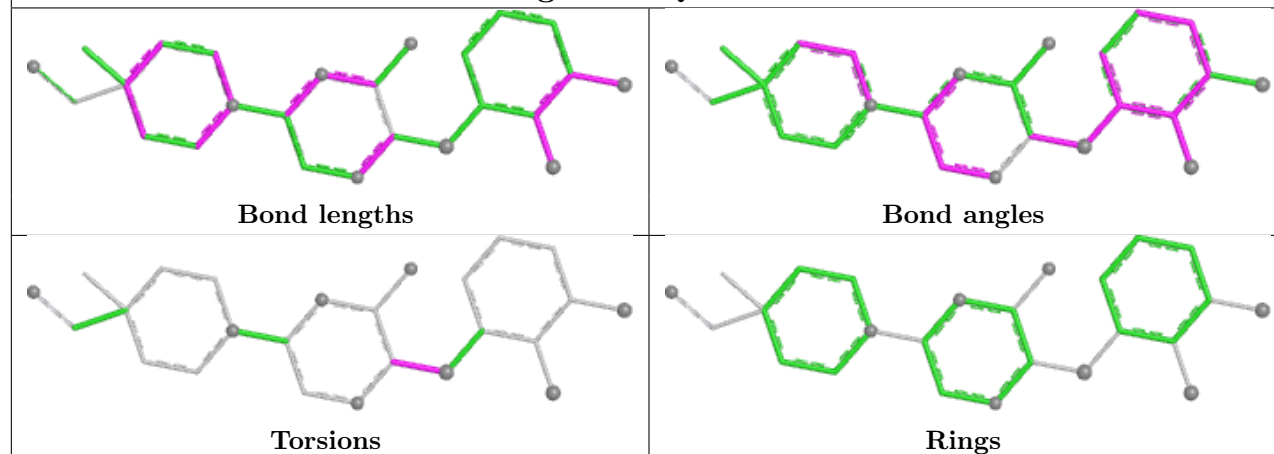
There are no ring outliers.

4 monomers are involved in 5 short contacts:

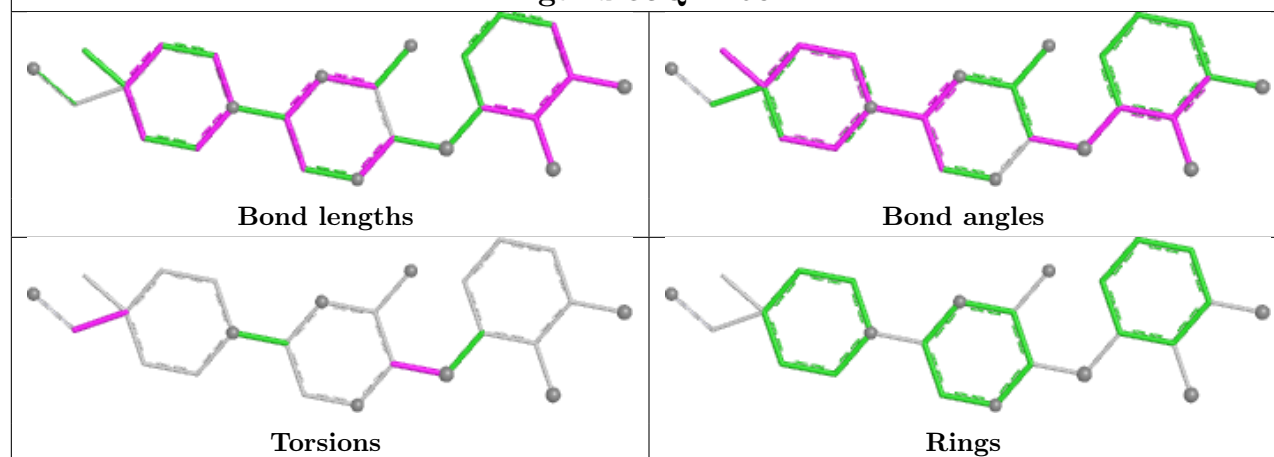
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	601	83Q	1	0
2	B	601	83Q	1	0
2	A	601	83Q	2	0
2	C	601	83Q	1	0

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.

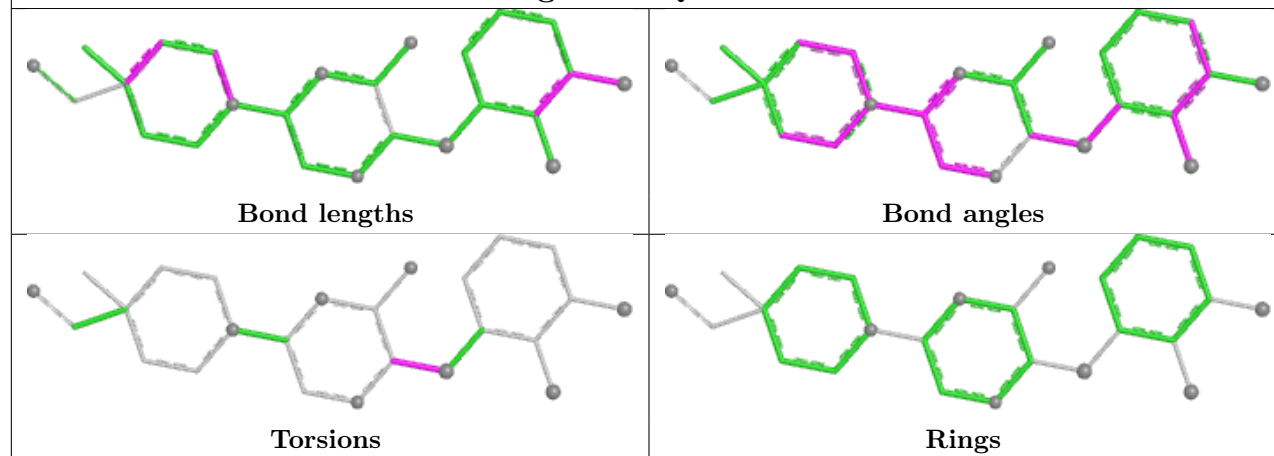
Ligand 83Q D 601

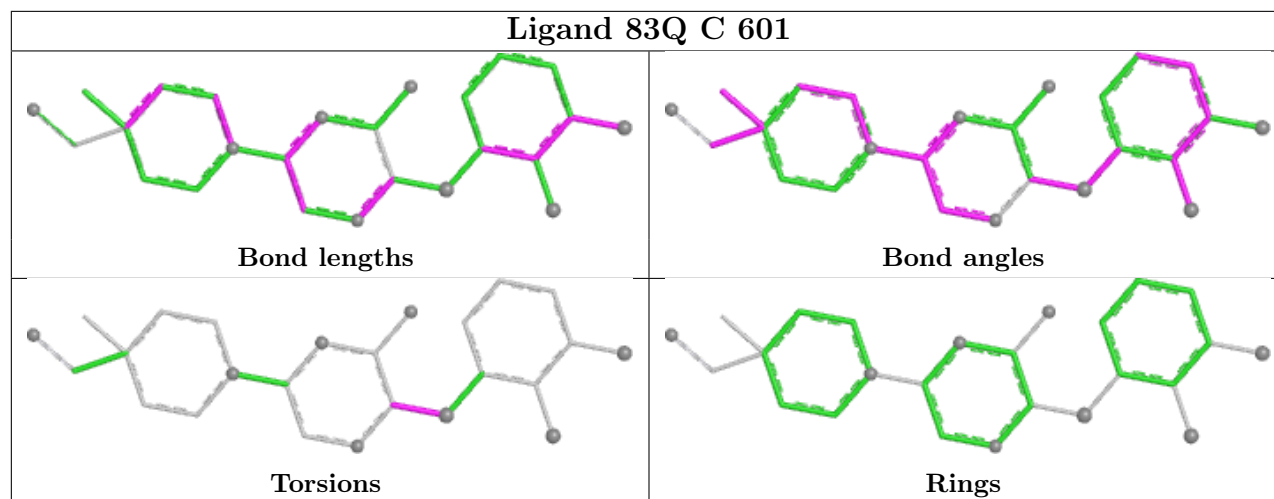


Ligand 83Q B 601



Ligand 83Q A 601





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	477/526 (90%)	-0.87	0 100 100	20, 40, 64, 90	0
1	B	480/526 (91%)	-0.87	0 100 100	22, 38, 63, 107	0
1	C	481/526 (91%)	-0.88	0 100 100	20, 38, 63, 111	0
1	D	487/526 (92%)	-0.86	0 100 100	19, 38, 66, 124	0
All	All	1925/2104 (91%)	-0.87	0 100 100	19, 38, 65, 124	0

There are no RSRZ outliers to report.

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](#)

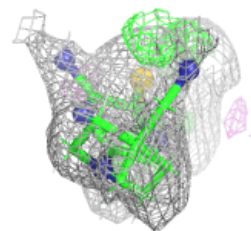
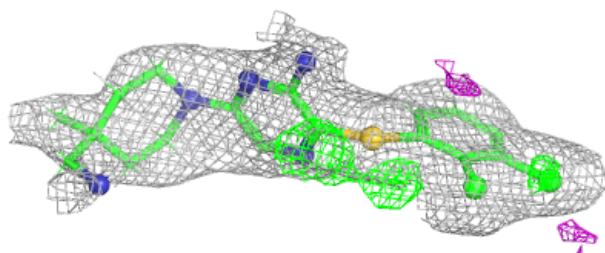
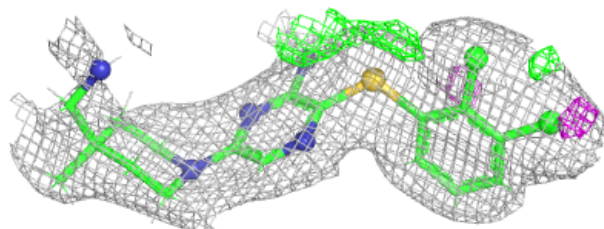
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	83Q	A	601	25/25	0.99	0.04	14,31,42,49	0
2	83Q	B	601	25/25	0.99	0.04	11,31,49,53	0
2	83Q	C	601	25/25	0.99	0.05	16,30,38,49	0
2	83Q	D	601	25/25	0.99	0.05	19,32,45,47	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.

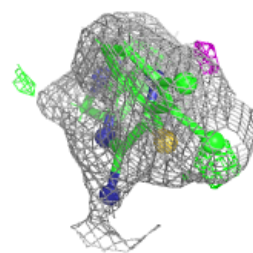
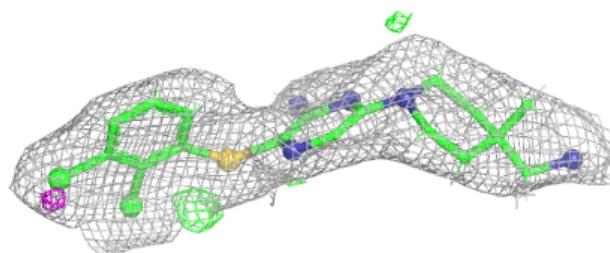
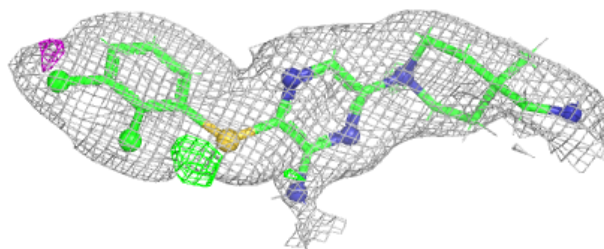
Electron density around 83Q A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

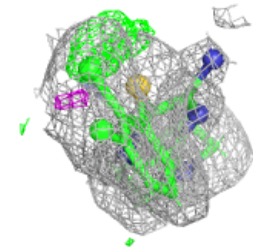
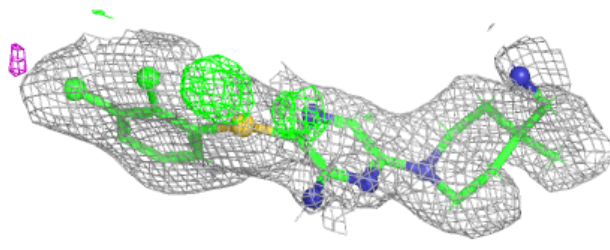
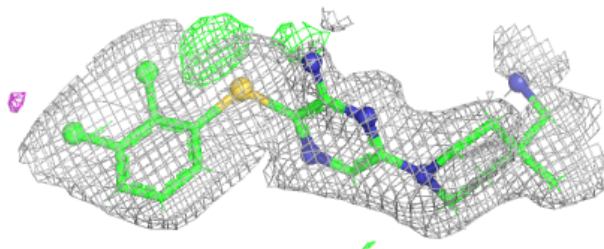


Electron density around 83Q B 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

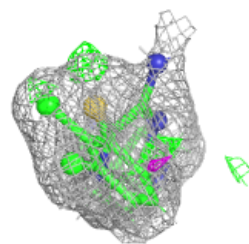
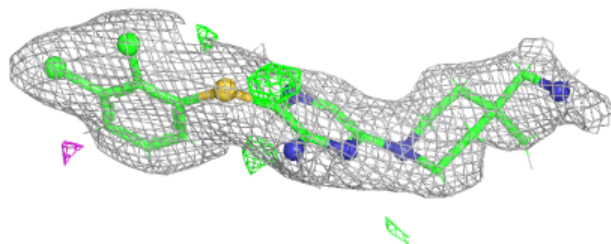
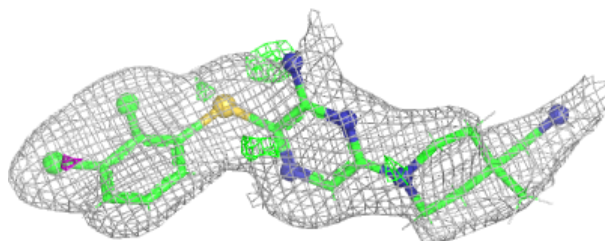
**Electron density around 83Q C 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 83Q D 601:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.