



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:31 PM UTC

PDB ID : 2BMC / pdb_00002bmc
Title : Aurora-2 T287D T288D complexed with PHA-680632
Authors : Cameron, A.D.; Izzo, G.; Sagliano, A.; Rusconi, L.; Storici, P.; Fancelli, D.; Berta, D.; Bindi, S.; Catana, C.; Forte, B.; Giordano, P.; Mantegani, S.; Meroni, M.; Moll, J.; Pittala, V.; Severino, D.; Tonani, R.; Varasi, M.; Vulpetti, A.; Vianello, P.
Deposited on : 2005-03-11
Resolution : 2.60 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)
Xtrriage (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

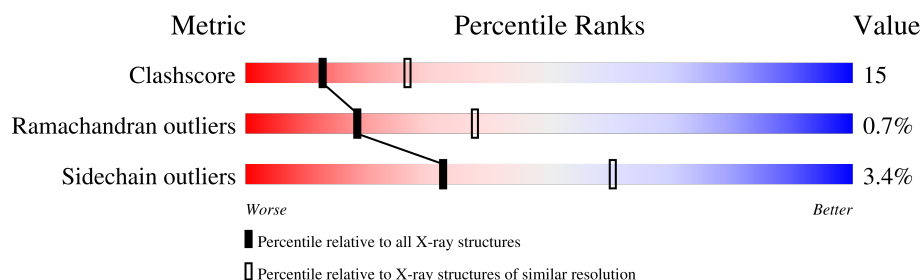
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)

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	306	
1	B	306	
1	C	306	
1	D	306	
1	E	306	
1	F	306	

2 Entry composition [i](#)

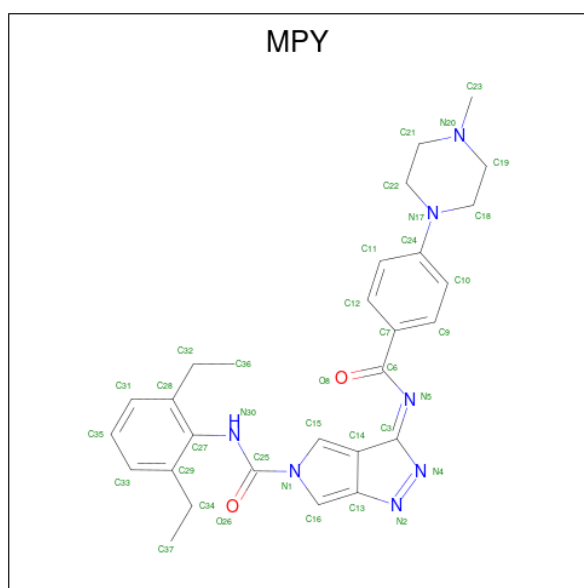
There are 3 unique types of molecules in this entry. The entry contains 13094 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERINE THREONINE-PROTEIN KINASE 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	0	0	0
			2105	1355	367	377	6			
1	B	258	Total	C	N	O	S	0	0	0
			2125	1367	373	379	6			
1	C	257	Total	C	N	O	S	0	0	0
			2116	1362	371	377	6			
1	D	256	Total	C	N	O	S	0	0	0
			2105	1355	367	377	6			
1	E	256	Total	C	N	O	S	0	0	0
			2105	1355	367	377	6			
1	F	257	Total	C	N	O	S	0	0	0
			2116	1361	371	378	6			

- Molecule 2 is (3E)-N-(2,6-DIETHYLPHENYL)-3-{[4-(4-METHYLPYPERAZIN-1-YL)BENZOYL]IMINO}PYRROLO[3,4-C]PYRAZOLE-5(3H)-CARBOXAMIDE (CCD ID: MPY) (formula: C₂₈H₃₁N₇O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			37	28	7	2		
2	B	1	Total	C	N	O	0	0
			37	28	7	2		
2	C	1	Total	C	N	O	0	0
			37	28	7	2		
2	D	1	Total	C	N	O	0	0
			37	28	7	2		
2	E	1	Total	C	N	O	0	0
			37	28	7	2		
2	F	1	Total	C	N	O	0	0
			37	28	7	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	45	Total	O	0	0
			45	45		
3	B	37	Total	O	0	0
			37	37		
3	C	26	Total	O	0	0
			26	26		
3	D	26	Total	O	0	0
			26	26		
3	E	29	Total	O	0	0
			29	29		
3	F	37	Total	O	0	0
			37	37		

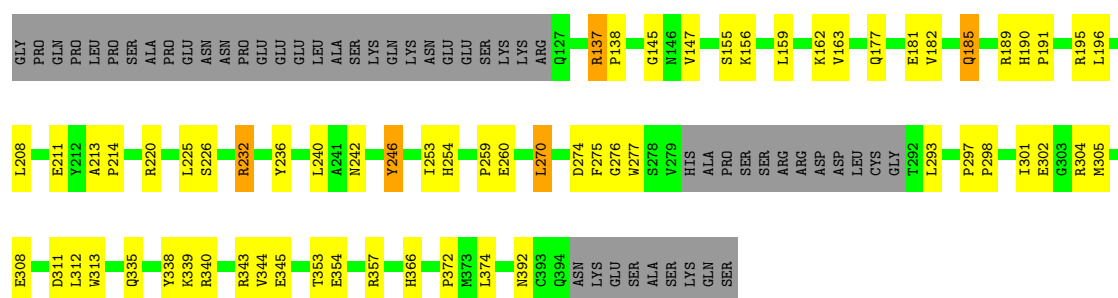
3 Residue-property plots

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. 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.

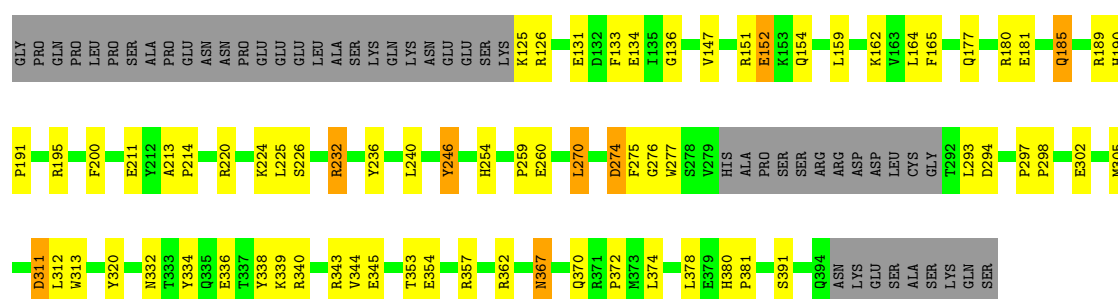
• Molecule 1: SERINE THREONINE-PROTEIN KINASE 6

Chain A: 



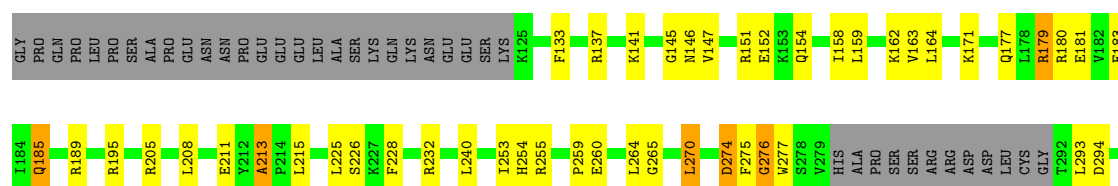
• Molecule 1: SERINE THREONINE-PROTEIN KINASE 6

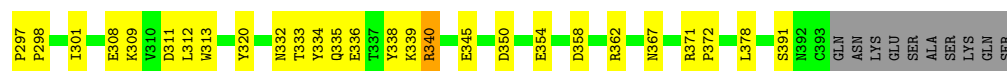
Chain B: 



• Molecule 1: SERINE THREONINE-PROTEIN KINASE 6

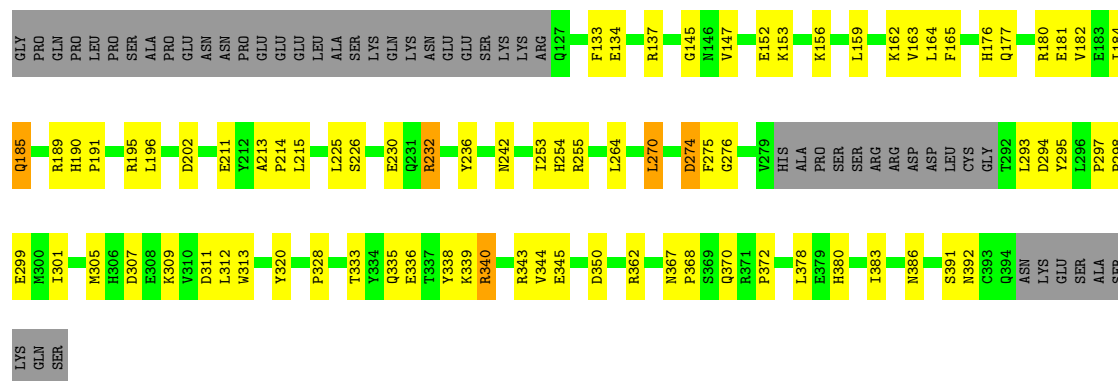
Chain C: 





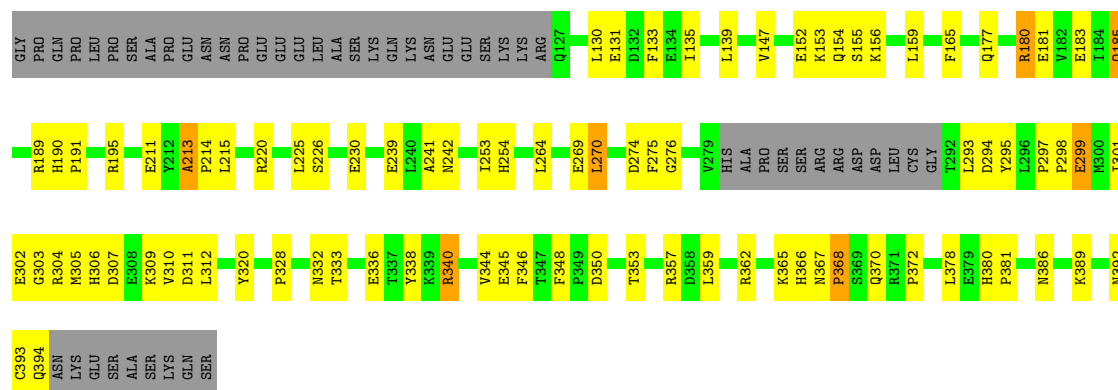
• Molecule 1: SERINE THREONINE-PROTEIN KINASE 6

Chain D: 58% 25% 16%



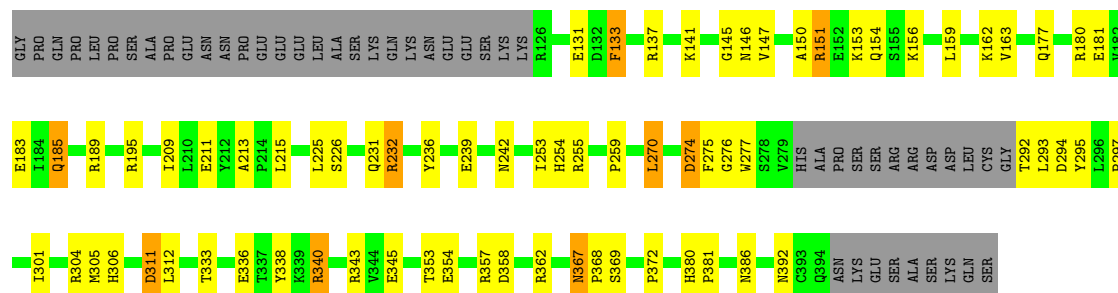
• Molecule 1: SERINE THREONINE-PROTEIN KINASE 6

Chain E: 55% 26% 16%



• Molecule 1: SERINE THREONINE-PROTEIN KINASE 6

Chain F: 60% 21% 16%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	74.02Å 101.22Å 101.49Å 115.73° 92.40° 101.54°	Depositor
Resolution (Å)	29.95 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (29.95-2.60)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNX 2000	Depositor
R, R_{free}	0.228 , 0.252	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13094	wwPDB-VP
Average B, all atoms (Å ²)	48.0	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: MPY

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	0.44	0/2157	0.93	5/2915 (0.2%)
1	B	0.41	0/2177	0.93	8/2940 (0.3%)
1	C	0.43	0/2168	0.96	10/2928 (0.3%)
1	D	0.40	0/2157	0.93	8/2915 (0.3%)
1	E	0.42	0/2157	0.95	5/2915 (0.2%)
1	F	0.41	0/2168	0.95	8/2929 (0.3%)
All	All	0.42	0/12984	0.94	44/17542 (0.3%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	213	ALA	CA-C-N	7.88	128.07	119.87
1	E	213	ALA	C-N-CA	7.88	128.07	119.87
1	A	274	ASP	N-CA-C	-7.55	99.93	110.35
1	B	274	ASP	N-CA-C	-6.99	100.71	110.35
1	D	274	ASP	N-CA-C	-6.74	101.05	110.35
1	B	294	ASP	N-CA-C	6.61	119.33	111.33
1	F	274	ASP	N-CA-C	-6.52	101.36	110.35
1	C	274	ASP	N-CA-C	-6.43	101.47	110.35
1	E	274	ASP	N-CA-C	-6.41	101.51	110.35
1	A	213	ALA	CA-C-N	6.24	125.94	119.82
1	A	213	ALA	C-N-CA	6.24	125.94	119.82
1	C	367	ASN	CA-C-N	6.14	127.51	119.84
1	C	367	ASN	C-N-CA	6.14	127.51	119.84
1	B	213	ALA	CA-C-N	6.08	126.19	119.87
1	B	213	ALA	C-N-CA	6.08	126.19	119.87
1	F	215	LEU	N-CA-C	5.80	120.05	112.92
1	E	135	ILE	N-CA-C	5.74	116.68	107.73
1	E	133	PHE	N-CA-C	5.67	118.80	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	213	ALA	CA-C-N	5.51	125.18	119.56
1	C	213	ALA	C-N-CA	5.51	125.18	119.56
1	B	214	PRO	N-CA-C	5.51	120.81	114.03
1	F	133	PHE	N-CA-C	5.50	118.51	109.72
1	F	367	ASN	CA-C-N	5.49	125.57	119.32
1	F	367	ASN	C-N-CA	5.49	125.57	119.32
1	C	350	ASP	N-CA-C	5.44	117.21	111.28
1	D	380	HIS	CA-C-N	5.43	126.63	119.84
1	D	380	HIS	C-N-CA	5.43	126.63	119.84
1	F	137	ARG	N-CA-C	5.39	117.34	109.84
1	A	214	PRO	N-CA-C	5.37	120.90	113.98
1	D	152	GLU	N-CA-C	-5.29	101.34	109.76
1	B	136	GLY	N-CA-C	5.29	119.22	112.13
1	C	133	PHE	N-CA-C	5.29	118.19	109.72
1	A	137	ARG	N-CA-C	5.23	117.30	110.08
1	D	134	GLU	N-CA-C	-5.23	99.25	108.20
1	B	152	GLU	N-CA-C	-5.16	102.01	109.59
1	C	205	ARG	N-CA-C	5.12	116.97	109.24
1	D	214	PRO	N-CA-C	5.09	120.29	114.03
1	D	133	PHE	N-CA-C	5.08	117.85	109.72
1	F	213	ALA	CA-C-N	5.08	124.79	119.82
1	F	213	ALA	C-N-CA	5.08	124.79	119.82
1	B	133	PHE	N-CA-C	5.05	117.80	109.72
1	C	137	ARG	CA-C-N	5.01	124.77	119.76
1	C	137	ARG	C-N-CA	5.01	124.77	119.76
1	D	137	ARG	N-CA-C	5.00	116.98	110.08

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	2105	0	2107	55	0
1	B	2125	0	2133	69	0
1	C	2116	0	2125	61	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2105	0	2107	74	0
1	E	2105	0	2107	99	0
1	F	2116	0	2120	71	0
2	A	37	0	31	2	0
2	B	37	0	31	2	0
2	C	37	0	31	3	0
2	D	37	0	31	2	0
2	E	37	0	31	1	0
2	F	37	0	31	1	0
3	A	45	0	0	0	0
3	B	37	0	0	0	0
3	C	26	0	0	0	0
3	D	26	0	0	0	0
3	E	29	0	0	2	0
3	F	37	0	0	0	0
All	All	13094	0	12885	392	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 15.

All (392) 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:F:151:ARG:HD3	1:F:156:LYS:HA	1.50	0.93
1:E:348:PHE:CE2	1:E:357:ARG:HA	2.15	0.81
1:B:367:ASN:HD22	1:B:370:GLN:HG3	1.46	0.80
1:C:253:ILE:HG21	1:C:308:GLU:HG2	1.63	0.80
1:E:333:THR:OG1	1:E:336:GLU:HG3	1.83	0.78
1:D:153:LYS:O	1:D:156:LYS:HE3	1.85	0.77
1:F:185:GLN:HG3	1:F:275:PHE:CD2	2.19	0.76
1:F:131:GLU:O	1:F:153:LYS:HE2	1.85	0.76
1:E:346:PHE:HE1	1:E:348:PHE:CE1	2.04	0.75
1:E:340:ARG:HH11	1:E:340:ARG:HB3	1.51	0.75
1:B:362:ARG:NH2	1:F:343:ARG:HB3	2.02	0.75
1:D:242:ASN:HD21	1:D:392:ASN:H	1.35	0.75
1:B:367:ASN:HD22	1:B:370:GLN:CG	2.00	0.74
1:F:242:ASN:HD21	1:F:392:ASN:H	1.36	0.74
1:E:348:PHE:HE2	1:E:357:ARG:HA	1.52	0.73
1:B:162:LYS:HE2	1:B:164:LEU:HD21	1.70	0.72
1:C:378:LEU:HD22	1:C:391:SER:HB2	1.72	0.71
1:D:242:ASN:ND2	1:D:392:ASN:H	1.90	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:MET:HE2	1:B:338:TYR:HD2	1.57	0.69
1:D:254:HIS:HE1	1:D:274:ASP:OD1	1.76	0.69
1:F:254:HIS:HD2	1:F:255:ARG:O	1.77	0.68
1:C:254:HIS:HD2	1:C:255:ARG:O	1.78	0.67
1:E:348:PHE:CD2	1:E:357:ARG:HG2	2.30	0.67
1:E:253:ILE:HD13	1:E:307:ASP:HB2	1.76	0.66
1:E:298:PRO:O	1:E:302:GLU:HG3	1.95	0.66
1:E:297:PRO:HB2	1:E:299:GLU:OE2	1.96	0.66
1:F:232:ARG:HH11	1:F:232:ARG:HB2	1.61	0.66
1:F:367:ASN:HD22	1:F:368:PRO:HD2	1.60	0.66
1:E:305:MET:HG3	1:E:306:HIS:H	1.60	0.66
1:E:242:ASN:ND2	1:E:392:ASN:H	1.94	0.65
1:B:232:ARG:HB2	1:B:232:ARG:HH11	1.61	0.65
1:E:353:THR:O	1:E:357:ARG:HG3	1.97	0.65
1:A:339:LYS:HD3	1:A:343:ARG:HH21	1.62	0.64
1:E:367:ASN:HB3	1:E:370:GLN:HG2	1.78	0.64
1:E:338:TYR:CE2	1:F:305:MET:HG3	2.33	0.64
1:D:309:LYS:HA	1:D:312:LEU:HD12	1.80	0.64
1:F:353:THR:O	1:F:357:ARG:HG3	1.98	0.64
1:F:141:LYS:HD2	1:F:146:ASN:OD1	1.98	0.63
1:D:232:ARG:HH11	1:D:232:ARG:HB2	1.62	0.63
1:B:254:HIS:HE1	1:B:274:ASP:OD1	1.82	0.63
1:A:185:GLN:HG3	1:A:275:PHE:CD2	2.34	0.63
1:F:147:VAL:HG22	1:F:162:LYS:HG3	1.81	0.62
1:B:232:ARG:HD2	1:B:236:TYR:CZ	2.34	0.62
1:E:367:ASN:HB3	1:E:370:GLN:CG	2.30	0.62
1:E:346:PHE:HE1	1:E:348:PHE:HE1	1.48	0.62
1:B:370:GLN:NE2	1:E:370:GLN:NE2	2.48	0.61
1:C:358:ASP:O	1:C:362:ARG:HG2	2.01	0.61
1:F:231:GLN:NE2	1:F:386:ASN:HD22	1.98	0.61
1:B:370:GLN:NE2	1:E:370:GLN:HE22	1.99	0.61
1:B:340:ARG:NH1	1:B:345:GLU:O	2.34	0.60
1:B:362:ARG:HG3	1:B:362:ARG:HH11	1.65	0.60
1:C:254:HIS:HE1	1:C:274:ASP:OD1	1.84	0.60
1:A:270:LEU:HD23	1:A:270:LEU:C	2.27	0.60
1:D:253:ILE:HD13	1:D:307:ASP:HB2	1.81	0.60
1:A:338:TYR:CD2	1:B:305:MET:HE2	2.37	0.60
1:B:339:LYS:HD3	1:B:343:ARG:HH21	1.67	0.60
1:A:246:TYR:CD1	1:A:246:TYR:C	2.79	0.60
1:B:225:LEU:O	1:B:226:SER:HB2	2.02	0.60
1:D:335:GLN:H	1:D:335:GLN:CD	2.09	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:253:ILE:HG22	1:F:254:HIS:N	2.17	0.60
1:C:151:ARG:NH1	1:C:158:ILE:HD11	2.17	0.60
1:E:242:ASN:HD21	1:E:392:ASN:H	1.46	0.60
1:B:152:GLU:OE1	1:B:154:GLN:HB2	2.01	0.59
1:E:253:ILE:HG21	1:E:307:ASP:HB3	1.83	0.59
1:E:153:LYS:O	1:E:156:LYS:HE3	2.01	0.59
1:F:367:ASN:HD22	1:F:368:PRO:CD	2.14	0.59
1:E:340:ARG:HB3	1:E:340:ARG:NH1	2.15	0.59
1:A:242:ASN:HD21	1:A:392:ASN:H	1.50	0.59
1:E:215:LEU:HB2	1:E:264:LEU:HB2	1.85	0.59
1:E:309:LYS:HD3	1:E:312:LEU:HD12	1.83	0.59
1:F:312:LEU:HD13	1:F:372:PRO:O	2.01	0.59
1:D:215:LEU:HB2	1:D:264:LEU:HB2	1.84	0.58
1:B:334:TYR:O	1:B:338:TYR:HD1	1.87	0.58
1:A:293:LEU:HD11	1:B:338:TYR:CE1	2.38	0.58
1:F:151:ARG:CB	1:F:151:ARG:HH11	2.17	0.58
1:A:340:ARG:NH1	1:A:345:GLU:O	2.37	0.58
1:E:305:MET:HB3	1:F:338:TYR:CZ	2.38	0.58
1:D:182:VAL:HG13	1:D:196:LEU:HD21	1.86	0.58
1:A:338:TYR:CE1	1:B:293:LEU:HD11	2.39	0.57
1:F:304:ARG:O	1:F:305:MET:HE2	2.04	0.57
1:A:177:GLN:HG2	1:A:181:GLU:OE2	2.04	0.57
1:B:190:HIS:CG	1:B:191:PRO:HD2	2.40	0.57
1:A:343:ARG:HB3	1:E:362:ARG:NH2	2.19	0.57
1:A:312:LEU:HD13	1:A:372:PRO:O	2.05	0.57
1:C:338:TYR:CZ	1:D:305:MET:HB3	2.40	0.57
1:B:353:THR:O	1:B:357:ARG:HG3	2.05	0.57
1:C:301:ILE:HD11	1:D:338:TYR:HA	1.87	0.57
1:D:344:VAL:HG12	1:D:344:VAL:O	2.04	0.57
1:E:346:PHE:CE1	1:E:348:PHE:HE1	2.22	0.57
1:F:293:LEU:HD22	1:F:301:ILE:HG21	1.87	0.56
1:E:309:LYS:HA	1:E:312:LEU:HD12	1.87	0.56
1:F:151:ARG:HH11	1:F:151:ARG:HB3	1.70	0.56
1:A:242:ASN:ND2	1:A:392:ASN:H	2.03	0.56
1:C:255:ARG:CZ	1:C:276:GLY:HA2	2.36	0.56
1:C:312:LEU:HD13	1:C:372:PRO:O	2.06	0.56
1:D:333:THR:HG23	1:D:336:GLU:OE2	2.05	0.56
1:D:190:HIS:CG	1:D:191:PRO:HD2	2.41	0.56
1:D:270:LEU:HD23	1:D:270:LEU:O	2.06	0.56
1:E:190:HIS:CG	1:E:191:PRO:HD2	2.41	0.56
1:E:362:ARG:HG3	1:E:362:ARG:HH11	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:362:ARG:HG3	1:D:362:ARG:HH11	1.71	0.55
1:C:185:GLN:HG3	1:C:277:TRP:CD1	2.41	0.55
1:A:232:ARG:HD2	1:A:236:TYR:CZ	2.42	0.55
1:A:338:TYR:CZ	1:B:293:LEU:HD11	2.41	0.55
1:D:230:GLU:HB3	1:D:386:ASN:ND2	2.22	0.55
1:E:312:LEU:HD13	1:E:372:PRO:O	2.06	0.55
1:F:232:ARG:HD2	1:F:236:TYR:CZ	2.42	0.54
1:D:253:ILE:HG22	1:D:254:HIS:N	2.22	0.54
1:E:346:PHE:CE1	1:E:348:PHE:CE1	2.90	0.54
1:F:293:LEU:O	1:F:295:TYR:N	2.40	0.54
1:B:131:GLU:N	1:B:131:GLU:OE1	2.36	0.54
1:D:270:LEU:HD23	1:D:270:LEU:C	2.33	0.54
1:E:241:ALA:HB3	1:E:378:LEU:HD21	1.89	0.54
1:E:298:PRO:HA	1:E:301:ILE:HG22	1.90	0.54
1:A:225:LEU:O	1:A:226:SER:HB2	2.06	0.54
1:D:340:ARG:NH1	1:D:345:GLU:O	2.41	0.54
1:D:242:ASN:HD21	1:D:392:ASN:N	2.05	0.54
1:D:367:ASN:C	1:D:367:ASN:HD22	2.15	0.54
1:F:195:ARG:HB3	1:F:211:GLU:HB2	1.89	0.54
1:E:155:SER:O	1:E:156:LYS:HB2	2.07	0.54
1:E:338:TYR:CD2	1:F:305:MET:HG3	2.43	0.54
1:A:190:HIS:CG	1:A:191:PRO:HD2	2.43	0.54
1:E:367:ASN:O	1:E:370:GLN:HB2	2.08	0.54
1:F:358:ASP:O	1:F:362:ARG:HG2	2.08	0.54
1:E:393:CYS:O	1:E:394:GLN:HB2	2.07	0.53
1:F:311:ASP:OD1	1:F:311:ASP:N	2.36	0.53
1:C:177:GLN:HG2	1:C:181:GLU:OE2	2.08	0.53
1:D:230:GLU:HB3	1:D:386:ASN:HD22	1.73	0.53
1:B:339:LYS:O	1:B:343:ARG:HG3	2.09	0.53
1:C:228:PHE:CG	1:C:232:ARG:HD2	2.43	0.53
1:C:335:GLN:HG3	1:C:339:LYS:HE3	1.90	0.53
1:F:141:LYS:HD2	1:F:146:ASN:CG	2.33	0.53
1:A:181:GLU:OE1	1:A:275:PHE:O	2.27	0.53
1:A:312:LEU:HD21	1:A:374:LEU:HD23	1.91	0.53
1:B:298:PRO:O	1:B:302:GLU:HG3	2.08	0.53
1:B:378:LEU:HD22	1:B:391:SER:HB2	1.90	0.53
1:C:225:LEU:O	1:C:226:SER:HB2	2.07	0.53
1:A:298:PRO:HA	1:A:301:ILE:HG22	1.90	0.52
1:C:270:LEU:HD23	1:C:270:LEU:O	2.08	0.52
1:C:159:LEU:HD12	1:C:159:LEU:C	2.34	0.52
1:B:367:ASN:ND2	1:B:370:GLN:HG3	2.20	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:VAL:HG23	2:A:1395:MPY:H341	1.90	0.52
1:A:338:TYR:HD2	1:B:305:MET:HE2	1.74	0.52
1:C:180:ARG:O	1:C:183:GLU:HB3	2.10	0.52
1:F:177:GLN:HG2	1:F:181:GLU:OE2	2.09	0.52
1:F:380:HIS:ND1	1:F:381:PRO:HD2	2.24	0.52
1:C:147:VAL:HG23	2:C:1394:MPY:H341	1.91	0.52
1:E:366:HIS:O	1:E:368:PRO:HD3	2.09	0.52
1:E:293:LEU:HD22	1:E:301:ILE:CD1	2.39	0.52
1:A:344:VAL:HG21	1:A:366:HIS:HB2	1.92	0.51
1:E:185:GLN:HG3	1:E:275:PHE:CD2	2.45	0.51
1:C:152:GLU:OE1	1:C:154:GLN:HB2	2.11	0.51
1:F:362:ARG:HG3	1:F:362:ARG:HH11	1.76	0.51
1:D:340:ARG:HB3	1:D:340:ARG:HH11	1.75	0.51
1:A:232:ARG:HH11	1:A:232:ARG:HB2	1.76	0.50
1:A:301:ILE:O	1:A:304:ARG:HG2	2.11	0.50
1:C:179:ARG:HD3	1:C:179:ARG:C	2.35	0.50
1:C:293:LEU:HD13	1:C:301:ILE:HD11	1.93	0.50
1:E:332:ASN:HB2	1:E:336:GLU:OE1	2.11	0.50
1:D:367:ASN:HD22	1:D:368:PRO:N	2.09	0.50
1:E:305:MET:HG3	1:E:306:HIS:N	2.25	0.50
1:E:253:ILE:HG22	1:E:254:HIS:N	2.26	0.50
1:E:344:VAL:HG12	1:E:344:VAL:O	2.10	0.50
1:D:232:ARG:HD2	1:D:236:TYR:CZ	2.47	0.50
1:E:270:LEU:C	1:E:270:LEU:HD23	2.37	0.50
1:A:340:ARG:HH11	1:A:340:ARG:HB3	1.76	0.50
1:B:185:GLN:HG3	1:B:275:PHE:CD2	2.47	0.50
1:B:362:ARG:HG3	1:B:362:ARG:NH1	2.26	0.50
1:D:367:ASN:C	1:D:367:ASN:ND2	2.68	0.50
1:B:232:ARG:HH11	1:B:232:ARG:CB	2.24	0.50
1:B:195:ARG:HB3	1:B:211:GLU:HB2	1.94	0.50
1:F:180:ARG:O	1:F:183:GLU:HB3	2.12	0.50
1:A:335:GLN:CD	1:A:335:GLN:H	2.19	0.50
1:F:225:LEU:O	1:F:226:SER:HB2	2.11	0.50
1:A:185:GLN:HG3	1:A:275:PHE:HD2	1.77	0.49
1:A:353:THR:O	1:A:357:ARG:HG3	2.12	0.49
1:C:181:GLU:CD	1:C:255:ARG:HH22	2.20	0.49
1:C:301:ILE:HG12	1:D:338:TYR:HD1	1.77	0.49
1:F:274:ASP:C	1:F:274:ASP:OD1	2.56	0.49
1:C:334:TYR:CE1	1:D:293:LEU:HD11	2.47	0.49
1:A:301:ILE:HG23	1:A:302:GLU:N	2.27	0.49
1:A:392:ASN:C	1:A:392:ASN:HD22	2.21	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:LEU:HD13	1:B:372:PRO:O	2.13	0.49
1:C:298:PRO:O	1:C:301:ILE:HG22	2.12	0.49
1:A:253:ILE:HG22	1:A:254:HIS:N	2.28	0.48
1:A:302:GLU:HA	1:A:305:MET:HG2	1.95	0.48
1:B:340:ARG:HD2	1:B:345:GLU:HG3	1.96	0.48
1:C:313:TRP:CG	1:D:297:PRO:HB3	2.47	0.48
1:A:195:ARG:HB3	1:A:211:GLU:HB2	1.96	0.48
1:C:297:PRO:HB3	1:D:313:TRP:CG	2.48	0.48
1:C:232:ARG:HG3	1:C:232:ARG:HH11	1.78	0.48
1:E:367:ASN:C	1:E:367:ASN:ND2	2.70	0.48
1:A:253:ILE:CD1	1:A:308:GLU:HG2	2.43	0.48
1:C:254:HIS:O	1:C:311:ASP:OD1	2.32	0.48
1:E:295:TYR:CE2	1:F:259:PRO:HD3	2.49	0.48
1:F:254:HIS:CD2	1:F:255:ARG:N	2.81	0.48
1:F:255:ARG:CZ	1:F:276:GLY:HA2	2.43	0.48
1:B:177:GLN:HG2	1:B:181:GLU:OE2	2.14	0.48
1:C:270:LEU:HD23	1:C:270:LEU:C	2.38	0.48
1:E:309:LYS:HD3	1:E:312:LEU:CD1	2.43	0.48
1:A:340:ARG:CZ	1:E:381:PRO:HG3	2.44	0.48
1:D:312:LEU:HD13	1:D:372:PRO:O	2.14	0.48
1:B:246:TYR:CD1	1:B:246:TYR:C	2.91	0.47
1:C:195:ARG:HB3	1:C:211:GLU:HB2	1.95	0.47
1:A:392:ASN:C	1:A:392:ASN:ND2	2.71	0.47
1:B:224:LYS:NZ	2:B:1395:MPY:H233	2.29	0.47
1:A:182:VAL:HG13	1:A:196:LEU:HD21	1.96	0.47
1:D:225:LEU:O	1:D:226:SER:HB2	2.14	0.47
1:D:253:ILE:HG21	1:D:307:ASP:HB3	1.96	0.47
1:C:141:LYS:HD2	1:C:146:ASN:CG	2.39	0.47
1:E:177:GLN:HG2	1:E:181:GLU:OE2	2.14	0.47
1:E:340:ARG:NH1	1:E:345:GLU:O	2.48	0.47
1:E:131:GLU:O	1:E:153:LYS:HD2	2.14	0.47
1:B:185:GLN:HG2	1:B:277:TRP:CD1	2.50	0.47
1:C:332:ASN:HB2	1:C:336:GLU:OE2	2.15	0.47
1:E:195:ARG:HG2	1:E:195:ARG:HH11	1.79	0.47
1:E:225:LEU:O	1:E:226:SER:HB2	2.15	0.47
1:F:340:ARG:NH1	1:F:345:GLU:O	2.48	0.47
1:A:159:LEU:HD12	1:A:159:LEU:C	2.40	0.47
1:D:254:HIS:CE1	1:D:274:ASP:OD1	2.62	0.47
1:B:381:PRO:HG3	1:F:340:ARG:CZ	2.45	0.47
1:E:195:ARG:HG2	1:E:195:ARG:NH1	2.29	0.46
1:E:180:ARG:O	1:E:183:GLU:HB3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:310:VAL:HG13	1:F:297:PRO:HG3	1.97	0.46
1:B:159:LEU:C	1:B:159:LEU:HD12	2.40	0.46
1:E:239:GLU:HB3	1:E:270:LEU:HD22	1.96	0.46
1:B:181:GLU:OE1	1:B:275:PHE:O	2.34	0.46
1:F:159:LEU:HD12	1:F:159:LEU:C	2.40	0.46
1:C:338:TYR:CD1	1:C:338:TYR:C	2.92	0.46
1:D:165:PHE:CD1	1:D:165:PHE:N	2.83	0.46
1:E:293:LEU:HD22	1:E:301:ILE:HD12	1.96	0.46
1:D:145:GLY:HA3	1:D:163:VAL:O	2.15	0.46
1:F:147:VAL:HG23	2:F:1395:MPY:H341	1.97	0.46
1:C:309:LYS:NZ	1:C:371:ARG:O	2.43	0.46
1:C:333:THR:HG23	1:C:336:GLU:OE2	2.16	0.46
1:B:240:LEU:HG	1:B:270:LEU:HD11	1.97	0.46
1:C:141:LYS:HD2	1:C:146:ASN:OD1	2.15	0.46
1:F:185:GLN:HG3	1:F:275:PHE:HD2	1.79	0.46
1:B:147:VAL:HG23	2:B:1395:MPY:H341	1.97	0.46
1:B:370:GLN:CD	1:E:370:GLN:NE2	2.74	0.46
1:C:254:HIS:CE1	1:C:274:ASP:OD1	2.67	0.45
1:D:336:GLU:O	1:D:340:ARG:HG2	2.15	0.45
1:F:367:ASN:HD22	1:F:368:PRO:N	2.14	0.45
1:A:298:PRO:O	1:A:301:ILE:HG22	2.16	0.45
1:E:320:TYR:CD2	1:E:328:PRO:HD3	2.51	0.45
1:A:147:VAL:CG2	2:A:1395:MPY:H341	2.44	0.45
1:C:145:GLY:HA3	1:C:163:VAL:O	2.17	0.45
1:D:162:LYS:HE2	1:D:164:LEU:HD21	1.99	0.45
1:E:307:ASP:C	1:E:309:LYS:H	2.25	0.45
1:E:139:LEU:HD23	1:E:139:LEU:HA	1.80	0.45
1:D:181:GLU:OE1	1:D:275:PHE:O	2.35	0.45
1:B:344:VAL:HG12	1:B:344:VAL:O	2.17	0.45
1:C:259:PRO:HD3	1:D:295:TYR:CE2	2.52	0.45
1:F:181:GLU:OE1	1:F:275:PHE:O	2.34	0.45
1:B:332:ASN:HB2	1:B:336:GLU:OE1	2.17	0.45
1:C:253:ILE:CG2	1:C:308:GLU:HG2	2.39	0.45
1:B:339:LYS:HD3	1:B:343:ARG:NH2	2.32	0.45
1:D:230:GLU:CB	1:D:386:ASN:HD22	2.30	0.45
1:D:378:LEU:HD22	1:D:391:SER:HB2	1.97	0.45
1:E:338:TYR:HE1	1:F:301:ILE:O	2.00	0.45
1:A:259:PRO:HD2	1:A:260:GLU:OE1	2.17	0.45
1:A:313:TRP:CG	1:B:297:PRO:HB3	2.52	0.45
1:B:370:GLN:CD	1:E:370:GLN:HE22	2.25	0.45
1:A:301:ILE:CG2	1:A:302:GLU:N	2.80	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:311:ASP:OD1	1:B:311:ASP:N	2.50	0.44
1:E:195:ARG:HB3	1:E:211:GLU:HB2	1.99	0.44
1:A:240:LEU:HD23	1:A:270:LEU:HD21	1.98	0.44
1:B:302:GLU:HA	1:B:305:MET:HG2	1.99	0.44
1:C:309:LYS:HD3	1:C:312:LEU:HD12	2.00	0.44
1:C:320:TYR:CD1	1:C:320:TYR:C	2.94	0.44
1:D:362:ARG:HG3	1:D:362:ARG:NH1	2.33	0.44
1:E:185:GLN:HG3	1:E:275:PHE:HD2	1.80	0.44
1:F:239:GLU:OE1	1:F:270:LEU:HD22	2.18	0.44
1:C:147:VAL:CG2	2:C:1394:MPY:H341	2.46	0.44
1:F:145:GLY:HA3	1:F:163:VAL:O	2.17	0.44
1:D:159:LEU:C	1:D:159:LEU:HD12	2.42	0.44
1:E:367:ASN:C	1:E:367:ASN:HD22	2.25	0.44
1:C:254:HIS:CD2	1:C:255:ARG:N	2.86	0.44
1:D:333:THR:OG1	1:D:336:GLU:HG3	2.18	0.44
1:E:230:GLU:CB	1:E:386:ASN:HD22	2.29	0.44
1:E:336:GLU:O	1:E:340:ARG:HG2	2.17	0.44
1:B:270:LEU:C	1:B:270:LEU:HD23	2.42	0.44
1:D:185:GLN:HG3	1:D:275:PHE:CD2	2.52	0.44
1:D:185:GLN:HE21	1:D:185:GLN:HB3	1.59	0.44
1:F:150:ALA:HB3	1:F:159:LEU:HD12	1.98	0.44
1:E:389:LYS:HE2	3:E:2007:HOH:O	2.17	0.44
1:C:340:ARG:NH1	1:C:345:GLU:O	2.51	0.44
1:E:165:PHE:CD1	1:E:165:PHE:N	2.86	0.44
1:E:362:ARG:HG3	1:E:362:ARG:NH1	2.32	0.44
1:E:367:ASN:CB	1:E:370:GLN:HG2	2.48	0.44
1:E:380:HIS:HA	1:E:381:PRO:HD3	1.91	0.44
1:A:155:SER:O	1:A:156:LYS:C	2.61	0.43
1:A:297:PRO:HB3	1:B:313:TRP:CG	2.53	0.43
1:E:253:ILE:CG2	1:E:254:HIS:N	2.81	0.43
1:B:270:LEU:HD23	1:B:270:LEU:O	2.18	0.43
1:D:147:VAL:HG23	2:D:1395:MPY:H341	2.00	0.43
1:E:254:HIS:O	1:E:311:ASP:OD1	2.36	0.43
1:F:133:PHE:HZ	1:F:209:ILE:HD13	1.83	0.43
1:C:215:LEU:HD12	1:C:265:GLY:O	2.18	0.43
1:D:177:GLN:HG2	1:D:181:GLU:OE2	2.19	0.43
1:D:339:LYS:HD3	1:D:343:ARG:HH21	1.83	0.43
1:E:301:ILE:HG12	1:F:338:TYR:HD2	1.84	0.43
1:C:274:ASP:OD1	1:C:274:ASP:C	2.61	0.43
1:D:367:ASN:HD22	1:D:368:PRO:CD	2.32	0.43
1:D:367:ASN:O	1:D:370:GLN:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:213:ALA:HA	1:E:214:PRO:HD2	1.87	0.43
1:E:359:LEU:HD13	1:E:380:HIS:CD2	2.54	0.43
1:A:145:GLY:HA3	1:A:163:VAL:O	2.18	0.43
1:C:151:ARG:HH12	1:C:158:ILE:HD11	1.83	0.43
1:D:213:ALA:O	2:D:1395:MPY:H12	2.18	0.43
1:D:242:ASN:HD22	1:D:242:ASN:HA	1.64	0.43
1:E:180:ARG:NH1	1:E:180:ARG:HG2	2.32	0.43
1:F:333:THR:HG23	1:F:336:GLU:OE2	2.18	0.43
1:A:220:ARG:HA	1:A:220:ARG:HD3	1.85	0.43
1:B:125:LYS:HD2	1:B:126:ARG:H	1.84	0.43
1:C:301:ILE:HG12	1:D:338:TYR:CD1	2.54	0.43
1:E:181:GLU:OE1	1:E:275:PHE:O	2.37	0.42
1:B:134:GLU:HB2	1:B:151:ARG:HB3	2.01	0.42
1:D:298:PRO:HA	1:D:301:ILE:HG12	2.01	0.42
1:E:147:VAL:HG23	2:E:1395:MPY:H341	2.01	0.42
1:F:292:THR:HB	1:F:293:LEU:H	1.72	0.42
1:F:306:HIS:CE1	1:F:369:SER:HB3	2.54	0.42
1:F:367:ASN:C	1:F:367:ASN:ND2	2.77	0.42
1:D:215:LEU:HB2	1:D:264:LEU:CB	2.48	0.42
1:E:301:ILE:C	1:E:303:GLY:H	2.27	0.42
1:E:365:LYS:HD3	1:E:370:GLN:HB3	2.01	0.42
1:B:380:HIS:HA	1:B:381:PRO:HD3	1.92	0.42
1:C:171:LYS:O	1:D:176:HIS:CD2	2.73	0.42
1:D:195:ARG:HH11	1:D:195:ARG:HG2	1.84	0.42
1:B:232:ARG:HH11	1:B:232:ARG:CG	2.32	0.42
1:C:162:LYS:HE2	1:C:164:LEU:HD21	2.02	0.42
1:D:297:PRO:HB2	1:D:299:GLU:OE2	2.19	0.42
1:F:185:GLN:HG2	1:F:277:TRP:CD1	2.54	0.42
1:F:362:ARG:HG3	1:F:362:ARG:NH1	2.33	0.42
1:B:165:PHE:CD1	1:B:165:PHE:N	2.87	0.42
1:B:200:PHE:CD1	1:B:200:PHE:N	2.87	0.42
1:B:320:TYR:CD1	1:B:320:TYR:C	2.98	0.42
1:D:195:ARG:HB3	1:D:211:GLU:HB2	2.00	0.42
1:E:305:MET:HB3	1:F:338:TYR:CE2	2.53	0.42
1:E:305:MET:CB	1:F:338:TYR:CE2	3.01	0.42
1:B:232:ARG:HD2	1:B:236:TYR:CE2	2.54	0.42
1:B:254:HIS:CE1	1:B:274:ASP:OD1	2.68	0.42
1:C:215:LEU:HB2	1:C:264:LEU:HB2	2.02	0.42
1:D:335:GLN:CD	1:D:335:GLN:N	2.74	0.42
1:F:195:ARG:HG2	1:F:195:ARG:HH11	1.83	0.42
1:D:190:HIS:CE1	1:D:191:PRO:HG2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:180:ARG:HG2	1:E:180:ARG:HH11	1.85	0.42
1:E:320:TYR:HD2	1:E:328:PRO:HD3	1.85	0.42
1:E:338:TYR:CD1	1:F:301:ILE:HD12	2.55	0.42
1:D:185:GLN:HG3	1:D:275:PHE:HD2	1.85	0.42
1:C:162:LYS:HB3	1:C:208:LEU:HB2	2.02	0.41
1:F:306:HIS:CB	1:F:368:PRO:HB2	2.50	0.41
1:D:145:GLY:HA2	1:D:165:PHE:CE1	2.55	0.41
1:D:180:ARG:O	1:D:184:ILE:HG13	2.20	0.41
1:D:195:ARG:HG2	1:D:195:ARG:NH1	2.35	0.41
1:E:305:MET:O	1:E:306:HIS:C	2.62	0.41
1:F:255:ARG:NH1	1:F:276:GLY:HA2	2.35	0.41
1:C:240:LEU:HG	1:C:270:LEU:HD11	2.03	0.41
1:D:254:HIS:HD2	1:D:255:ARG:O	2.02	0.41
1:E:269:GLU:HB2	3:E:2022:HOH:O	2.19	0.41
1:C:259:PRO:HD2	1:C:260:GLU:OE1	2.21	0.41
1:E:220:ARG:HD2	1:E:220:ARG:HA	1.72	0.41
1:F:270:LEU:HD23	1:F:270:LEU:O	2.21	0.41
1:B:274:ASP:OD1	1:B:274:ASP:C	2.64	0.41
1:C:213:ALA:O	2:C:1394:MPY:H12	2.20	0.41
1:E:159:LEU:C	1:E:159:LEU:HD12	2.45	0.41
1:A:240:LEU:CD2	1:A:270:LEU:HD21	2.50	0.41
1:B:362:ARG:HH22	1:F:343:ARG:HB3	1.78	0.41
1:C:275:PHE:O	1:C:277:TRP:N	2.51	0.41
1:F:367:ASN:HD22	1:F:367:ASN:C	2.28	0.41
1:A:162:LYS:HB3	1:A:208:LEU:HB2	2.02	0.41
1:B:220:ARG:HD3	1:B:220:ARG:HA	1.97	0.41
1:B:259:PRO:HD2	1:B:260:GLU:OE1	2.21	0.41
1:D:307:ASP:C	1:D:309:LYS:H	2.28	0.41
1:F:242:ASN:ND2	1:F:392:ASN:H	2.12	0.41
1:F:380:HIS:CE1	1:F:381:PRO:HD2	2.55	0.41
1:A:185:GLN:HG2	1:A:277:TRP:CD1	2.56	0.41
1:C:179:ARG:HD3	1:C:179:ARG:HA	1.79	0.41
1:E:270:LEU:HD23	1:E:270:LEU:O	2.21	0.41
1:D:320:TYR:CD2	1:D:328:PRO:HD3	2.57	0.40
1:E:301:ILE:O	1:E:304:ARG:HG2	2.21	0.40
1:F:253:ILE:CG2	1:F:254:HIS:N	2.82	0.40
1:D:383:ILE:HD13	1:D:383:ILE:HA	1.93	0.40
1:F:270:LEU:HD23	1:F:270:LEU:C	2.46	0.40
1:D:202:ASP:OD1	1:D:202:ASP:C	2.65	0.40
1:E:380:HIS:ND1	1:E:381:PRO:HD2	2.37	0.40
1:A:293:LEU:CD1	1:B:338:TYR:CE1	3.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ARG:HA	1:A:138:PRO:HD3	1.85	0.40
1:B:312:LEU:HD21	1:B:374:LEU:HD23	2.04	0.40
1:C:147:VAL:HG22	1:C:162:LYS:HG3	2.02	0.40
1:E:130:LEU:HA	1:E:130:LEU:HD23	1.89	0.40
1:F:253:ILE:HG22	1:F:254:HIS:H	1.87	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	252/306 (82%)	244 (97%)	7 (3%)	1 (0%)	30	51
1	B	254/306 (83%)	246 (97%)	7 (3%)	1 (0%)	30	51
1	C	253/306 (83%)	242 (96%)	9 (4%)	2 (1%)	16	34
1	D	252/306 (82%)	243 (96%)	7 (3%)	2 (1%)	16	34
1	E	252/306 (82%)	241 (96%)	8 (3%)	3 (1%)	10	23
1	F	253/306 (83%)	241 (95%)	10 (4%)	2 (1%)	16	34
All	All	1516/1836 (83%)	1457 (96%)	48 (3%)	11 (1%)	18	38

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	294	ASP
1	E	294	ASP
1	F	294	ASP
1	C	294	ASP
1	E	154	GLN
1	F	154	GLN
1	C	276	GLY

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Mol	Chain	Res	Type
1	E	276	GLY
1	A	276	GLY
1	D	276	GLY
1	B	276	GLY

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	227/271 (84%)	220 (97%)	7 (3%)	35	63
1	B	229/271 (84%)	220 (96%)	9 (4%)	28	55
1	C	228/271 (84%)	222 (97%)	6 (3%)	40	68
1	D	227/271 (84%)	220 (97%)	7 (3%)	35	63
1	E	227/271 (84%)	218 (96%)	9 (4%)	28	55
1	F	228/271 (84%)	220 (96%)	8 (4%)	32	59
All	All	1366/1626 (84%)	1320 (97%)	46 (3%)	32	60

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	GLN
1	A	189	ARG
1	A	232	ARG
1	A	246	TYR
1	A	270	LEU
1	A	311	ASP
1	A	354	GLU
1	B	180	ARG
1	B	185	GLN
1	B	189	ARG
1	B	232	ARG
1	B	246	TYR
1	B	270	LEU
1	B	311	ASP

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Mol	Chain	Res	Type
1	B	354	GLU
1	B	367	ASN
1	C	179	ARG
1	C	185	GLN
1	C	189	ARG
1	C	270	LEU
1	C	340	ARG
1	C	354	GLU
1	D	185	GLN
1	D	189	ARG
1	D	232	ARG
1	D	270	LEU
1	D	311	ASP
1	D	340	ARG
1	D	350	ASP
1	E	152	GLU
1	E	180	ARG
1	E	185	GLN
1	E	189	ARG
1	E	270	LEU
1	E	299	GLU
1	E	340	ARG
1	E	350	ASP
1	E	368	PRO
1	F	151	ARG
1	F	185	GLN
1	F	189	ARG
1	F	232	ARG
1	F	270	LEU
1	F	311	ASP
1	F	340	ARG
1	F	354	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	ASN
1	A	335	GLN
1	A	367	ASN
1	A	392	ASN
1	B	146	ASN
1	B	185	GLN

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Mol	Chain	Res	Type
1	B	248	HIS
1	B	254	HIS
1	B	367	ASN
1	B	370	GLN
1	B	392	ASN
1	C	176	HIS
1	C	242	ASN
1	C	254	HIS
1	C	367	ASN
1	C	392	ASN
1	D	185	GLN
1	D	242	ASN
1	D	248	HIS
1	D	254	HIS
1	D	367	ASN
1	D	386	ASN
1	D	392	ASN
1	D	394	GLN
1	E	154	GLN
1	E	177	GLN
1	E	185	GLN
1	E	231	GLN
1	E	242	ASN
1	E	367	ASN
1	E	370	GLN
1	E	386	ASN
1	E	392	ASN
1	F	176	HIS
1	F	177	GLN
1	F	242	ASN
1	F	254	HIS
1	F	367	ASN
1	F	386	ASN
1	F	392	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](#)

6 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	MPY	F	1395	-	41,41,41	2.88	17 (41%)	46,58,58	3.07	11 (23%)
2	MPY	B	1395	-	41,41,41	2.93	18 (43%)	46,58,58	3.10	12 (26%)
2	MPY	C	1394	-	41,41,41	2.97	19 (46%)	46,58,58	3.07	15 (32%)
2	MPY	E	1395	-	41,41,41	3.04	21 (51%)	46,58,58	3.08	14 (30%)
2	MPY	D	1395	-	41,41,41	3.03	20 (48%)	46,58,58	3.10	14 (30%)
2	MPY	A	1395	-	41,41,41	2.92	20 (48%)	46,58,58	2.98	12 (26%)

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	MPY	F	1395	-	-	3/23/43/43	0/5/5/5
2	MPY	B	1395	-	-	1/23/43/43	0/5/5/5
2	MPY	C	1394	-	-	1/23/43/43	0/5/5/5
2	MPY	E	1395	-	-	1/23/43/43	0/5/5/5
2	MPY	D	1395	-	-	1/23/43/43	0/5/5/5
2	MPY	A	1395	-	-	1/23/43/43	0/5/5/5

All (115) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1394	MPY	C16-C13	8.74	1.47	1.37
2	B	1395	MPY	C16-C13	8.73	1.47	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1395	MPY	C16-C13	8.45	1.47	1.37
2	F	1395	MPY	C16-C13	8.38	1.47	1.37
2	E	1395	MPY	C16-C13	8.28	1.47	1.37
2	A	1395	MPY	C16-C13	7.95	1.46	1.37
2	D	1395	MPY	C16-N1	7.71	1.50	1.37
2	B	1395	MPY	C16-N1	7.37	1.49	1.37
2	F	1395	MPY	C16-N1	7.37	1.49	1.37
2	A	1395	MPY	C16-N1	7.32	1.49	1.37
2	C	1394	MPY	C16-N1	7.13	1.49	1.37
2	E	1395	MPY	C16-N1	7.12	1.49	1.37
2	E	1395	MPY	C15-N1	6.87	1.48	1.37
2	D	1395	MPY	C15-N1	6.40	1.48	1.37
2	A	1395	MPY	C15-C14	6.30	1.47	1.38
2	C	1394	MPY	C15-C14	6.11	1.47	1.38
2	E	1395	MPY	C15-C14	5.99	1.47	1.38
2	D	1395	MPY	C15-C14	5.99	1.47	1.38
2	F	1395	MPY	C15-C14	5.98	1.47	1.38
2	F	1395	MPY	C15-N1	5.88	1.47	1.37
2	A	1395	MPY	C15-N1	5.86	1.47	1.37
2	C	1394	MPY	C15-N1	5.85	1.47	1.37
2	B	1395	MPY	C15-N1	5.83	1.47	1.37
2	B	1395	MPY	C15-C14	5.82	1.47	1.38
2	F	1395	MPY	C27-C28	4.29	1.47	1.40
2	A	1395	MPY	C3-N5	4.26	1.41	1.33
2	F	1395	MPY	C14-C13	-4.06	1.35	1.43
2	E	1395	MPY	C3-N5	3.96	1.40	1.33
2	E	1395	MPY	C27-C28	3.92	1.46	1.40
2	B	1395	MPY	C3-N5	3.87	1.40	1.33
2	B	1395	MPY	C14-C13	-3.83	1.35	1.43
2	D	1395	MPY	C3-N5	3.81	1.40	1.33
2	C	1394	MPY	C27-C28	3.77	1.46	1.40
2	D	1395	MPY	C27-C28	3.68	1.46	1.40
2	A	1395	MPY	C27-C28	3.68	1.46	1.40
2	A	1395	MPY	C14-C13	-3.64	1.35	1.43
2	D	1395	MPY	C11-C24	3.63	1.46	1.39
2	E	1395	MPY	C9-C7	3.61	1.44	1.39
2	E	1395	MPY	C11-C24	3.57	1.46	1.39
2	C	1394	MPY	C18-N17	3.54	1.52	1.46
2	B	1395	MPY	C11-C24	3.52	1.45	1.39
2	C	1394	MPY	C11-C24	3.42	1.45	1.39
2	F	1395	MPY	C3-N5	3.32	1.39	1.33
2	B	1395	MPY	C27-C28	3.30	1.45	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1394	MPY	C14-C13	-3.26	1.36	1.43
2	F	1395	MPY	C25-N30	3.25	1.43	1.37
2	D	1395	MPY	C14-C13	-3.24	1.36	1.43
2	B	1395	MPY	C25-N30	3.24	1.43	1.37
2	E	1395	MPY	C14-C13	-3.24	1.36	1.43
2	C	1394	MPY	C12-C7	3.23	1.44	1.39
2	B	1395	MPY	C9-C7	3.21	1.44	1.39
2	C	1394	MPY	C3-N5	3.21	1.39	1.33
2	D	1395	MPY	C18-N17	3.21	1.52	1.46
2	E	1395	MPY	C18-N17	3.17	1.52	1.46
2	D	1395	MPY	C12-C7	3.16	1.44	1.39
2	D	1395	MPY	C9-C7	3.14	1.44	1.39
2	B	1395	MPY	C27-C29	3.13	1.45	1.40
2	C	1394	MPY	C22-N17	3.10	1.52	1.46
2	A	1395	MPY	C11-C24	3.09	1.45	1.39
2	D	1395	MPY	C25-N30	3.09	1.42	1.37
2	F	1395	MPY	C11-C24	3.04	1.45	1.39
2	A	1395	MPY	C27-C29	3.03	1.45	1.40
2	E	1395	MPY	C12-C7	3.02	1.44	1.39
2	C	1394	MPY	C9-C7	3.02	1.44	1.39
2	A	1395	MPY	C18-N17	3.02	1.51	1.46
2	C	1394	MPY	C27-C29	3.02	1.45	1.40
2	E	1395	MPY	C10-C24	3.02	1.45	1.39
2	D	1395	MPY	C22-N17	2.97	1.51	1.46
2	E	1395	MPY	C22-N17	2.97	1.51	1.46
2	C	1394	MPY	C25-N30	2.96	1.42	1.37
2	F	1395	MPY	C27-C29	2.95	1.45	1.40
2	E	1395	MPY	C24-N17	2.90	1.46	1.38
2	F	1395	MPY	C12-C7	2.86	1.43	1.39
2	C	1394	MPY	C10-C24	2.86	1.44	1.39
2	A	1395	MPY	C9-C7	2.83	1.43	1.39
2	A	1395	MPY	C25-N30	2.77	1.42	1.37
2	D	1395	MPY	C21-N20	2.73	1.53	1.46
2	F	1395	MPY	C9-C7	2.69	1.43	1.39
2	B	1395	MPY	C12-C7	2.68	1.43	1.39
2	D	1395	MPY	C10-C24	2.66	1.44	1.39
2	B	1395	MPY	C18-N17	2.65	1.51	1.46
2	D	1395	MPY	C12-C11	2.64	1.43	1.38
2	C	1394	MPY	C6-N5	2.62	1.43	1.37
2	A	1395	MPY	C24-N17	2.60	1.46	1.38
2	A	1395	MPY	C6-N5	2.60	1.43	1.37
2	B	1395	MPY	C6-N5	2.59	1.43	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1395	MPY	C27-C29	2.58	1.44	1.40
2	C	1394	MPY	C24-N17	2.49	1.45	1.38
2	D	1395	MPY	C24-N17	2.48	1.45	1.38
2	B	1395	MPY	C21-N20	2.46	1.52	1.46
2	C	1394	MPY	C21-N20	2.45	1.52	1.46
2	E	1395	MPY	C25-N30	2.44	1.41	1.37
2	D	1395	MPY	C6-N5	2.43	1.42	1.37
2	F	1395	MPY	C21-N20	2.40	1.52	1.46
2	F	1395	MPY	C6-N5	2.40	1.42	1.37
2	D	1395	MPY	C27-C29	2.38	1.44	1.40
2	A	1395	MPY	C12-C7	2.37	1.43	1.39
2	A	1395	MPY	C10-C24	2.37	1.43	1.39
2	A	1395	MPY	C22-N17	2.35	1.50	1.46
2	A	1395	MPY	C7-C6	2.34	1.53	1.49
2	D	1395	MPY	C33-C29	2.32	1.43	1.39
2	B	1395	MPY	C24-N17	2.32	1.45	1.38
2	E	1395	MPY	C12-C11	2.32	1.42	1.38
2	F	1395	MPY	C10-C24	2.31	1.43	1.39
2	B	1395	MPY	C10-C24	2.31	1.43	1.39
2	C	1394	MPY	C12-C11	2.29	1.42	1.38
2	E	1395	MPY	C21-N20	2.28	1.51	1.46
2	E	1395	MPY	C7-C6	2.28	1.52	1.49
2	A	1395	MPY	C21-N20	2.28	1.51	1.46
2	E	1395	MPY	C6-N5	2.24	1.42	1.37
2	A	1395	MPY	C31-C28	2.24	1.43	1.39
2	E	1395	MPY	C33-C29	2.15	1.43	1.39
2	F	1395	MPY	C24-N17	2.13	1.44	1.38
2	B	1395	MPY	C22-N17	2.03	1.50	1.46
2	F	1395	MPY	C22-N17	2.02	1.50	1.46

All (78) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1395	MPY	C13-C16-N1	-16.49	97.53	106.01
2	F	1395	MPY	C13-C16-N1	-16.44	97.55	106.01
2	D	1395	MPY	C13-C16-N1	-16.29	97.63	106.01
2	E	1395	MPY	C13-C16-N1	-16.18	97.69	106.01
2	C	1394	MPY	C13-C16-N1	-16.13	97.71	106.01
2	A	1395	MPY	C13-C16-N1	-15.43	98.07	106.01
2	E	1395	MPY	N5-C3-N4	-5.58	119.88	128.28
2	F	1395	MPY	N5-C3-N4	-5.24	120.39	128.28
2	D	1395	MPY	N5-C3-N4	-5.16	120.51	128.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1394	MPY	N5-C3-N4	-5.04	120.70	128.28
2	A	1395	MPY	N5-C3-N4	-5.03	120.71	128.28
2	B	1395	MPY	N5-C3-N4	-4.90	120.90	128.28
2	E	1395	MPY	C14-C3-N5	4.63	130.41	122.78
2	A	1395	MPY	C21-N20-C19	4.31	116.41	109.54
2	D	1395	MPY	C21-N20-C19	4.26	116.33	109.54
2	C	1394	MPY	C16-N1-C15	4.17	113.38	108.92
2	F	1395	MPY	C14-C3-N5	4.10	129.54	122.78
2	F	1395	MPY	C21-N20-C19	4.06	116.01	109.54
2	B	1395	MPY	C21-N20-C19	4.03	115.96	109.54
2	D	1395	MPY	C14-C3-N5	4.01	129.40	122.78
2	F	1395	MPY	C16-N1-C15	3.99	113.18	108.92
2	A	1395	MPY	C14-C3-N5	3.95	129.30	122.78
2	E	1395	MPY	C16-N1-C15	3.93	113.11	108.92
2	B	1395	MPY	C14-C3-N5	3.91	129.24	122.78
2	C	1394	MPY	C14-C3-N5	3.91	129.23	122.78
2	B	1395	MPY	C16-N1-C15	3.73	112.91	108.92
2	D	1395	MPY	C16-N1-C15	3.72	112.90	108.92
2	D	1395	MPY	C27-N30-C25	3.62	129.50	122.43
2	A	1395	MPY	C16-N1-C15	3.61	112.78	108.92
2	B	1395	MPY	C27-N30-C25	3.55	129.37	122.43
2	E	1395	MPY	C27-N30-C25	3.41	129.10	122.43
2	B	1395	MPY	C34-C29-C27	3.41	125.79	121.38
2	A	1395	MPY	C27-N30-C25	3.29	128.85	122.43
2	C	1394	MPY	C34-C29-C27	3.27	125.61	121.38
2	C	1394	MPY	C21-N20-C19	3.26	114.73	109.54
2	F	1395	MPY	C34-C29-C27	3.25	125.59	121.38
2	F	1395	MPY	C27-N30-C25	3.12	128.52	122.43
2	A	1395	MPY	C34-C29-C27	3.10	125.38	121.38
2	E	1395	MPY	C21-N20-C19	3.03	114.36	109.54
2	F	1395	MPY	C14-C15-N1	-2.98	104.15	108.73
2	E	1395	MPY	C34-C29-C27	2.91	125.14	121.38
2	D	1395	MPY	C22-C21-N20	2.90	115.52	110.86
2	F	1395	MPY	C32-C28-C27	2.89	125.12	121.38
2	C	1394	MPY	C22-C21-N20	2.88	115.48	110.86
2	E	1395	MPY	C14-C15-N1	-2.87	104.32	108.73
2	B	1395	MPY	C29-C27-N30	2.86	122.89	119.06
2	C	1394	MPY	C29-C27-N30	2.84	122.86	119.06
2	A	1395	MPY	C29-C27-N30	2.81	122.81	119.06
2	C	1394	MPY	C22-N17-C18	2.78	117.81	111.57
2	D	1395	MPY	C34-C29-C27	2.75	124.94	121.38
2	C	1394	MPY	C14-C15-N1	-2.70	104.58	108.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1395	MPY	C22-C21-N20	2.70	115.20	110.86
2	A	1395	MPY	C14-C15-N1	-2.69	104.60	108.73
2	A	1395	MPY	C22-C21-N20	2.68	115.17	110.86
2	D	1395	MPY	C14-C15-N1	-2.63	104.69	108.73
2	C	1394	MPY	C21-C22-N17	2.54	116.12	110.78
2	C	1394	MPY	C27-N30-C25	2.53	127.38	122.43
2	B	1395	MPY	C14-C15-N1	-2.53	104.84	108.73
2	B	1395	MPY	C22-C21-N20	2.51	114.89	110.86
2	D	1395	MPY	C22-N17-C18	2.45	117.09	111.57
2	E	1395	MPY	C29-C27-N30	2.42	122.30	119.06
2	F	1395	MPY	C29-C27-N30	2.38	122.25	119.06
2	E	1395	MPY	C32-C28-C27	2.37	124.44	121.38
2	D	1395	MPY	C29-C27-N30	2.34	122.20	119.06
2	B	1395	MPY	C32-C28-C27	2.34	124.40	121.38
2	F	1395	MPY	C18-C19-N20	2.31	114.56	110.86
2	A	1395	MPY	C32-C28-C27	2.27	124.31	121.38
2	C	1394	MPY	C19-C18-N17	2.26	115.54	110.78
2	C	1394	MPY	C32-C28-C27	2.26	124.30	121.38
2	E	1395	MPY	C21-C22-N17	2.23	115.48	110.78
2	C	1394	MPY	C12-C7-C6	2.21	124.44	120.45
2	D	1395	MPY	C19-C18-N17	2.21	115.43	110.78
2	D	1395	MPY	C12-C7-C6	2.19	124.40	120.45
2	B	1395	MPY	C18-C19-N20	2.18	114.37	110.86
2	E	1395	MPY	C22-N17-C18	2.16	116.44	111.57
2	A	1395	MPY	C22-N17-C18	2.12	116.33	111.57
2	E	1395	MPY	C19-C18-N17	2.05	115.11	110.78
2	D	1395	MPY	C32-C28-C27	2.04	124.02	121.38

There are no chirality outliers.

All (8) torsion outliers are listed below:

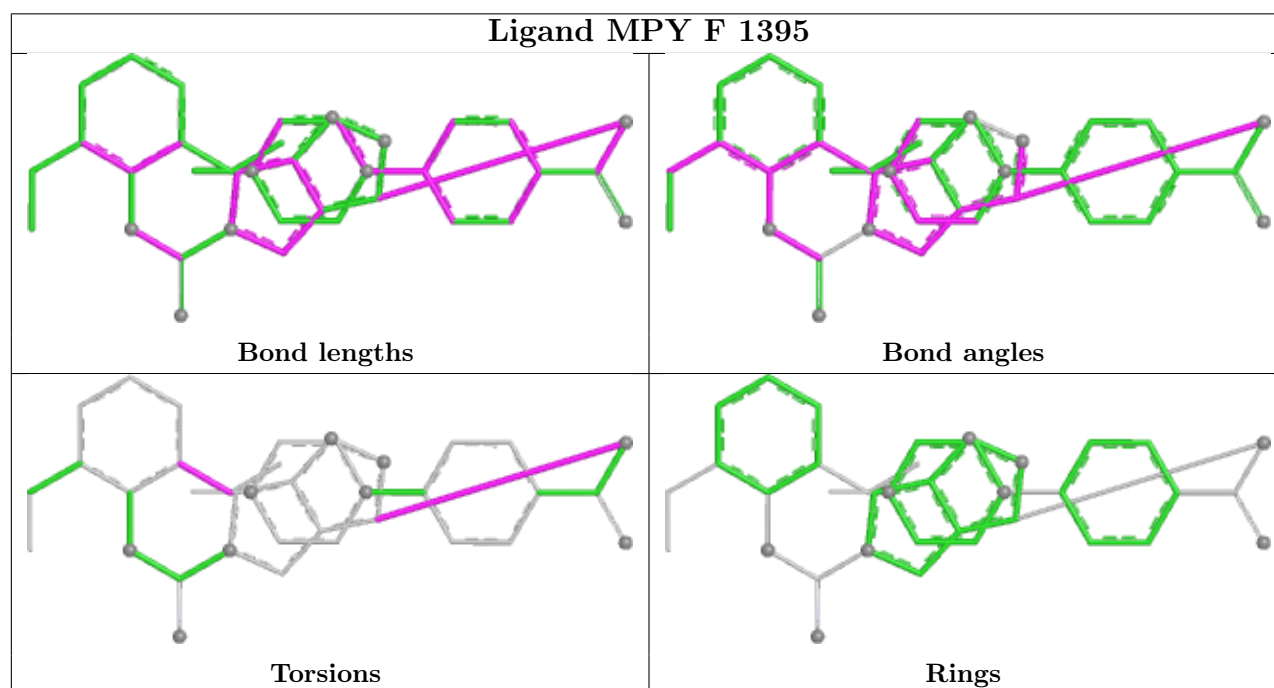
Mol	Chain	Res	Type	Atoms
2	A	1395	MPY	C14-C3-N5-C6
2	B	1395	MPY	C14-C3-N5-C6
2	C	1394	MPY	C14-C3-N5-C6
2	D	1395	MPY	C14-C3-N5-C6
2	E	1395	MPY	C14-C3-N5-C6
2	F	1395	MPY	C14-C3-N5-C6
2	F	1395	MPY	C31-C28-C32-C36
2	F	1395	MPY	C27-C28-C32-C36

There are no ring outliers.

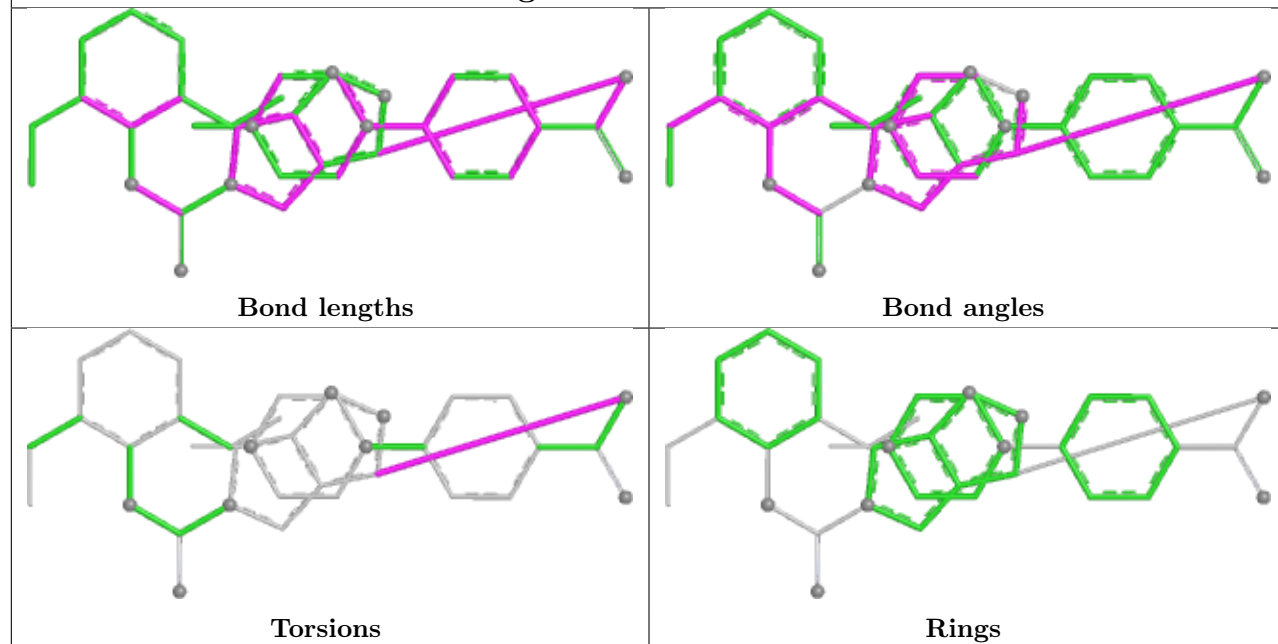
6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1395	MPY	1	0
2	B	1395	MPY	2	0
2	C	1394	MPY	3	0
2	E	1395	MPY	1	0
2	D	1395	MPY	2	0
2	A	1395	MPY	2	0

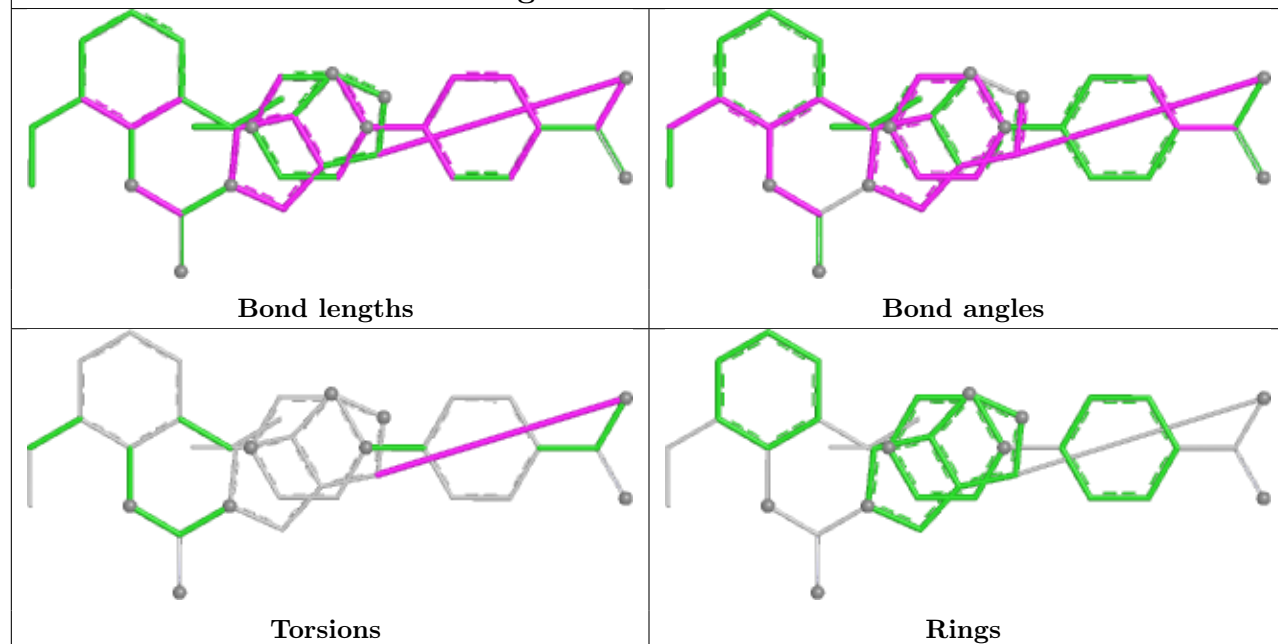
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



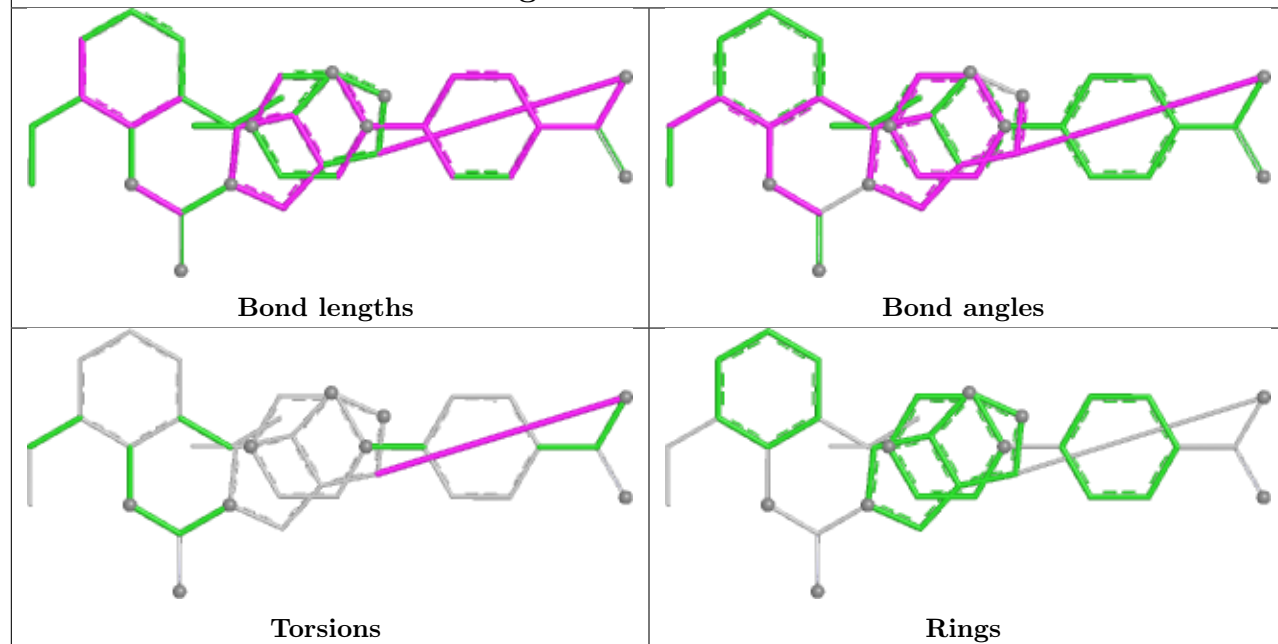
Ligand MPY B 1395



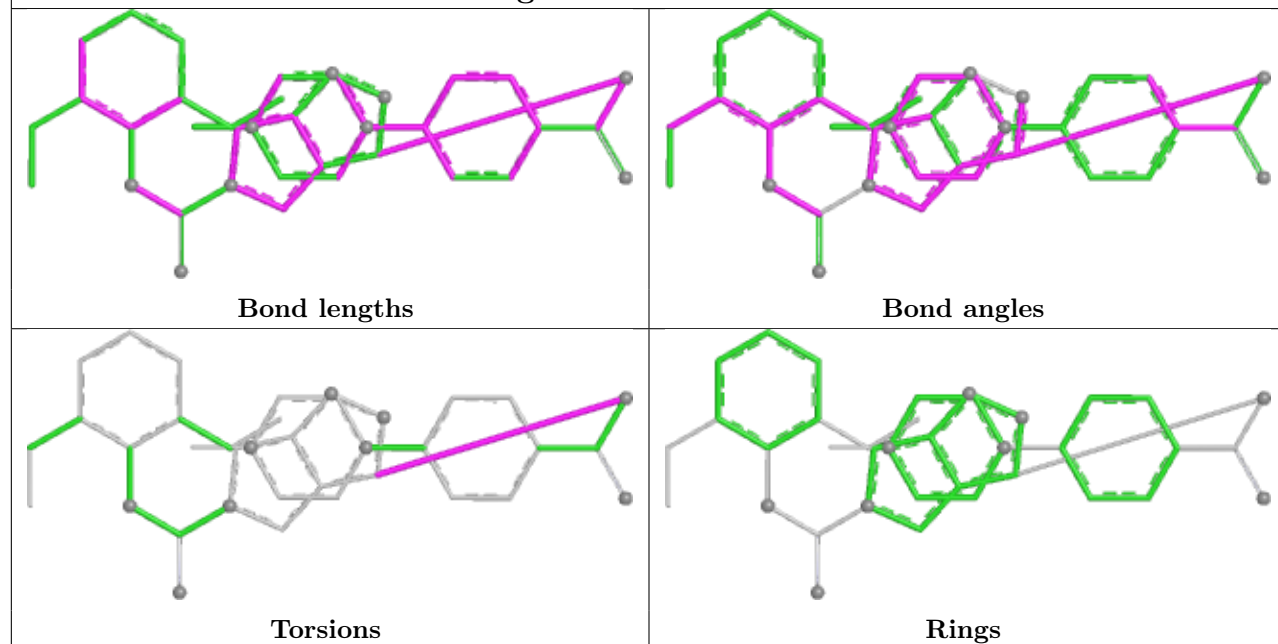
Ligand MPY C 1394

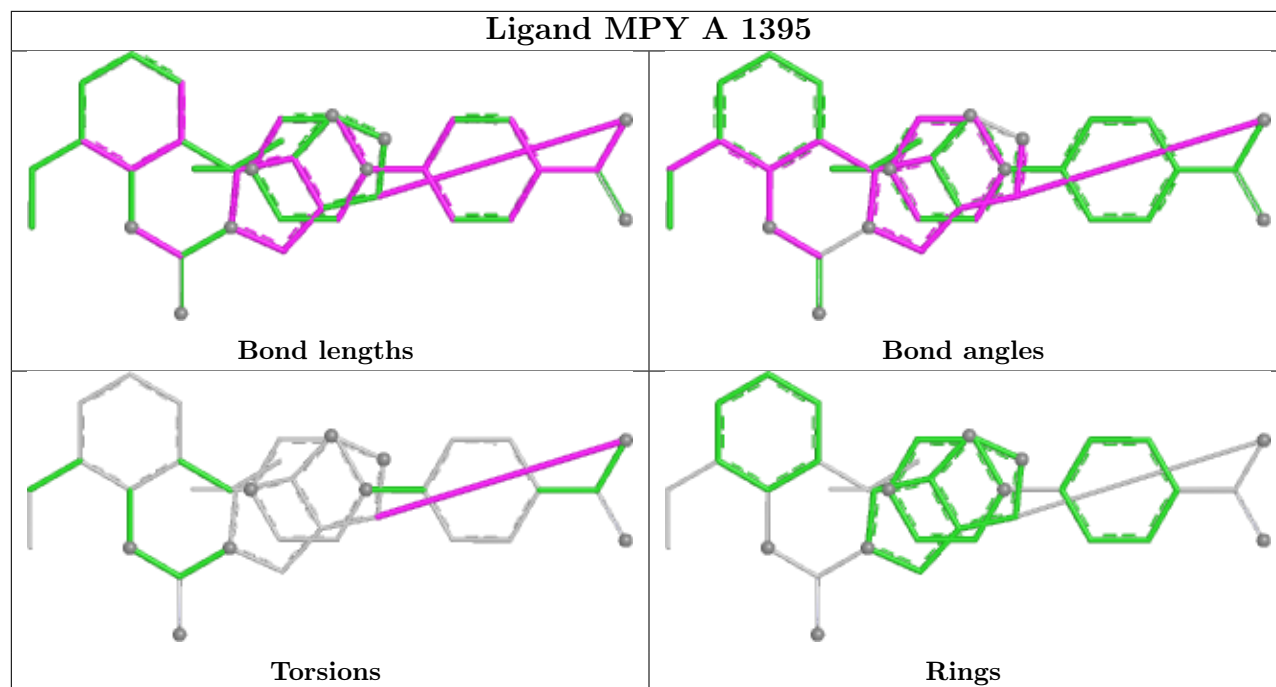


Ligand MPY E 1395



Ligand MPY D 1395





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.