



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:50 AM UTC

PDB ID : 6Q21 / pdb_00006q21
Title : MOLECULAR SWITCH FOR SIGNAL TRANSDUCTION: STRUCTURAL DIFFERENCES BETWEEN ACTIVE AND INACTIVE FORMS OF PROTOONCOGENIC RAS PROTEINS
Authors : Kim, S.-H.
Deposited on : 1992-07-27
Resolution : 1.95 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

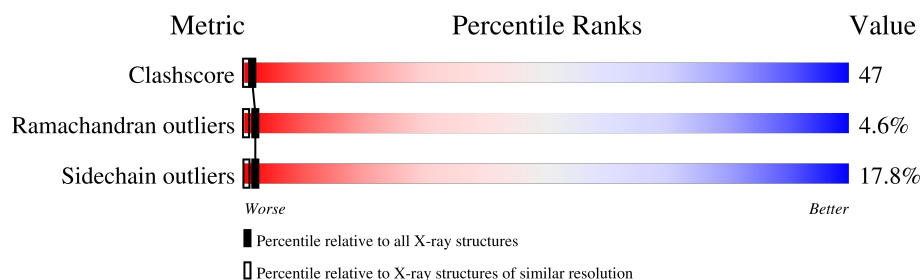
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	171	
1	B	171	
1	C	171	
1	D	171	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6057 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.

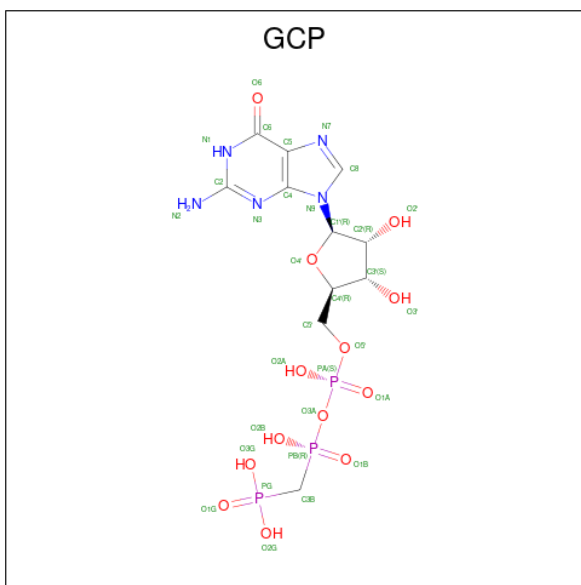
- Molecule 1 is a protein called C-H-RAS P21 PROTEIN CATALYTIC DOMAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1368	854	238	269	7			
1	B	171	Total	C	N	O	S	0	0	0
			1368	854	238	269	7			
1	C	171	Total	C	N	O	S	0	0	0
			1368	854	238	269	7			
1	D	171	Total	C	N	O	S	0	0	0
			1368	854	238	269	7			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PHOSPHOMETHYLPHOSPHONIC ACID GUANYLATE ESTER (CCD ID: GCP) (formula: C₁₁H₁₈N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 32	C 11	N 5	O 13	P 3	0	0
3	B	1	Total 32	C 11	N 5	O 13	P 3	0	0
3	C	1	Total 32	C 11	N 5	O 13	P 3	0	0
3	D	1	Total 32	C 11	N 5	O 13	P 3	0	0

- Molecule 4 is water.

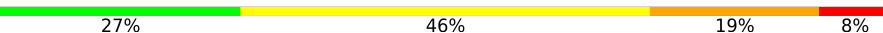
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	83	Total O 83 83	0	0
4	B	93	Total O 93 93	0	0
4	C	111	Total O 111 111	0	0
4	D	166	Total O 166 166	0	0

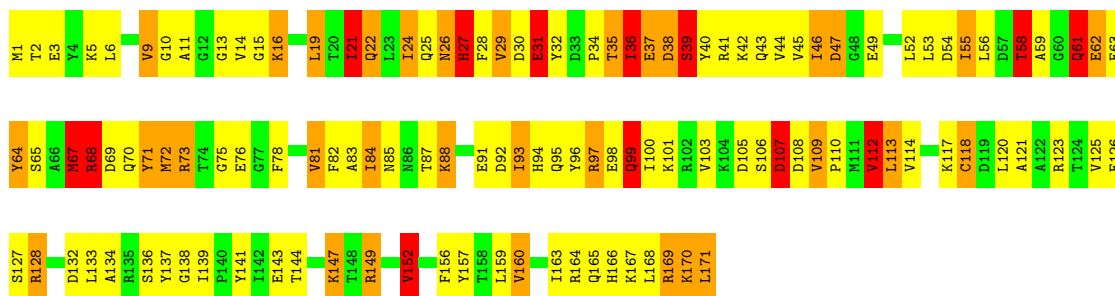
3 Residue-property plots [i](#)

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. 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. 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.

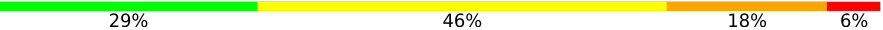
Note EDS was not executed.

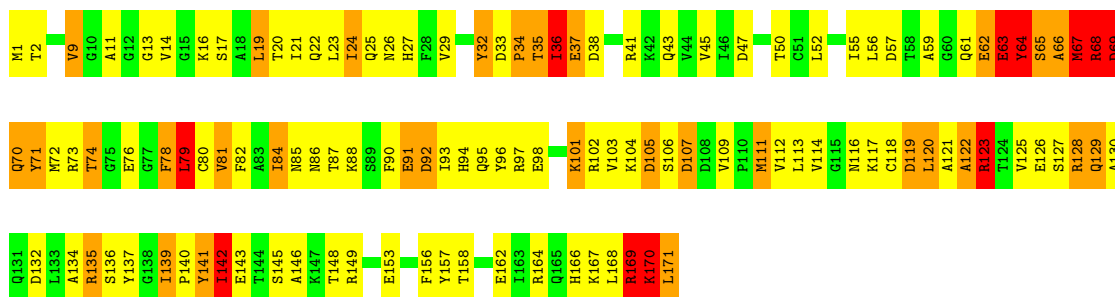
• Molecule 1: C-H-RAS P21 PROTEIN CATALYTIC DOMAIN

Chain A: 

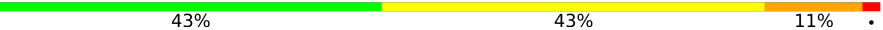


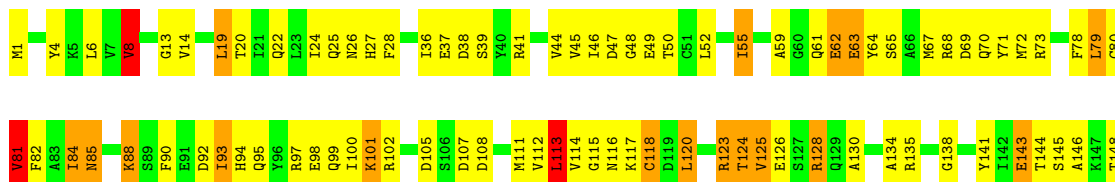
• Molecule 1: C-H-RAS P21 PROTEIN CATALYTIC DOMAIN

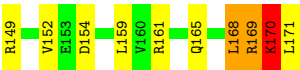
Chain B: 



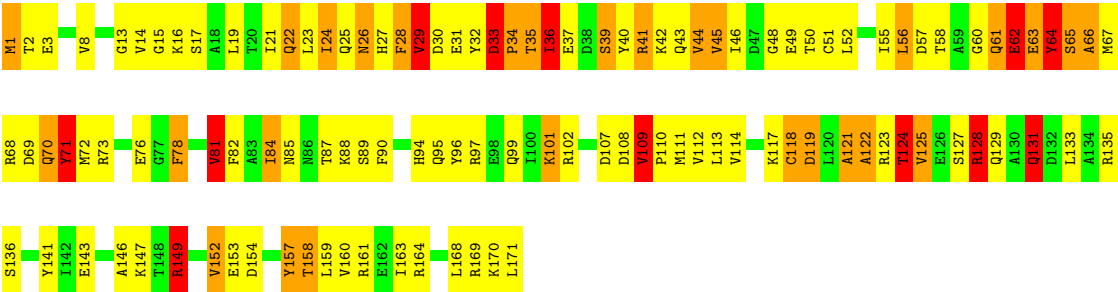
• Molecule 1: C-H-RAS P21 PROTEIN CATALYTIC DOMAIN

Chain C: 





● Molecule 1: C-H-RAS P21 PROTEIN CATALYTIC DOMAIN



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 1 21	Depositor
Cell constants a, b, c, α , β , γ	41.98Å 79.12Å 128.50Å 90.00° 90.00° 117.48°	Depositor
Resolution (Å)	(Not available) – 1.95	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.95)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.190 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6057	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.18	0/1387	1.98	44/1868 (2.4%)
1	B	1.20	1/1387 (0.1%)	1.94	42/1868 (2.2%)
1	C	1.30	2/1387 (0.1%)	1.86	28/1868 (1.5%)
1	D	1.30	3/1387 (0.2%)	2.01	50/1868 (2.7%)
All	All	1.24	6/5548 (0.1%)	1.95	164/7472 (2.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	2	0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	118	CYS	CA-CB	8.55	1.68	1.53
1	B	141	TYR	CA-C	-6.13	1.44	1.52
1	C	113	LEU	CA-C	-5.74	1.45	1.52
1	D	45	VAL	CA-C	-5.57	1.45	1.52
1	C	118	CYS	CA-CB	5.32	1.62	1.53
1	D	109	VAL	CA-C	-5.08	1.48	1.52

All (164) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	107	ASP	CA-CB-CG	-15.05	97.55	112.60
1	A	107	ASP	N-CA-CB	11.03	127.13	110.70
1	A	72	MET	N-CA-C	-10.32	95.21	110.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	109	VAL	CB-CA-C	9.69	119.94	110.16
1	B	170	LYS	N-CA-C	-9.61	98.84	111.71
1	B	67	MET	N-CA-C	-9.03	100.58	111.69
1	A	55	ILE	N-CA-C	8.85	120.69	107.51
1	A	149	ARG	N-CA-CB	8.83	125.41	110.49
1	D	44	VAL	CA-C-N	-8.81	111.59	123.14
1	D	44	VAL	C-N-CA	-8.81	111.59	123.14
1	D	66	ALA	N-CA-C	-8.68	96.70	109.63
1	D	119	ASP	CA-CB-CG	8.32	120.92	112.60
1	D	62	GLU	N-CA-C	-8.27	99.19	110.68
1	B	168	LEU	N-CA-CB	-8.25	97.79	111.74
1	C	8	VAL	CB-CA-C	8.16	122.16	110.98
1	B	169	ARG	CA-C-N	8.06	137.44	121.81
1	B	169	ARG	C-N-CA	8.06	137.44	121.81
1	D	152	VAL	N-CA-C	7.90	118.68	110.62
1	D	119	ASP	N-CA-C	7.82	119.81	111.28
1	C	125	VAL	CB-CA-C	7.65	120.94	110.99
1	A	64	TYR	N-CA-C	7.42	126.61	110.80
1	B	169	ARG	N-CA-CB	7.30	122.83	110.49
1	C	93	ILE	N-CA-CB	7.26	121.93	110.54
1	C	55	ILE	N-CA-C	7.17	118.45	107.99
1	B	36	ILE	N-CA-C	-7.16	104.87	111.67
1	C	14	VAL	N-CA-C	7.14	120.02	111.09
1	B	94	HIS	CA-CB-CG	-7.12	106.68	113.80
1	B	70	GLN	N-CA-CB	7.12	121.55	110.44
1	B	57	ASP	CA-CB-CG	-7.07	105.53	112.60
1	C	81	VAL	N-CA-C	7.04	119.59	108.89
1	D	28	PHE	CA-C-N	-6.99	113.95	122.90
1	D	28	PHE	C-N-CA	-6.99	113.95	122.90
1	B	24	ILE	N-CA-C	6.91	117.66	110.62
1	D	33	ASP	CB-CA-C	6.89	123.75	110.17
1	D	121	ALA	N-CA-C	-6.83	98.12	109.46
1	B	169	ARG	O-C-N	6.72	131.53	122.59
1	B	111	MET	N-CA-CB	-6.70	100.47	111.66
1	D	64	TYR	O-C-N	6.68	131.48	122.59
1	C	27	HIS	CA-CB-CG	-6.62	107.18	113.80
1	C	124	THR	N-CA-CB	-6.62	101.49	111.15
1	A	28	PHE	N-CA-C	6.61	120.28	109.76
1	C	28	PHE	CA-CB-CG	-6.59	107.21	113.80
1	B	119	ASP	CA-CB-CG	-6.53	106.07	112.60
1	C	28	PHE	CA-C-N	-6.51	114.66	123.12
1	C	28	PHE	C-N-CA	-6.51	114.66	123.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	ILE	CB-CA-C	6.50	121.55	112.05
1	B	107	ASP	CA-C-N	-6.50	113.78	122.42
1	B	107	ASP	C-N-CA	-6.50	113.78	122.42
1	B	139	ILE	O-C-N	6.45	125.76	121.69
1	B	67	MET	N-CA-CB	6.41	119.75	110.20
1	A	71	TYR	CB-CA-C	6.40	123.15	110.42
1	A	67	MET	N-CA-C	6.33	124.29	110.80
1	B	19	LEU	O-C-N	6.33	128.62	122.03
1	A	39	SER	N-CA-C	6.33	120.01	109.95
1	A	34	PRO	CA-C-N	6.32	133.61	121.54
1	A	34	PRO	C-N-CA	6.32	133.61	121.54
1	D	125	VAL	CB-CA-C	6.31	119.19	110.99
1	C	55	ILE	N-CA-CB	-6.28	103.86	111.21
1	A	170	LYS	CA-C-N	6.27	132.98	121.70
1	A	170	LYS	C-N-CA	6.27	132.98	121.70
1	D	131	GLN	CB-CG-CD	6.27	123.25	112.60
1	B	123	ARG	CB-CA-C	-6.23	99.25	109.53
1	A	29	VAL	N-CA-C	6.23	117.70	107.24
1	D	81	VAL	N-CA-C	6.21	118.33	108.89
1	D	62	GLU	N-CA-CB	6.20	118.26	110.45
1	A	31	GLU	N-CA-CB	-6.19	102.12	111.15
1	C	27	HIS	N-CA-CB	6.18	121.94	110.99
1	D	124	THR	CA-C-N	-6.17	114.86	123.13
1	D	124	THR	C-N-CA	-6.17	114.86	123.13
1	B	63	GLU	O-C-N	6.16	131.20	123.19
1	C	38	ASP	CA-CB-CG	-6.10	106.50	112.60
1	B	91	GLU	N-CA-C	6.07	118.86	111.82
1	A	112	VAL	CB-CA-C	6.06	119.18	110.33
1	A	152	VAL	N-CA-C	6.03	116.81	110.72
1	A	112	VAL	N-CA-CB	6.02	120.31	111.52
1	D	26	ASN	CA-CB-CG	-5.99	106.61	112.60
1	A	36	ILE	N-CA-CB	5.92	121.01	111.23
1	B	71	TYR	N-CA-C	-5.88	105.60	112.89
1	A	58	THR	N-CA-CB	-5.86	102.29	110.38
1	B	119	ASP	CA-C-O	5.85	126.14	119.18
1	A	62	GLU	N-CA-C	5.84	117.91	107.80
1	A	110	PRO	N-CA-CB	5.84	107.14	103.23
1	B	71	TYR	N-CA-CB	5.84	119.92	110.40
1	A	61	GLN	N-CA-C	5.83	123.21	110.80
1	B	74	THR	N-CA-CB	-5.82	101.52	110.20
1	B	78	PHE	N-CA-C	5.82	118.44	109.07
1	B	64	TYR	CB-CA-C	5.81	121.97	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	19	LEU	CB-CA-C	-5.79	101.75	110.90
1	D	71	TYR	CA-C-N	-5.77	111.97	120.28
1	D	71	TYR	C-N-CA	-5.77	111.97	120.28
1	D	70	GLN	CA-C-N	5.76	128.47	120.29
1	D	70	GLN	C-N-CA	5.76	128.47	120.29
1	A	121	ALA	N-CA-C	-5.75	103.40	110.65
1	A	81	VAL	N-CA-C	5.74	117.86	108.97
1	D	64	TYR	CA-C-N	5.71	132.44	121.54
1	D	64	TYR	C-N-CA	5.71	132.44	121.54
1	A	71	TYR	N-CA-CB	5.68	120.10	110.49
1	D	157	TYR	CA-C-N	-5.66	111.28	120.72
1	D	157	TYR	C-N-CA	-5.66	111.28	120.72
1	C	82	PHE	CA-C-N	-5.64	114.75	122.93
1	C	82	PHE	C-N-CA	-5.64	114.75	122.93
1	D	114	VAL	N-CA-C	5.64	115.99	107.75
1	D	78	PHE	CA-CB-CG	-5.62	108.18	113.80
1	B	79	LEU	N-CA-CB	5.60	119.11	110.21
1	B	76	GLU	CB-CG-CD	-5.59	103.10	112.60
1	C	19	LEU	N-CA-CB	5.58	118.16	110.07
1	B	68	ARG	N-CA-C	-5.58	98.92	110.80
1	A	47	ASP	CA-CB-CG	-5.56	107.04	112.60
1	A	27	HIS	CA-C-N	5.54	129.43	121.50
1	A	27	HIS	C-N-CA	5.54	129.43	121.50
1	D	149	ARG	CD-NE-CZ	5.51	132.12	124.40
1	B	142	ILE	CA-C-O	5.51	126.54	120.64
1	C	154	ASP	CB-CA-C	-5.48	102.27	110.88
1	B	92	ASP	CB-CA-C	5.46	120.81	109.95
1	A	19	LEU	O-C-N	5.45	127.70	122.03
1	B	65	SER	N-CA-C	-5.42	101.83	109.96
1	D	128	ARG	NE-CZ-NH2	5.41	124.07	119.20
1	A	84	ILE	N-CA-CB	-5.41	99.70	111.05
1	A	35	THR	CA-C-N	5.40	131.70	121.97
1	A	35	THR	C-N-CA	5.40	131.70	121.97
1	C	62	GLU	N-CA-C	-5.40	101.37	109.15
1	C	143	GLU	N-CA-CB	5.40	119.23	110.43
1	D	111	MET	CA-C-N	-5.39	115.45	122.94
1	D	111	MET	C-N-CA	-5.39	115.45	122.94
1	D	26	ASN	N-CA-C	5.39	119.80	113.23
1	B	35	THR	CA-C-N	-5.36	116.26	122.63
1	B	35	THR	C-N-CA	-5.36	116.26	122.63
1	D	128	ARG	CB-CG-CD	5.35	123.61	111.30
1	B	45	VAL	CA-C-N	-5.35	115.38	122.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	VAL	C-N-CA	-5.35	115.38	122.98
1	D	108	ASP	CA-CB-CG	-5.34	107.26	112.60
1	C	94	HIS	CA-CB-CG	-5.33	108.47	113.80
1	D	84	ILE	CA-CB-CG2	5.32	119.55	110.50
1	D	33	ASP	CA-C-N	5.29	126.04	120.11
1	D	33	ASP	C-N-CA	5.29	126.04	120.11
1	A	109	VAL	N-CA-C	5.29	120.31	108.88
1	D	45	VAL	CA-C-N	-5.29	115.48	122.98
1	D	45	VAL	C-N-CA	-5.29	115.48	122.98
1	B	169	ARG	CA-C-O	5.27	128.05	120.51
1	D	50	THR	N-CA-CB	5.25	119.00	110.43
1	D	29	VAL	N-CA-CB	5.25	117.35	111.21
1	D	64	TYR	CA-C-O	5.22	127.97	120.51
1	A	93	ILE	CB-CA-C	-5.19	105.23	111.88
1	C	81	VAL	CB-CA-C	5.18	119.30	110.95
1	C	90	PHE	N-CA-C	-5.15	105.67	111.28
1	D	36	ILE	N-CA-CB	5.15	116.87	111.00
1	A	132	ASP	CA-CB-CG	5.13	117.73	112.60
1	B	70	GLN	CB-CG-CD	5.12	121.30	112.60
1	B	142	ILE	CB-CA-C	5.12	119.65	110.65
1	B	107	ASP	N-CA-CB	5.10	117.97	110.33
1	D	34	PRO	CA-N-CD	-5.09	104.88	112.00
1	C	125	VAL	N-CA-CB	5.08	118.30	111.90
1	C	80	CYS	CA-C-N	-5.08	115.06	122.68
1	C	80	CYS	C-N-CA	-5.08	115.06	122.68
1	A	99	GLN	CB-CA-C	-5.08	102.22	110.85
1	C	123	ARG	N-CA-CB	5.08	117.50	110.04
1	D	1	MET	CA-C-N	-5.07	115.58	122.93
1	D	1	MET	C-N-CA	-5.07	115.58	122.93
1	A	170	LYS	O-C-N	5.06	128.35	121.97
1	A	71	TYR	N-CA-C	-5.05	100.03	110.80
1	A	46	ILE	CA-C-N	-5.03	115.10	122.19
1	A	46	ILE	C-N-CA	-5.03	115.10	122.19
1	A	170	LYS	N-CA-C	-5.01	103.09	113.31
1	D	157	TYR	N-CA-CB	5.00	118.69	110.39

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	64	TYR	CA
1	A	107	ASP	CA

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	1368	0	1357	136	0
1	B	1368	0	1355	156	0
1	C	1368	0	1355	73	0
1	D	1368	0	1356	162	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	32	0	14	6	0
3	B	32	0	14	2	0
3	C	32	0	14	4	0
3	D	32	0	14	5	0
4	A	83	0	0	18	0
4	B	93	0	0	14	0
4	C	111	0	0	16	0
4	D	166	0	0	58	0
All	All	6057	0	5479	517	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 47.

All (517) 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:33:ASP:HB3	1:D:34:PRO:HD3	1.18	1.14
1:B:41:ARG:HD2	1:D:135:ARG:HD3	1.17	1.09
1:D:34:PRO:HB2	1:D:61:GLN:HG3	1.38	1.05
1:B:72:MET:HE3	1:B:103:VAL:HG21	1.37	1.02
1:A:126:GLU:HB3	1:A:128:ARG:HE	1.24	1.02
1:D:13:GLY:H	3:D:180:GCP:H3B1	1.27	0.98
1:A:35:THR:HB	1:A:59:ALA:HB2	1.44	0.97
1:C:13:GLY:H	3:C:180:GCP:H3B1	1.31	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:127:SER:H	1:D:128:ARG:HH21	1.04	0.93
1:B:41:ARG:HD2	1:D:135:ARG:CD	1.98	0.92
1:D:157:TYR:O	4:D:1110:HOH:O	1.90	0.90
1:B:114:VAL:HA	4:B:1101:HOH:O	1.69	0.90
1:C:64:TYR:HB3	1:C:67:MET:HE2	1.52	0.88
1:D:127:SER:H	1:D:128:ARG:NH2	1.70	0.88
1:A:127:SER:H	1:A:128:ARG:NH2	1.72	0.87
1:B:72:MET:HE3	1:B:103:VAL:CG2	2.05	0.86
1:A:29:VAL:HG11	1:A:32:TYR:HB2	1.55	0.86
1:D:33:ASP:CB	1:D:34:PRO:HD3	2.05	0.85
1:D:68:ARG:HA	1:D:71:TYR:CZ	2.11	0.85
1:C:63:GLU:N	4:C:1055:HOH:O	2.09	0.84
1:D:82:PHE:HA	4:D:1174:HOH:O	1.76	0.84
1:C:128:ARG:HH11	1:C:128:ARG:HB2	1.42	0.84
1:A:108:ASP:HA	4:A:953:HOH:O	1.77	0.84
1:D:170:LYS:HD2	1:D:171:LEU:H	1.43	0.83
1:B:41:ARG:CD	1:D:135:ARG:HD3	2.06	0.83
1:B:13:GLY:H	3:B:180:GCP:H3B1	1.43	0.82
1:D:160:VAL:HB	4:D:1110:HOH:O	1.79	0.82
1:D:149:ARG:HA	4:D:1206:HOH:O	1.78	0.82
1:D:61:GLN:HB3	4:D:1021:HOH:O	1.79	0.82
1:D:128:ARG:H	1:D:128:ARG:HE	1.24	0.82
1:A:169:ARG:CZ	1:A:170:LYS:HE3	2.09	0.82
1:A:41:ARG:NH2	1:B:135:ARG:HG3	1.95	0.81
1:B:64:TYR:HB3	1:B:67:MET:HE2	1.62	0.81
1:A:68:ARG:HG3	1:A:68:ARG:HH11	1.46	0.81
1:A:126:GLU:HB3	1:A:128:ARG:NE	1.96	0.80
1:B:117:LYS:HB3	1:B:120:LEU:HD22	1.63	0.79
1:B:130:ALA:HB1	1:B:141:TYR:CE1	2.18	0.79
1:A:149:ARG:HB2	1:A:152:VAL:CG1	2.14	0.78
1:A:30:ASP:C	1:A:31:GLU:HG3	2.09	0.78
1:B:126:GLU:HB2	1:B:129:GLN:HG2	1.65	0.77
4:A:1210:HOH:O	1:B:135:ARG:HD2	1.84	0.77
1:A:101:LYS:NZ	1:A:107:ASP:HB2	1.99	0.77
1:A:126:GLU:CB	1:A:128:ARG:HE	1.96	0.77
1:C:117:LYS:NZ	4:C:1100:HOH:O	2.18	0.77
1:D:67:MET:HA	4:D:913:HOH:O	1.83	0.77
1:A:41:ARG:HD2	1:B:135:ARG:NH2	2.00	0.76
1:D:33:ASP:HB3	1:D:34:PRO:CD	2.08	0.76
1:D:62:GLU:HB2	4:D:1022:HOH:O	1.85	0.76
1:D:152:VAL:HG21	4:D:1206:HOH:O	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:97:ARG:HD3	4:D:865:HOH:O	1.86	0.75
1:A:108:ASP:HB3	1:A:166:HIS:NE2	2.02	0.75
1:C:118:CYS:SG	4:C:1053:HOH:O	2.44	0.75
1:A:2:THR:OG1	1:B:169:ARG:NH2	2.19	0.75
1:D:67:MET:HG3	4:D:1220:HOH:O	1.86	0.74
1:C:116:ASN:OD1	1:C:144:THR:HG23	1.87	0.74
1:D:13:GLY:N	3:D:180:GCP:H3B1	2.03	0.73
1:B:69:ASP:O	1:B:73:ARG:HG3	1.88	0.73
1:A:147:LYS:HB2	1:A:147:LYS:NZ	2.03	0.73
1:A:149:ARG:HB2	1:A:152:VAL:HG11	1.70	0.73
1:D:63:GLU:HG2	1:D:64:TYR:CD1	2.24	0.73
1:A:61:GLN:HG3	1:A:62:GLU:HG3	1.70	0.73
1:C:84:ILE:HG23	1:C:123:ARG:HB2	1.70	0.73
1:A:94:HIS:O	1:A:98:GLU:HG2	1.88	0.73
1:A:82:PHE:HB3	1:A:93:ILE:HD11	1.70	0.73
1:D:66:ALA:HB2	4:D:1237:HOH:O	1.88	0.72
1:A:147:LYS:HB2	1:A:147:LYS:HZ3	1.54	0.72
1:A:70:GLN:HG2	1:A:73:ARG:HE	1.54	0.72
1:B:90:PHE:HZ	1:B:130:ALA:HA	1.54	0.71
1:D:8:VAL:HG11	4:D:1016:HOH:O	1.90	0.71
1:D:95:GLN:O	1:D:99:GLN:HG3	1.89	0.71
1:B:64:TYR:O	1:B:67:MET:HG2	1.90	0.71
1:B:72:MET:CB	1:B:103:VAL:HG21	2.20	0.71
1:C:168:LEU:O	1:C:170:LYS:N	2.17	0.71
1:C:165:GLN:NE2	1:C:169:ARG:HH21	1.88	0.71
1:A:41:ARG:HD2	1:B:135:ARG:CZ	2.19	0.71
1:A:35:THR:HG23	1:A:38:ASP:OD1	1.89	0.71
1:C:39:SER:HB3	1:C:41:ARG:NH2	2.05	0.71
1:D:133:LEU:CB	4:D:1191:HOH:O	2.38	0.70
1:B:20:THR:O	1:B:24:ILE:HG12	1.90	0.70
1:C:6:LEU:HG	1:C:159:LEU:HD23	1.73	0.70
1:A:35:THR:HG22	1:A:36:ILE:N	2.07	0.70
1:B:67:MET:O	1:B:69:ASP:N	2.25	0.69
1:C:48:GLY:N	4:C:1149:HOH:O	2.16	0.69
1:B:43:GLN:HE22	1:D:141:TYR:H	1.39	0.69
1:A:16:LYS:NZ	3:A:180:GCP:O2G	2.26	0.69
1:D:68:ARG:HD2	1:D:71:TYR:OH	1.92	0.69
1:B:37:GLU:CG	1:B:56:LEU:HD11	2.22	0.69
1:B:37:GLU:HG2	1:B:56:LEU:HD11	1.75	0.69
1:A:24:ILE:HG22	1:A:25:GLN:N	2.08	0.68
1:D:72:MET:HE1	1:D:99:GLN:NE2	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:LEU:O	4:B:1101:HOH:O	2.11	0.68
1:D:164:ARG:HD2	4:D:1065:HOH:O	1.93	0.68
1:B:142:ILE:HD12	1:B:143:GLU:N	2.08	0.68
1:A:30:ASP:HB2	4:A:1127:HOH:O	1.92	0.68
1:A:156:PHE:O	1:A:160:VAL:HG12	1.93	0.68
1:A:81:VAL:HG11	4:A:1141:HOH:O	1.93	0.68
1:B:134:ALA:HB1	1:B:139:ILE:O	1.94	0.68
1:B:166:HIS:HD2	1:B:170:LYS:NZ	1.91	0.67
1:B:166:HIS:CD2	1:B:170:LYS:HG2	2.30	0.67
1:D:128:ARG:H	1:D:128:ARG:NE	1.90	0.67
1:D:170:LYS:HD2	1:D:171:LEU:N	2.08	0.67
1:C:47:ASP:OD1	1:C:161:ARG:NH2	2.25	0.67
1:A:35:THR:CB	1:A:59:ALA:HB2	2.24	0.67
1:A:94:HIS:ND1	1:A:95:GLN:OE1	2.25	0.67
1:B:69:ASP:HB3	1:B:73:ARG:HH12	1.59	0.67
1:B:72:MET:CE	1:B:103:VAL:HG11	2.25	0.66
1:C:24:ILE:HD11	1:C:55:ILE:HD12	1.78	0.66
1:A:45:VAL:HA	1:A:49:GLU:O	1.95	0.66
1:D:127:SER:N	1:D:128:ARG:HH21	1.85	0.66
1:D:89:SER:HB2	4:D:1174:HOH:O	1.95	0.65
1:D:133:LEU:HB3	4:D:1191:HOH:O	1.94	0.65
1:B:72:MET:HB2	1:B:103:VAL:HG21	1.78	0.65
1:B:17:SER:O	1:B:21:ILE:HG12	1.97	0.65
1:D:49:GLU:OE1	1:D:164:ARG:NH2	2.27	0.65
1:A:16:LYS:NZ	3:A:180:GCP:O2B	2.30	0.64
1:D:67:MET:O	1:D:70:GLN:HG3	1.98	0.64
1:C:84:ILE:HG23	1:C:123:ARG:CB	2.27	0.64
1:D:61:GLN:HB2	4:D:813:HOH:O	1.98	0.64
1:D:17:SER:O	1:D:21:ILE:HD13	1.98	0.64
1:D:37:GLU:HG3	1:D:58:THR:HA	1.79	0.64
1:A:11:ALA:O	1:A:14:VAL:HG13	1.98	0.64
1:C:128:ARG:NH1	4:C:1140:HOH:O	2.30	0.64
1:D:147:LYS:N	4:D:1059:HOH:O	2.30	0.64
1:D:169:ARG:HD3	1:D:169:ARG:N	2.12	0.63
1:A:35:THR:HA	4:A:1242:HOH:O	1.98	0.63
1:C:100:ILE:HG21	1:C:111:MET:HE2	1.79	0.63
1:D:61:GLN:NE2	4:D:1042:HOH:O	2.30	0.63
1:B:149:ARG:HH12	1:B:153:GLU:HG3	1.64	0.63
1:C:88:LYS:HE2	1:C:92:ASP:OD2	1.98	0.63
1:B:123:ARG:NE	1:B:125:VAL:O	2.24	0.62
1:D:63:GLU:HG2	1:D:64:TYR:N	2.11	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:24:ILE:HG22	1:D:25:GLN:N	2.14	0.62
1:C:20:THR:O	1:C:24:ILE:HG12	1.99	0.62
1:D:112:VAL:HG23	1:D:159:LEU:HD13	1.80	0.62
1:B:142:ILE:HD12	1:B:143:GLU:H	1.61	0.62
1:D:35:THR:HB	4:D:1015:HOH:O	1.99	0.62
1:A:45:VAL:HG23	1:A:49:GLU:O	1.99	0.62
1:B:145:SER:HB3	1:B:148:THR:OG1	1.99	0.62
1:A:44:VAL:HG22	1:A:45:VAL:N	2.15	0.62
1:A:169:ARG:NH2	1:A:170:LYS:HE3	2.15	0.62
1:C:134:ALA:O	1:C:138:GLY:N	2.33	0.61
1:A:58:THR:HG23	1:A:59:ALA:N	2.15	0.61
1:A:168:LEU:HD12	4:A:1240:HOH:O	2.00	0.61
1:D:42:LYS:HE3	4:D:889:HOH:O	2.01	0.61
1:D:161:ARG:N	4:D:1110:HOH:O	2.10	0.61
1:B:72:MET:HB3	1:B:103:VAL:HG21	1.83	0.60
1:B:90:PHE:HD2	1:B:125:VAL:CG2	2.13	0.60
1:B:166:HIS:HA	1:B:170:LYS:HB2	1.83	0.60
1:A:61:GLN:HA	1:A:61:GLN:OE1	2.00	0.60
1:D:28:PHE:CD2	1:D:147:LYS:HA	2.36	0.60
1:D:123:ARG:HH22	1:D:143:GLU:CD	2.09	0.60
1:B:171:LEU:HB3	4:B:1238:HOH:O	2.01	0.60
1:C:69:ASP:OD1	1:C:73:ARG:NH2	2.34	0.60
1:D:48:GLY:N	4:D:916:HOH:O	2.32	0.60
1:D:170:LYS:HB3	4:D:1160:HOH:O	2.02	0.60
1:C:1:MET:HB2	1:C:50:THR:O	2.02	0.60
1:D:101:LYS:HE3	1:D:107:ASP:OD1	2.01	0.60
1:D:90:PHE:CZ	4:D:1191:HOH:O	2.52	0.60
1:A:149:ARG:O	1:A:152:VAL:HG12	2.02	0.59
1:D:44:VAL:HG22	1:D:45:VAL:N	2.17	0.59
1:A:5:LYS:O	1:A:76:GLU:HB2	2.03	0.59
1:A:68:ARG:HE	1:A:96:TYR:HE1	1.50	0.59
1:A:26:ASN:O	1:A:26:ASN:ND2	2.35	0.59
1:B:61:GLN:HA	4:B:1076:HOH:O	2.01	0.59
1:A:29:VAL:HG12	1:A:29:VAL:O	2.02	0.59
1:C:161:ARG:NH1	4:C:1092:HOH:O	2.27	0.59
1:D:131:GLN:NE2	1:D:135:ARG:NH2	2.51	0.59
1:D:21:ILE:O	1:D:25:GLN:N	2.36	0.58
1:D:15:GLY:O	1:D:19:LEU:HB2	2.03	0.58
1:D:34:PRO:HB2	1:D:61:GLN:CG	2.23	0.58
1:B:167:LYS:HA	1:B:167:LYS:HE2	1.86	0.58
1:B:13:GLY:N	3:B:180:GCP:H3B1	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:21:ILE:HG21	1:D:29:VAL:HG13	1.85	0.58
1:C:52:LEU:HD13	1:C:52:LEU:C	2.29	0.58
1:B:90:PHE:CZ	1:B:130:ALA:HA	2.38	0.58
1:D:2:THR:HB	1:D:49:GLU:OE2	2.04	0.57
1:B:97:ARG:HD3	1:B:137:TYR:CZ	2.40	0.57
1:D:94:HIS:HB2	4:D:837:HOH:O	2.03	0.57
1:A:41:ARG:CZ	1:B:135:ARG:HG3	2.34	0.57
1:C:130:ALA:HB1	1:C:141:TYR:CE1	2.39	0.57
1:D:22:GLN:NE2	1:D:149:ARG:HD3	2.19	0.57
1:A:29:VAL:CG1	1:A:32:TYR:HB2	2.30	0.57
1:D:62:GLU:HA	1:D:62:GLU:OE1	2.04	0.57
1:A:127:SER:H	1:A:128:ARG:HH21	1.51	0.57
1:B:116:ASN:O	1:B:117:LYS:HB2	2.04	0.57
1:D:3:GLU:HG2	1:D:52:LEU:HB3	1.85	0.57
4:A:971:HOH:O	1:C:45:VAL:HG21	2.05	0.57
1:A:101:LYS:HZ2	1:A:107:ASP:HB2	1.68	0.56
1:C:59:ALA:HB1	4:C:808:HOH:O	2.04	0.56
1:D:85:ASN:HA	1:D:124:THR:HB	1.85	0.56
1:B:41:ARG:HB3	1:D:135:ARG:CZ	2.35	0.56
1:B:142:ILE:N	4:B:1101:HOH:O	2.14	0.56
1:A:70:GLN:HG2	1:A:73:ARG:NE	2.20	0.56
1:B:78:PHE:HB2	4:B:951:HOH:O	2.05	0.56
1:B:101:LYS:HE2	1:B:107:ASP:OD1	2.05	0.56
1:C:61:GLN:HB2	4:C:808:HOH:O	2.05	0.56
1:B:86:ASN:OD1	1:B:88:LYS:N	2.38	0.56
1:B:81:VAL:HG12	1:B:82:PHE:N	2.21	0.56
1:D:22:GLN:NE2	1:D:149:ARG:CD	2.69	0.56
1:D:22:GLN:HG3	1:D:146:ALA:O	2.06	0.56
1:A:171:LEU:HD23	4:A:1240:HOH:O	2.06	0.56
1:B:126:GLU:HB2	1:B:129:GLN:CG	2.34	0.56
1:C:126:GLU:HA	1:C:126:GLU:OE1	2.06	0.56
1:D:34:PRO:CB	1:D:61:GLN:HG3	2.24	0.56
1:A:123:ARG:NE	4:A:1201:HOH:O	2.28	0.56
1:A:67:MET:HB2	4:A:1232:HOH:O	2.04	0.55
1:A:169:ARG:NH1	4:A:1106:HOH:O	2.39	0.55
1:C:13:GLY:O	4:C:1100:HOH:O	2.18	0.55
1:D:69:ASP:O	1:D:73:ARG:HB2	2.06	0.55
1:A:21:ILE:HD11	1:A:29:VAL:HG23	1.89	0.55
1:B:62:GLU:HG2	1:B:63:GLU:HG3	1.89	0.55
1:B:112:VAL:HG22	1:B:140:PRO:HG2	1.89	0.55
1:D:128:ARG:HE	1:D:128:ARG:N	2.00	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:GLN:O	1:B:63:GLU:N	2.40	0.54
1:D:136:SER:HA	4:D:1108:HOH:O	2.05	0.54
1:B:69:ASP:HA	1:B:72:MET:HE2	1.89	0.54
1:D:22:GLN:HE21	1:D:22:GLN:CA	2.20	0.54
1:A:72:MET:O	1:A:73:ARG:HB2	2.05	0.54
1:A:118:CYS:SG	4:A:1061:HOH:O	2.55	0.54
1:C:117:LYS:HG2	3:C:180:GCP:C6	2.37	0.54
1:A:29:VAL:CG1	1:A:32:TYR:H	2.20	0.54
1:B:127:SER:OG	1:B:128:ARG:NH2	2.40	0.54
1:A:24:ILE:C	1:A:26:ASN:H	2.16	0.54
1:B:90:PHE:HD2	1:B:125:VAL:HG22	1.72	0.54
1:B:81:VAL:HG12	1:B:82:PHE:H	1.74	0.53
1:B:145:SER:O	1:B:149:ARG:N	2.40	0.53
1:C:149:ARG:NH2	4:C:1048:HOH:O	2.40	0.53
1:D:97:ARG:NH1	1:D:101:LYS:HD3	2.22	0.53
1:B:157:TYR:CD1	1:B:157:TYR:N	2.76	0.53
1:D:127:SER:N	1:D:128:ARG:HE	2.07	0.53
1:A:13:GLY:H	3:A:180:GCP:H3B1	1.73	0.53
1:A:24:ILE:HG22	1:A:25:GLN:CD	2.33	0.53
1:A:44:VAL:HG22	1:A:45:VAL:H	1.71	0.53
1:C:13:GLY:N	3:C:180:GCP:H3B1	2.12	0.53
1:B:84:ILE:C	1:B:84:ILE:HD12	2.33	0.53
1:D:122:ALA:N	4:D:1073:HOH:O	2.24	0.53
1:B:104:LYS:O	1:B:105:ASP:HB2	2.09	0.53
1:D:146:ALA:HB3	4:D:1059:HOH:O	2.08	0.53
1:C:36:ILE:O	1:C:37:GLU:HB2	2.08	0.53
1:D:28:PHE:HA	4:D:901:HOH:O	2.09	0.53
1:D:25:GLN:HB2	1:D:27:HIS:CD2	2.44	0.53
1:D:43:GLN:HA	1:D:51:CYS:O	2.09	0.53
1:A:35:THR:CG2	1:A:36:ILE:N	2.72	0.53
1:B:61:GLN:NE2	1:B:64:TYR:CE2	2.77	0.52
1:A:9:VAL:HG13	1:A:10:GLY:N	2.23	0.52
1:A:105:ASP:OD1	1:A:105:ASP:O	2.27	0.52
1:C:128:ARG:HH11	1:C:128:ARG:CB	2.19	0.52
1:C:101:LYS:O	1:C:105:ASP:N	2.42	0.52
1:A:11:ALA:O	1:A:14:VAL:HG22	2.09	0.52
1:B:66:ALA:O	1:B:69:ASP:HB2	2.10	0.52
1:D:63:GLU:HA	4:D:1193:HOH:O	2.10	0.52
1:B:149:ARG:HH12	1:B:153:GLU:CG	2.23	0.52
1:C:168:LEU:C	1:C:170:LYS:H	2.14	0.52
1:D:70:GLN:O	1:D:73:ARG:HB3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:SER:C	1:A:40:TYR:CD1	2.88	0.52
1:C:145:SER:O	1:C:149:ARG:N	2.40	0.52
1:B:78:PHE:CD1	1:B:78:PHE:N	2.78	0.52
1:B:166:HIS:HD2	1:B:170:LYS:HZ3	1.57	0.51
1:B:33:ASP:O	1:B:34:PRO:O	2.28	0.51
1:B:85:ASN:ND2	1:B:120:LEU:HD23	2.25	0.51
1:C:111:MET:C	1:C:112:VAL:HG23	2.35	0.51
1:D:121:ALA:HA	4:D:1073:HOH:O	2.09	0.51
1:B:111:MET:HG3	4:B:951:HOH:O	2.09	0.51
1:C:20:THR:OG1	4:C:1058:HOH:O	2.19	0.51
1:C:85:ASN:HD22	1:C:85:ASN:H	1.58	0.51
1:A:68:ARG:HG3	1:A:68:ARG:NH1	2.22	0.51
1:D:16:LYS:NZ	3:D:180:GCP:O1G	2.35	0.51
1:B:111:MET:O	1:B:112:VAL:HG23	2.10	0.51
1:A:103:VAL:HG12	1:A:103:VAL:O	2.10	0.51
1:B:47:ASP:OD2	1:B:164:ARG:NH1	2.41	0.51
1:C:72:MET:SD	1:C:99:GLN:HG2	2.51	0.51
1:D:27:HIS:CD2	1:D:27:HIS:N	2.78	0.51
1:B:126:GLU:O	1:B:129:GLN:HG3	2.11	0.51
1:D:90:PHE:CE1	4:D:1191:HOH:O	2.64	0.50
1:A:113:LEU:O	1:A:141:TYR:HA	2.11	0.50
1:C:113:LEU:HB3	1:C:141:TYR:HD1	1.77	0.50
1:A:88:LYS:O	1:A:92:ASP:OD2	2.30	0.50
1:D:127:SER:H	1:D:128:ARG:CZ	2.22	0.50
1:C:65:SER:HB3	4:C:1055:HOH:O	2.12	0.50
1:B:117:LYS:C	1:B:119:ASP:H	2.20	0.49
1:A:46:ILE:HD11	1:A:157:TYR:CD1	2.46	0.49
1:A:137:TYR:HB2	1:A:139:ILE:HG12	1.93	0.49
1:D:170:LYS:O	1:D:171:LEU:HB2	2.11	0.49
1:B:102:ARG:HD2	4:B:1105:HOH:O	2.11	0.49
1:B:145:SER:HB3	1:B:148:THR:HG1	1.78	0.49
1:A:112:VAL:HG13	1:A:159:LEU:HB2	1.95	0.49
1:C:70:GLN:NE2	1:C:73:ARG:HH12	2.10	0.49
1:A:83:ALA:HB1	1:A:85:ASN:OD1	2.12	0.49
1:A:123:ARG:HD3	1:A:125:VAL:O	2.12	0.49
1:B:37:GLU:HG3	1:B:56:LEU:HD11	1.92	0.49
1:A:127:SER:N	1:A:128:ARG:NH2	2.52	0.49
1:B:69:ASP:O	1:B:73:ARG:NH1	2.46	0.49
1:D:102:ARG:HD2	4:D:843:HOH:O	2.12	0.49
1:A:36:ILE:HG23	1:A:64:TYR:HB3	1.93	0.49
1:D:141:TYR:OH	1:D:143:GLU:OE2	2.29	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:THR:HG22	1:A:36:ILE:H	1.76	0.48
1:A:22:GLN:O	1:A:26:ASN:HA	2.12	0.48
1:A:108:ASP:HB3	1:A:166:HIS:CE1	2.48	0.48
1:B:117:LYS:O	1:B:119:ASP:N	2.46	0.48
1:A:30:ASP:OD1	1:A:31:GLU:OE2	2.31	0.48
1:B:126:GLU:HB2	1:B:129:GLN:OE1	2.13	0.48
1:B:61:GLN:C	1:B:63:GLU:N	2.72	0.48
1:B:85:ASN:ND2	1:B:120:LEU:CD2	2.76	0.48
1:B:101:LYS:O	1:B:105:ASP:N	2.47	0.48
1:B:2:THR:N	4:B:1085:HOH:O	2.45	0.48
1:D:61:GLN:HB3	1:D:62:GLU:H	1.42	0.48
1:D:68:ARG:HA	1:D:71:TYR:CE1	2.47	0.48
1:B:37:GLU:O	1:B:38:ASP:OD1	2.31	0.48
1:B:121:ALA:O	1:B:122:ALA:HB2	2.13	0.48
1:C:81:VAL:HB	1:C:114:VAL:HB	1.96	0.48
1:D:170:LYS:CG	1:D:171:LEU:N	2.77	0.48
1:B:90:PHE:HD1	1:B:90:PHE:O	1.96	0.47
1:D:154:ASP:O	1:D:158:THR:N	2.45	0.47
1:A:97:ARG:NH2	4:A:974:HOH:O	2.47	0.47
1:A:47:ASP:OD2	1:A:164:ARG:NH1	2.36	0.47
1:A:167:LYS:NZ	4:A:995:HOH:O	2.47	0.47
1:D:129:GLN:HG2	4:D:842:HOH:O	2.14	0.47
1:C:145:SER:HB3	1:C:148:THR:OG1	2.13	0.47
1:D:66:ALA:HA	4:D:1083:HOH:O	2.13	0.47
1:D:123:ARG:NH2	1:D:143:GLU:OE2	2.42	0.47
1:A:46:ILE:O	1:A:47:ASP:HB2	2.13	0.47
1:D:41:ARG:HB2	1:D:41:ARG:CZ	2.44	0.47
1:D:64:TYR:HE2	1:D:96:TYR:CZ	2.33	0.47
1:D:109:VAL:HA	1:D:110:PRO:HD3	1.80	0.47
1:D:39:SER:HA	1:D:55:ILE:O	2.15	0.47
1:B:130:ALA:HB1	1:B:141:TYR:CZ	2.49	0.47
3:D:180:GCP:H3B2	4:D:872:HOH:O	2.14	0.47
1:A:37:GLU:HG2	1:A:56:LEU:HD11	1.96	0.47
1:A:67:MET:N	1:A:67:MET:HE3	2.30	0.47
1:D:60:GLY:HA2	4:D:949:HOH:O	2.15	0.47
1:D:117:LYS:C	1:D:119:ASP:N	2.71	0.47
1:B:170:LYS:O	1:B:171:LEU:HB2	2.15	0.47
1:B:11:ALA:O	1:B:14:VAL:HG13	2.14	0.47
1:B:130:ALA:CB	1:B:141:TYR:CZ	2.98	0.47
1:D:89:SER:CB	4:D:1174:HOH:O	2.58	0.47
1:B:157:TYR:N	1:B:157:TYR:HD1	2.13	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:4:TYR:OH	1:C:49:GLU:OE1	2.24	0.46
1:D:44:VAL:CG2	1:D:45:VAL:N	2.78	0.46
1:A:67:MET:HB3	1:A:71:TYR:HE2	1.80	0.46
1:B:166:HIS:HD2	1:B:170:LYS:HZ2	1.62	0.46
1:C:116:ASN:O	1:C:117:LYS:HB2	2.15	0.46
1:D:28:PHE:CD1	1:D:28:PHE:O	2.68	0.46
1:B:79:LEU:HD13	1:B:114:VAL:HG21	1.97	0.46
1:B:149:ARG:NH1	1:B:153:GLU:OE2	2.48	0.46
1:D:36:ILE:HD12	4:D:1109:HOH:O	2.15	0.46
1:B:149:ARG:NH1	1:B:153:GLU:HG3	2.30	0.46
1:D:170:LYS:CD	1:D:171:LEU:N	2.77	0.46
1:C:78:PHE:O	1:C:111:MET:HA	2.15	0.46
1:B:171:LEU:HD22	1:B:171:LEU:HA	1.44	0.46
1:C:41:ARG:NH2	4:C:863:HOH:O	2.49	0.46
1:D:133:LEU:CD2	4:D:1191:HOH:O	2.64	0.46
1:D:102:ARG:HG2	4:D:843:HOH:O	2.15	0.46
1:B:128:ARG:H	1:B:128:ARG:NE	2.14	0.46
1:B:97:ARG:NH1	1:B:98:GLU:OE1	2.49	0.46
1:C:46:ILE:O	1:C:49:GLU:N	2.44	0.46
1:C:108:ASP:OD2	1:C:169:ARG:NH1	2.47	0.46
1:D:149:ARG:HD2	4:D:1206:HOH:O	2.16	0.46
1:A:21:ILE:HD12	1:A:27:HIS:O	2.16	0.45
1:D:61:GLN:C	1:D:63:GLU:H	2.24	0.45
1:D:160:VAL:N	4:D:1110:HOH:O	2.47	0.45
1:A:164:ARG:NH1	4:A:1064:HOH:O	2.47	0.45
3:C:180:GCP:O4'	4:C:1100:HOH:O	2.21	0.45
1:D:22:GLN:HE21	1:D:22:GLN:HA	1.81	0.45
1:B:36:ILE:HD13	1:B:36:ILE:N	2.31	0.45
1:B:68:ARG:HG3	1:B:71:TYR:OH	2.16	0.45
1:C:85:ASN:OD1	1:C:120:LEU:HD23	2.17	0.45
1:A:81:VAL:HG12	1:A:82:PHE:N	2.31	0.45
1:B:69:ASP:HB3	1:B:73:ARG:NH1	2.30	0.45
1:B:69:ASP:C	1:B:73:ARG:NH1	2.74	0.45
1:D:135:ARG:NH1	4:D:1094:HOH:O	2.28	0.45
1:B:33:ASP:HA	1:B:34:PRO:HD2	1.73	0.45
1:A:117:LYS:N	1:A:144:THR:O	2.50	0.45
1:A:133:LEU:HG	1:A:137:TYR:CE2	2.51	0.45
1:A:169:ARG:C	1:A:170:LYS:HG2	2.41	0.45
1:D:87:THR:OG1	1:D:124:THR:HG23	2.17	0.45
1:A:24:ILE:HD11	1:A:55:ILE:HD12	1.99	0.45
1:B:41:ARG:HD3	4:B:1175:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:THR:O	1:B:91:GLU:HG3	2.17	0.45
1:B:24:ILE:HD11	1:B:55:ILE:HD12	1.98	0.45
1:B:68:ARG:NH2	1:B:96:TYR:OH	2.50	0.45
1:D:1:MET:CE	1:D:43:GLN:HG3	2.47	0.45
1:D:61:GLN:C	1:D:63:GLU:N	2.75	0.45
1:B:158:THR:O	1:B:162:GLU:HG2	2.17	0.44
1:C:116:ASN:HA	1:C:144:THR:HG22	1.99	0.44
1:A:42:LYS:O	1:A:52:LEU:HD12	2.18	0.44
1:D:40:TYR:CD1	1:D:40:TYR:N	2.85	0.44
1:D:66:ALA:O	4:D:913:HOH:O	2.21	0.44
1:B:35:THR:O	1:B:59:ALA:HB2	2.17	0.44
1:B:142:ILE:HD12	4:B:1208:HOH:O	2.16	0.44
1:B:1:MET:HE1	4:D:941:HOH:O	2.16	0.44
1:C:78:PHE:CD1	1:C:78:PHE:N	2.86	0.44
1:D:27:HIS:H	1:D:27:HIS:HD2	1.66	0.44
1:D:31:GLU:C	1:D:33:ASP:N	2.73	0.44
1:D:44:VAL:HG22	1:D:45:VAL:H	1.82	0.44
1:A:169:ARG:C	1:A:171:LEU:N	2.73	0.44
1:B:113:LEU:O	1:B:141:TYR:HA	2.17	0.44
1:D:109:VAL:HG13	4:D:1069:HOH:O	2.18	0.44
1:A:1:MET:HE1	1:A:43:GLN:OE1	2.17	0.44
1:A:16:LYS:O	1:A:19:LEU:HB2	2.17	0.44
1:A:41:ARG:HA	1:A:53:LEU:O	2.17	0.44
1:B:32:TYR:CD1	1:B:32:TYR:C	2.95	0.44
1:B:170:LYS:HA	1:B:170:LYS:HD3	1.44	0.44
1:C:141:TYR:CE2	1:C:143:GLU:HG3	2.52	0.44
1:A:22:GLN:HE21	1:A:22:GLN:CA	2.30	0.44
1:D:23:LEU:HD12	1:D:153:GLU:HG3	2.00	0.44
1:A:78:PHE:CG	1:A:100:ILE:HD13	2.52	0.43
1:A:3:GLU:HA	1:A:52:LEU:O	2.18	0.43
1:B:16:LYS:O	1:B:19:LEU:HB2	2.17	0.43
1:B:69:ASP:C	1:B:73:ARG:HH11	2.26	0.43
1:B:111:MET:HE3	1:B:111:MET:HB2	1.73	0.43
1:B:169:ARG:HG2	4:B:1226:HOH:O	2.18	0.43
1:A:5:LYS:HD3	1:A:54:ASP:HB3	2.01	0.43
1:B:90:PHE:CE2	1:B:125:VAL:HG13	2.54	0.43
1:B:141:TYR:CE2	1:B:143:GLU:CG	3.01	0.43
1:B:166:HIS:HA	1:B:170:LYS:HG2	2.01	0.43
1:D:63:GLU:CG	1:D:64:TYR:N	2.72	0.43
1:D:67:MET:HB3	1:D:70:GLN:HG3	2.00	0.43
1:A:106:SER:OG	1:A:108:ASP:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:LEU:HA	1:A:120:LEU:HD23	1.76	0.43
1:B:92:ASP:O	1:B:95:GLN:HB2	2.19	0.43
1:D:136:SER:HB2	4:D:1108:HOH:O	2.18	0.43
1:A:6:LEU:N	1:A:6:LEU:CD1	2.81	0.43
1:B:120:LEU:HD12	1:B:120:LEU:HA	1.78	0.43
1:C:171:LEU:HG	1:C:171:LEU:O	2.18	0.43
1:D:78:PHE:CD1	1:D:78:PHE:N	2.87	0.43
1:D:168:LEU:HD22	1:D:168:LEU:N	2.33	0.43
1:A:21:ILE:HD11	1:A:29:VAL:CG2	2.49	0.43
1:B:166:HIS:CD2	1:B:170:LYS:HZ2	2.37	0.43
1:A:134:ALA:O	1:A:138:GLY:N	2.52	0.43
1:B:23:LEU:HD22	1:B:156:PHE:CD2	2.54	0.43
1:B:72:MET:CE	1:B:103:VAL:CG1	2.96	0.43
1:B:97:ARG:HD3	1:B:137:TYR:CE1	2.54	0.43
1:B:137:TYR:HD1	1:B:137:TYR:HA	1.70	0.43
1:D:161:ARG:NH1	4:D:1068:HOH:O	2.52	0.43
1:A:41:ARG:HG2	1:B:135:ARG:HH21	1.84	0.42
1:A:165:GLN:NE2	4:A:828:HOH:O	2.51	0.42
1:B:24:ILE:O	1:B:24:ILE:HG22	2.19	0.42
1:D:133:LEU:HB2	4:D:1191:HOH:O	2.12	0.42
1:B:32:TYR:C	1:B:32:TYR:HD1	2.26	0.42
1:B:61:GLN:HG3	1:B:63:GLU:OE1	2.19	0.42
1:C:85:ASN:HD22	1:C:85:ASN:N	2.15	0.42
1:D:19:LEU:HD13	1:D:146:ALA:HB2	2.01	0.42
1:A:6:LEU:N	1:A:6:LEU:HD12	2.34	0.42
1:A:68:ARG:HD3	1:A:99:GLN:OE1	2.19	0.42
1:A:68:ARG:NE	1:A:96:TYR:HE1	2.14	0.42
1:B:67:MET:C	1:B:69:ASP:N	2.77	0.42
1:C:135:ARG:HD2	4:D:992:HOH:O	2.19	0.42
1:A:35:THR:CG2	1:A:36:ILE:H	2.31	0.42
1:B:27:HIS:HD2	4:B:1192:HOH:O	2.03	0.42
1:B:128:ARG:HG3	4:B:1209:HOH:O	2.18	0.42
1:D:67:MET:HA	4:D:1158:HOH:O	2.19	0.42
1:D:14:VAL:CG2	1:D:81:VAL:HG22	2.49	0.42
1:D:14:VAL:HG23	1:D:81:VAL:HG22	2.01	0.42
1:D:63:GLU:CG	1:D:64:TYR:CD1	3.00	0.42
1:B:22:GLN:HG3	1:B:146:ALA:O	2.19	0.42
1:C:97:ARG:HG2	4:C:1087:HOH:O	2.19	0.42
1:D:24:ILE:HG12	1:D:40:TYR:HB3	2.01	0.42
1:A:45:VAL:HG23	1:A:49:GLU:N	2.35	0.42
1:A:107:ASP:O	1:A:107:ASP:OD2	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:GLY:O	1:B:117:LYS:NZ	2.50	0.42
1:B:70:GLN:HG3	1:B:73:ARG:HB2	2.02	0.42
1:A:141:TYR:OH	1:A:143:GLU:OE2	2.35	0.42
1:D:68:ARG:HA	1:D:71:TYR:CE2	2.53	0.42
1:A:58:THR:CG2	1:A:59:ALA:N	2.78	0.42
1:A:14:VAL:HG23	1:A:15:GLY:N	2.35	0.41
1:B:70:GLN:CG	1:B:70:GLN:O	2.68	0.41
1:C:168:LEU:HD12	1:C:168:LEU:HA	1.60	0.41
1:D:17:SER:HB3	1:D:32:TYR:CD2	2.55	0.41
1:D:170:LYS:HG3	1:D:171:LEU:N	2.35	0.41
1:D:56:LEU:HD22	1:D:57:ASP:N	2.35	0.41
1:A:29:VAL:HG13	1:A:32:TYR:H	1.83	0.41
1:A:71:TYR:O	1:A:75:GLY:N	2.48	0.41
1:A:137:TYR:O	4:A:1037:HOH:O	2.22	0.41
1:C:85:ASN:C	1:C:85:ASN:ND2	2.77	0.41
1:A:163:ILE:O	1:A:164:ARG:C	2.64	0.41
1:B:97:ARG:O	1:B:101:LYS:HD3	2.21	0.41
1:A:123:ARG:NH2	4:A:1201:HOH:O	2.51	0.41
1:B:41:ARG:CB	1:D:135:ARG:CZ	2.99	0.41
1:B:90:PHE:C	1:B:90:PHE:CD1	2.99	0.41
1:B:117:LYS:O	1:B:120:LEU:HB2	2.21	0.41
1:D:14:VAL:HG11	4:D:1174:HOH:O	2.21	0.41
1:D:67:MET:O	1:D:71:TYR:CE1	2.74	0.41
1:D:67:MET:HB3	4:D:1158:HOH:O	2.19	0.41
1:C:113:LEU:HD23	1:C:113:LEU:HA	1.82	0.41
1:C:116:ASN:HA	1:C:144:THR:CG2	2.50	0.41
1:A:30:ASP:HB3	3:A:180:GCP:O2'	2.21	0.41
1:A:35:THR:HB	3:A:180:GCP:O1G	2.21	0.41
1:B:166:HIS:CD2	1:B:170:LYS:NZ	2.80	0.41
1:D:27:HIS:CD2	1:D:27:HIS:H	2.39	0.41
1:B:9:VAL:HG12	1:B:80:CYS:HA	2.02	0.41
1:C:8:VAL:HB	1:C:79:LEU:HB2	2.03	0.41
1:C:84:ILE:HD13	1:C:115:GLY:C	2.46	0.41
1:C:98:GLU:OE1	1:C:101:LYS:NZ	2.53	0.41
1:C:101:LYS:HD3	1:C:107:ASP:HA	2.03	0.41
1:D:52:LEU:HD12	1:D:52:LEU:HA	1.83	0.41
1:D:160:VAL:O	1:D:163:ILE:HB	2.20	0.41
1:A:169:ARG:C	1:A:171:LEU:H	2.15	0.41
1:D:97:ARG:O	1:D:97:ARG:CG	2.66	0.41
1:D:97:ARG:O	1:D:101:LYS:HG3	2.21	0.41
1:A:149:ARG:CB	1:A:152:VAL:HG11	2.46	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:HIS:HA	1:B:170:LYS:CB	2.49	0.40
1:D:136:SER:CA	4:D:1108:HOH:O	2.67	0.40
1:D:160:VAL:CA	4:D:1110:HOH:O	2.68	0.40
1:A:87:THR:HG22	1:A:91:GLU:OE2	2.21	0.40
1:B:82:PHE:CD1	1:B:82:PHE:C	2.98	0.40
1:D:34:PRO:HB3	3:D:180:GCP:O2G	2.21	0.40
1:D:76:GLU:O	1:D:110:PRO:HG2	2.21	0.40
1:B:127:SER:O	1:B:130:ALA:N	2.53	0.40
1:C:22:GLN:HG3	1:C:146:ALA:O	2.21	0.40
1:A:24:ILE:C	1:A:26:ASN:N	2.79	0.40
1:A:61:GLN:HB2	3:A:180:GCP:O3G	2.21	0.40
1:B:132:ASP:O	1:B:135:ARG:HB2	2.21	0.40
1:B:24:ILE:C	1:B:25:GLN:HG3	2.47	0.40
1:C:25:GLN:NE2	4:C:1142:HOH:O	2.25	0.40
1:C:68:ARG:HA	1:C:71:TYR:CE2	2.56	0.40
1:D:46:ILE:HG22	1:D:164:ARG:HH22	1.86	0.40
1:D:97:ARG:NH1	4:D:865:HOH:O	2.45	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	169/171 (99%)	135 (80%)	25 (15%)	9 (5%)	1	0
1	B	169/171 (99%)	150 (89%)	8 (5%)	11 (6%)	1	0
1	C	169/171 (99%)	159 (94%)	7 (4%)	3 (2%)	6	1
1	D	169/171 (99%)	149 (88%)	12 (7%)	8 (5%)	2	0
All	All	676/684 (99%)	593 (88%)	52 (8%)	31 (5%)	2	0

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	GLU
1	A	67	MET
1	A	69	ASP
1	A	73	ARG
1	B	34	PRO
1	B	62	GLU
1	B	64	TYR
1	B	69	ASP
1	B	122	ALA
1	C	63	GLU
1	C	169	ARG
1	D	65	SER
1	D	122	ALA
1	A	26	ASN
1	A	68	ARG
1	B	66	ALA
1	B	106	SER
1	B	118	CYS
1	D	63	GLU
1	A	65	SER
1	B	169	ARG
1	C	170	LYS
1	D	30	ASP
1	D	33	ASP
1	D	64	TYR
1	A	24	ILE
1	D	61	GLN
1	D	118	CYS
1	B	68	ARG
1	A	36	ILE
1	B	105	ASP

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	149/149 (100%)	118 (79%)	31 (21%)	1 0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	149/149 (100%)	121 (81%)	28 (19%)	1	0
1	C	149/149 (100%)	127 (85%)	22 (15%)	3	0
1	D	149/149 (100%)	124 (83%)	25 (17%)	2	0
All	All	596/596 (100%)	490 (82%)	106 (18%)	2	0

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	VAL
1	A	16	LYS
1	A	21	ILE
1	A	22	GLN
1	A	27	HIS
1	A	31	GLU
1	A	36	ILE
1	A	37	GLU
1	A	38	ASP
1	A	39	SER
1	A	58	THR
1	A	61	GLN
1	A	67	MET
1	A	68	ARG
1	A	84	ILE
1	A	88	LYS
1	A	97	ARG
1	A	99	GLN
1	A	107	ASP
1	A	109	VAL
1	A	112	VAL
1	A	113	LEU
1	A	114	VAL
1	A	118	CYS
1	A	128	ARG
1	A	136	SER
1	A	147	LYS
1	A	152	VAL
1	A	160	VAL
1	A	169	ARG
1	A	171	LEU
1	B	9	VAL
1	B	26	ASN

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Mol	Chain	Res	Type
1	B	29	VAL
1	B	32	TYR
1	B	36	ILE
1	B	37	GLU
1	B	50	THR
1	B	52	LEU
1	B	63	GLU
1	B	65	SER
1	B	67	MET
1	B	69	ASP
1	B	74	THR
1	B	79	LEU
1	B	81	VAL
1	B	84	ILE
1	B	93	ILE
1	B	101	LYS
1	B	109	VAL
1	B	120	LEU
1	B	123	ARG
1	B	128	ARG
1	B	129	GLN
1	B	135	ARG
1	B	136	SER
1	B	142	ILE
1	B	170	LYS
1	B	171	LEU
1	C	8	VAL
1	C	19	LEU
1	C	26	ASN
1	C	44	VAL
1	C	62	GLU
1	C	79	LEU
1	C	81	VAL
1	C	84	ILE
1	C	85	ASN
1	C	88	LYS
1	C	93	ILE
1	C	95	GLN
1	C	101	LYS
1	C	102	ARG
1	C	113	LEU
1	C	120	LEU

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Mol	Chain	Res	Type
1	C	124	THR
1	C	125	VAL
1	C	128	ARG
1	C	152	VAL
1	C	168	LEU
1	C	170	LYS
1	D	22	GLN
1	D	24	ILE
1	D	26	ASN
1	D	29	VAL
1	D	35	THR
1	D	36	ILE
1	D	39	SER
1	D	41	ARG
1	D	56	LEU
1	D	62	GLU
1	D	64	TYR
1	D	65	SER
1	D	71	TYR
1	D	81	VAL
1	D	84	ILE
1	D	88	LYS
1	D	101	LYS
1	D	109	VAL
1	D	113	LEU
1	D	124	THR
1	D	125	VAL
1	D	128	ARG
1	D	131	GLN
1	D	149	ARG
1	D	158	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	26	ASN
1	B	25	GLN
1	B	27	HIS
1	B	43	GLN
1	B	61	GLN
1	B	70	GLN

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Mol	Chain	Res	Type
1	B	116	ASN
1	B	131	GLN
1	B	166	HIS
1	C	61	GLN
1	C	70	GLN
1	C	85	ASN
1	C	94	HIS
1	C	150	GLN
1	C	165	GLN
1	D	22	GLN
1	D	26	ASN
1	D	70	GLN
1	D	95	GLN
1	D	99	GLN
1	D	131	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GCP	A	180	2	32,34,34	1.40	1 (3%)	49,54,54	1.36	7 (14%)
3	GCP	C	180	2	32,34,34	1.06	3 (9%)	49,54,54	1.59	5 (10%)
3	GCP	B	180	2	32,34,34	1.14	2 (6%)	49,54,54	1.39	5 (10%)
3	GCP	D	180	2	32,34,34	1.60	4 (12%)	49,54,54	1.02	3 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GCP	A	180	2	-	0/19/38/38	0/3/3/3
3	GCP	C	180	2	-	2/19/38/38	0/3/3/3
3	GCP	B	180	2	-	1/19/38/38	0/3/3/3
3	GCP	D	180	2	-	0/19/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	180	GCP	PG-O1G	6.33	1.63	1.50
3	D	180	GCP	PG-O3G	5.24	1.66	1.55
3	B	180	GCP	PG-O3G	5.13	1.66	1.55
3	D	180	GCP	PG-O1G	4.99	1.60	1.50
3	C	180	GCP	PG-O1G	2.78	1.55	1.50
3	D	180	GCP	PB-O2B	-2.75	1.49	1.56
3	C	180	GCP	C4-N3	2.18	1.39	1.34
3	B	180	GCP	PB-O3A	-2.16	1.56	1.58
3	C	180	GCP	PG-O3G	-2.16	1.50	1.55
3	D	180	GCP	PB-O3A	-2.13	1.56	1.58

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	180	GCP	O1B-PB-C3B	5.79	124.31	109.05
3	B	180	GCP	O3G-PG-O1G	-4.82	99.93	112.39
3	C	180	GCP	O2B-PB-C3B	-4.23	89.20	106.73
3	B	180	GCP	O1B-PB-C3B	4.23	120.20	109.05
3	C	180	GCP	O2B-PB-O1B	4.23	123.72	109.95
3	C	180	GCP	O2G-PG-O1G	-4.17	101.60	112.39
3	A	180	GCP	O1B-PB-C3B	-4.13	98.18	109.05
3	B	180	GCP	O3G-PG-C3B	3.84	115.70	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	180	GCP	O2B-PB-O1B	3.73	122.09	109.95
3	A	180	GCP	O2B-PB-C3B	3.66	121.89	106.73
3	D	180	GCP	O2B-PB-O1B	3.66	121.85	109.95
3	B	180	GCP	O2B-PB-C3B	-3.42	92.59	106.73
3	A	180	GCP	O3G-PG-C3B	2.79	113.16	106.40
3	D	180	GCP	O2G-PG-O1G	-2.58	105.72	112.39
3	B	180	GCP	O2B-PB-O1B	2.57	118.30	109.95
3	A	180	GCP	O3G-PG-O1G	-2.36	106.27	112.39
3	A	180	GCP	O2G-PG-C3B	2.26	111.89	106.40
3	C	180	GCP	O3G-PG-C3B	2.22	111.80	106.40
3	A	180	GCP	PB-O3A-PA	-2.11	125.48	132.37
3	D	180	GCP	O3G-PG-C3B	2.01	111.28	106.40

There are no chirality outliers.

All (3) torsion outliers are listed below:

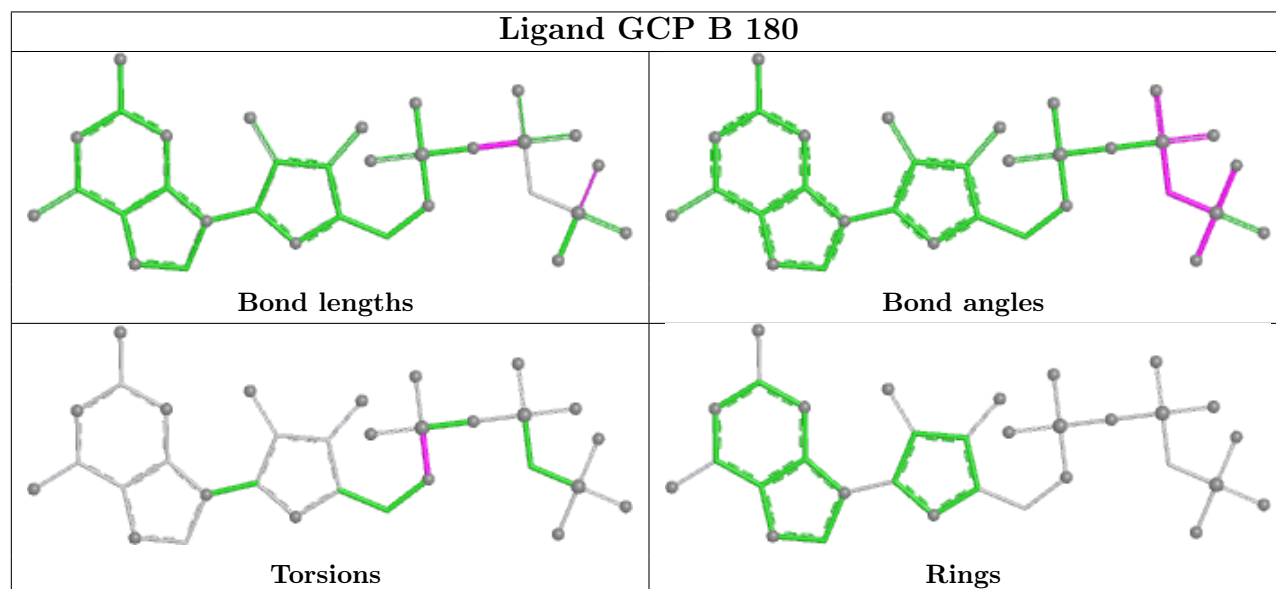
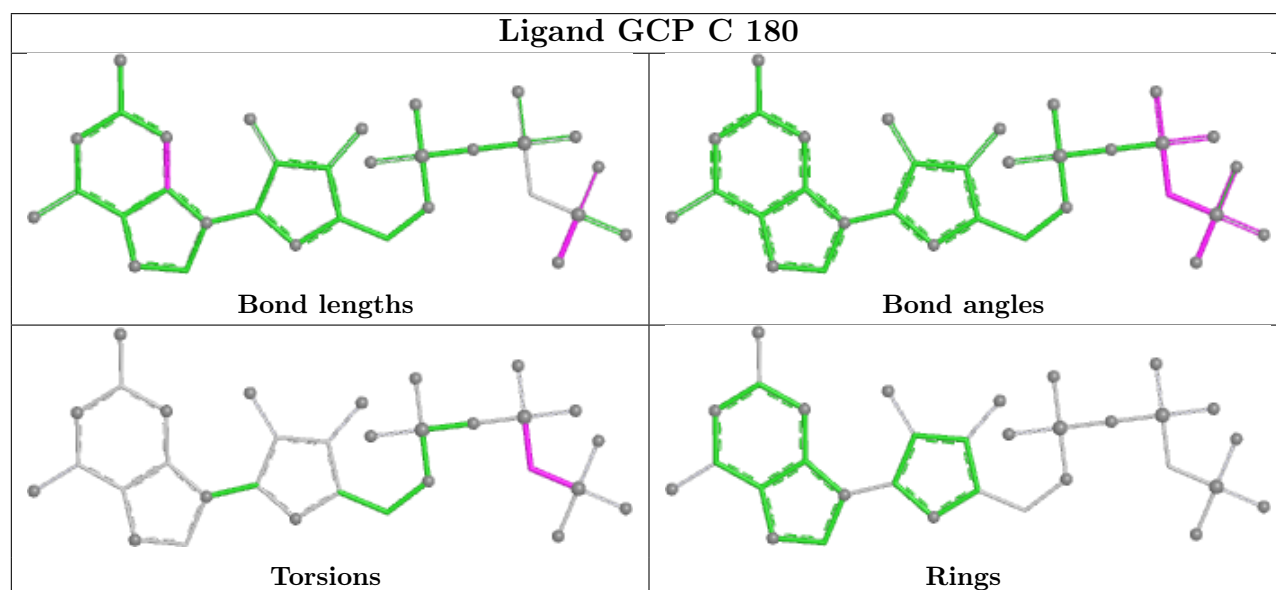
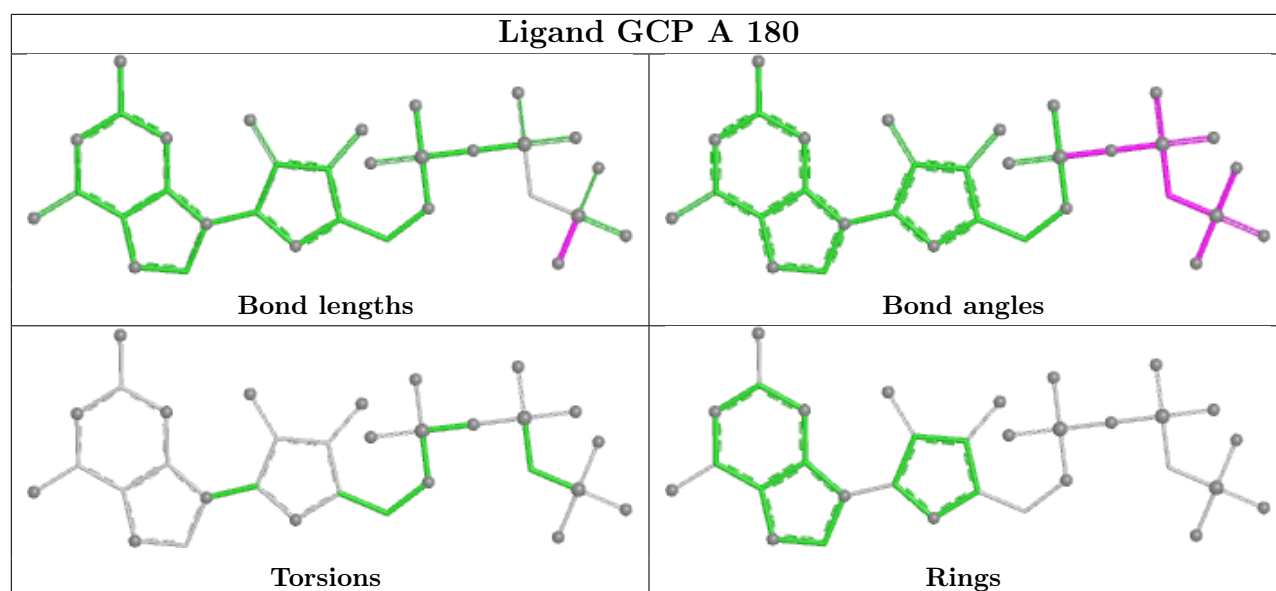
Mol	Chain	Res	Type	Atoms
3	B	180	GCP	C5'-O5'-PA-O1A
3	C	180	GCP	PG-C3B-PB-O2B
3	C	180	GCP	PB-C3B-PG-O1G

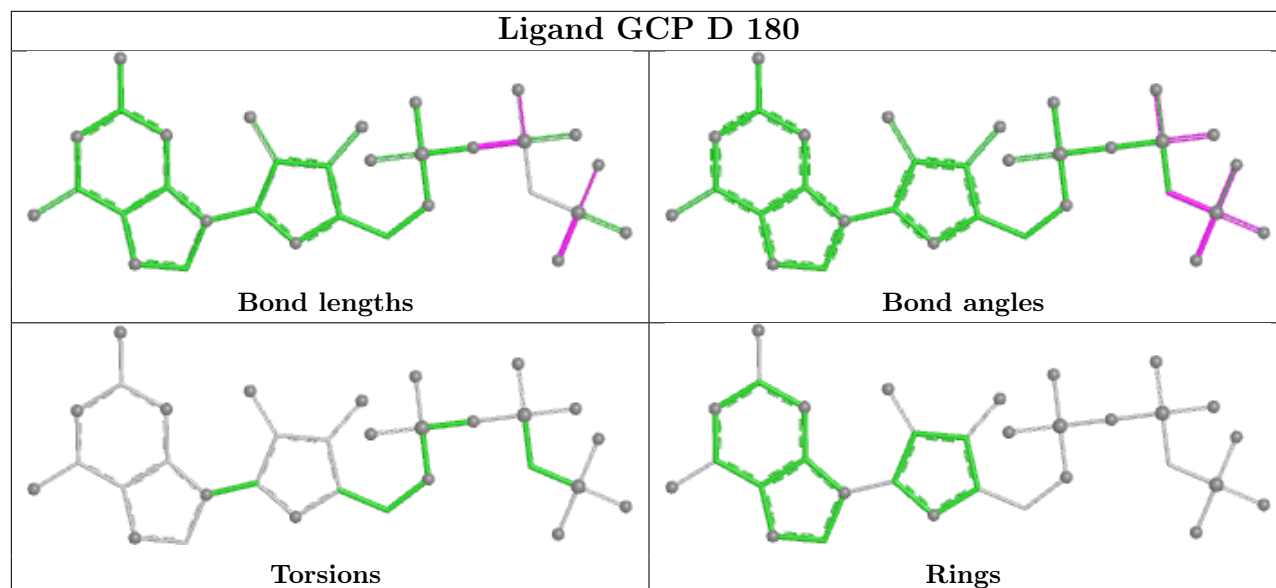
There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	180	GCP	6	0
3	C	180	GCP	4	0
3	B	180	GCP	2	0
3	D	180	GCP	5	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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.