



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 3, 2023 – 05:57 am GMT

PDB ID : 1E6V
Title : Methyl-coenzyme M reductase from Methanopyrus kandleri
Authors : Grabarse, W.; Ermler, U.
Deposited on : 2000-08-23
Resolution : 2.70 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

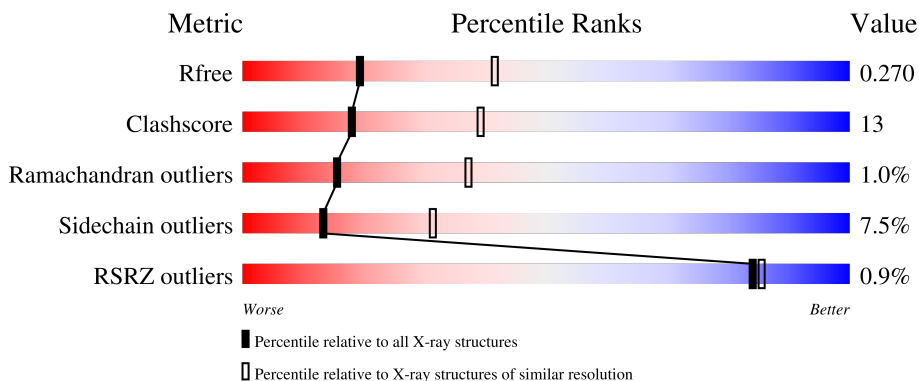
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 19402 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 METHYL-COENZYME M REDUCTASE I ALPHA SUB-UNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	545	4251	2704	712	807	28	18	0	0
1	D	545	4251	2704	712	807	28	18	0	0

- Molecule 2 is a protein called METHYL-COENZYME M REDUCTASE I BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	436	3316	2089	557	646	24	19	0	0
2	E	436	3316	2089	557	646	24	19	0	0

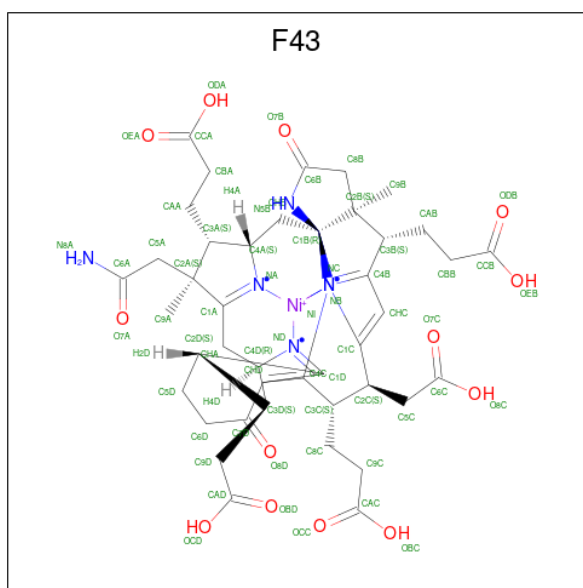
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	49	VAL	ILE	cloning artifact	UNP Q49601
B	98	GLN	GLU	cloning artifact	UNP Q49601
B	220	GLN	GLU	cloning artifact	UNP Q49601
E	49	VAL	ILE	cloning artifact	UNP Q49601
E	98	GLN	GLU	cloning artifact	UNP Q49601
E	220	GLN	GLU	cloning artifact	UNP Q49601

- Molecule 3 is a protein called METHYL-COENZYME M REDUCTASE I GAMMA SUB-UNIT.

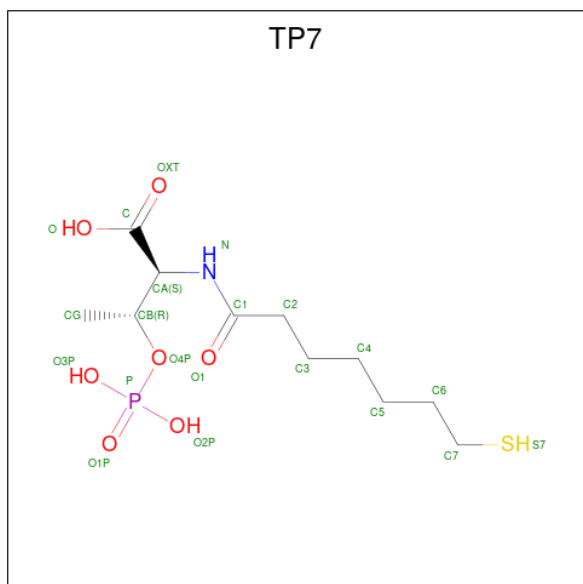
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	248	2044	1280	362	395	7	32	0	0
3	F	248	2044	1280	362	395	7	32	0	0

- Molecule 4 is FACTOR 430 (three-letter code: F43) (formula: $C_{42}H_{51}N_6NiO_{13}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
4	A	1	Total	C	N	Ni	O	0	0
			62	42	6	1	13		
4	D	1	Total	C	N	Ni	O	0	0
			62	42	6	1	13		

- Molecule 5 is Coenzyme B (three-letter code: TP7) (formula: $C_{11}H_{22}NO_7PS$).



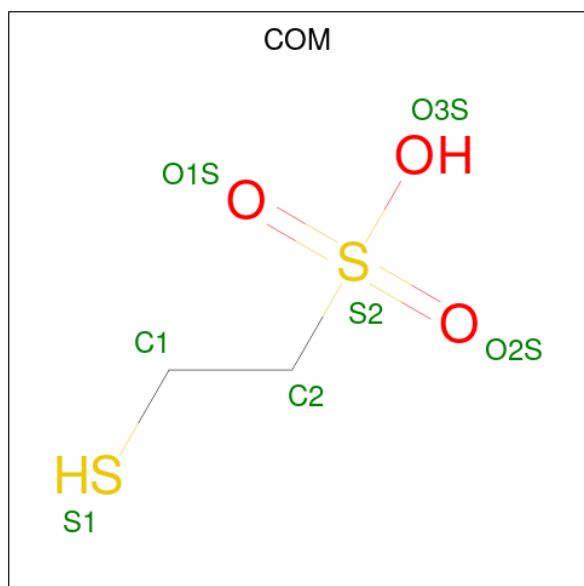
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
5	A	1	Total	C	N	O	P	S	0	0
			21	11	1	7	1	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
5	D	1	21	11	1	7	1	1	0	0

- Molecule 6 is 1-THIOETHANESULFONIC ACID (three-letter code: COM) (formula: C₂H₆O₃S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
6	A	1	7	2	3	2	0	0
6	D	1	7	2	3	2	0	0

SEQUENCE-PLOTS INFOmissingINFO

3 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.52Å 115.74Å 268.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 29.93 – 2.70	Depositor EDS
% Data completeness (in resolution range)	62.9 (30.00-2.70) 62.9 (29.93-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.32 (at 2.68Å)	Xtrriage
Refinement program	CNS 0.3	Depositor
R, R_{free}	0.239 , 0.278 0.235 , 0.270	Depositor DCC
R_{free} test set	2196 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtrriage
Anisotropy	0.745	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 28.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	19402	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

4 Model quality i

4.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: F43, TP7, COM

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	0.30	0/4352	0.51	2/5897 (0.0%)
1	D	0.30	0/4352	0.51	2/5897 (0.0%)
2	B	0.29	0/3374	0.51	1/4573 (0.0%)
2	E	0.29	0/3374	0.51	1/4573 (0.0%)
3	C	0.27	0/2089	0.47	1/2827 (0.0%)
3	F	0.27	0/2089	0.47	1/2827 (0.0%)
All	All	0.29	0/19630	0.50	8/26594 (0.0%)

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	0	5
1	D	0	5
2	B	0	3
2	E	0	3
3	C	0	2
3	F	0	2
All	All	0	20

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	121	ARG	NE-CZ-NH1	-10.49	115.06	120.30
2	E	121	ARG	NE-CZ-NH1	-10.44	115.08	120.30
1	D	228	ARG	NE-CZ-NH1	-8.47	116.06	120.30
1	A	228	ARG	NE-CZ-NH1	-8.36	116.12	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	20	ARG	NE-CZ-NH1	-6.17	117.22	120.30
3	F	20	ARG	NE-CZ-NH1	-6.10	117.25	120.30
1	D	54	ARG	NE-CZ-NH1	-5.49	117.55	120.30
1	A	54	ARG	NE-CZ-NH1	-5.49	117.56	120.30

There are no chirality outliers.

All (20) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	212	ARG	Sidechain
1	A	228	ARG	Sidechain
1	A	274	ARG	Sidechain
1	A	447	TYR	Sidechain
1	A	54	ARG	Sidechain
2	B	121	ARG	Sidechain
2	B	129	TYR	Sidechain
2	B	334	TYR	Sidechain
3	C	19	ARG	Sidechain
3	C	20	ARG	Sidechain
1	D	212	ARG	Sidechain
1	D	228	ARG	Sidechain
1	D	274	ARG	Sidechain
1	D	447	TYR	Sidechain
1	D	54	ARG	Sidechain
2	E	121	ARG	Sidechain
2	E	129	TYR	Sidechain
2	E	334	TYR	Sidechain
3	F	19	ARG	Sidechain
3	F	20	ARG	Sidechain

4.2 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	4251	0	4101	139	1
1	D	4251	0	4101	140	1
2	B	3316	0	3254	96	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	3316	0	3254	97	2
3	C	2044	0	1985	69	0
3	F	2044	0	1985	71	0
4	A	62	0	43	5	0
4	D	62	0	43	6	0
5	A	21	0	19	1	0
5	D	21	0	19	1	0
6	A	7	0	5	1	0
6	D	7	0	5	1	0
All	All	19402	0	18814	511	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (511) 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)
2:B:118:PRO:HG2	2:B:121:ARG:HG3	1.48	0.95
2:E:118:PRO:HG2	2:E:121:ARG:HG3	1.47	0.94
1:D:309:GLU:HG3	1:D:528:PRO:HD2	1.58	0.85
1:A:309:GLU:HG3	1:A:528:PRO:HD2	1.58	0.85
2:E:286:LYS:HD2	2:E:296:LYS:HB3	1.59	0.83
1:D:338:THR:O	1:D:342:PRO:HD2	1.78	0.83
1:A:338:THR:O	1:A:342:PRO:HD2	1.79	0.83
2:B:286:LYS:HD2	2:B:296:LYS:HB3	1.60	0.82
1:A:339:ALA:CB	1:A:403:GLN:HB2	2.11	0.81
1:D:339:ALA:CB	1:D:403:GLN:HB2	2.11	0.81
1:D:335:GLN:HA	1:D:338:THR:OG1	1.85	0.77
1:A:335:GLN:HA	1:A:338:THR:OG1	1.85	0.77
1:A:296:MET:HG2	1:A:306:VAL:HG11	1.67	0.77
1:D:451:LEU:HD12	2:E:359:GLY:HA3	1.67	0.77
1:D:400:GLY:H	1:D:403:GLN:HE21	1.30	0.77
1:A:400:GLY:H	1:A:403:GLN:HE21	1.30	0.76
1:A:428:ASN:HD21	3:C:246:THR:HG21	1.49	0.76
1:D:454:GLN:NE2	3:F:245:ARG:HH22	1.83	0.76
1:D:296:MET:HG2	1:D:306:VAL:HG11	1.67	0.76
1:A:451:LEU:HD12	2:B:359:GLY:HA3	1.67	0.75
1:A:454:GLN:NE2	3:C:245:ARG:HH22	1.83	0.75
1:D:428:ASN:HD21	3:F:246:THR:HG21	1.49	0.74
1:D:443:ARG:NH1	3:F:101:HIS:HD2	1.85	0.74
1:D:363:ILE:HD12	1:D:363:ILE:H	1.53	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:ILE:H	1:A:363:ILE:HD12	1.53	0.73
1:A:443:ARG:NH1	3:C:101:HIS:HD2	1.85	0.73
3:C:33:ARG:HH11	3:C:149:PRO:HD3	1.54	0.72
1:A:139:LEU:HD11	1:A:209:GLN:HE21	1.55	0.72
2:B:31:MET:HE2	2:E:121:ARG:HA	1.73	0.71
3:C:17:GLU:OE1	3:C:17:GLU:HA	1.91	0.71
1:D:139:LEU:HD11	1:D:209:GLN:HE21	1.55	0.71
3:F:33:ARG:HH11	3:F:149:PRO:HD3	1.54	0.71
2:B:129:TYR:CE1	2:B:181:ASP:HA	2.27	0.70
3:C:19:ARG:HH11	3:C:104:ILE:HD11	1.57	0.70
3:F:17:GLU:HA	3:F:17:GLU:OE1	1.91	0.70
1:D:58:HIS:ND1	1:D:59:TYR:N	2.40	0.69
1:A:58:HIS:ND1	1:A:59:TYR:N	2.40	0.69
2:B:194:ARG:NH2	2:B:220:GLN:HE22	1.90	0.69
6:A:1555:COM:O1S	1:D:446:PHE:HB2	1.93	0.69
2:B:121:ARG:HA	2:E:31:MET:HE2	1.75	0.68
2:E:194:ARG:NH2	2:E:220:GLN:HE22	1.90	0.68
2:E:129:TYR:CE1	2:E:181:ASP:HA	2.26	0.68
3:C:19:ARG:NH2	3:C:221:TYR:HE1	1.92	0.68
3:F:19:ARG:NH2	3:F:221:TYR:HE1	1.92	0.68
1:D:443:ARG:HH11	3:F:101:HIS:HD2	1.41	0.68
1:A:341:TYR:CD1	1:A:342:PRO:HD3	2.29	0.68
1:A:446:PHE:HB2	6:D:1555:COM:O1S	1.94	0.67
1:D:341:TYR:CD1	1:D:342:PRO:HD3	2.29	0.67
1:A:443:ARG:HH11	3:C:101:HIS:HD2	1.41	0.67
3:F:129:VAL:HG22	3:F:130:ARG:H	1.60	0.67
3:F:19:ARG:HH11	3:F:104:ILE:HD11	1.57	0.67
1:A:274:ARG:NH1	2:B:381:ILE:HD13	2.10	0.67
2:B:304:ASN:HD21	2:B:349:ASP:H	1.43	0.67
3:C:129:VAL:HG22	3:C:130:ARG:H	1.60	0.66
3:F:252:PRO:O	3:F:253:GLU:HB2	1.94	0.66
1:A:549:ILE:HA	1:D:538:ARG:HG2	1.78	0.66
3:C:252:PRO:O	3:C:253:GLU:HB2	1.94	0.66
3:C:122:SER:HB3	4:D:1553:F43:O8C	1.96	0.65
1:D:143:LEU:HD13	1:D:164:CYS:SG	2.36	0.65
2:E:304:ASN:HD21	2:E:349:ASP:H	1.43	0.65
1:D:274:ARG:NH1	2:E:381:ILE:HD13	2.10	0.65
2:E:214:CYS:O	2:E:218:THR:HG23	1.97	0.65
1:A:143:LEU:HD13	1:A:164:CYS:SG	2.36	0.65
4:A:1553:F43:O8C	3:F:122:SER:HB3	1.96	0.65
1:A:65:VAL:HG21	1:D:149:VAL:O	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:129:TYR:CE2	2:B:182:VAL:HG23	2.33	0.64
1:D:338:THR:O	1:D:342:PRO:CD	2.45	0.64
1:A:149:VAL:O	1:D:65:VAL:HG21	1.98	0.64
1:A:538:ARG:HG2	1:D:549:ILE:HA	1.78	0.64
2:B:214:CYS:O	2:B:218:THR:HG23	1.96	0.64
1:A:34:TRP:HE1	1:A:40:LYS:HB3	1.63	0.64
2:E:129:TYR:CE2	2:E:182:VAL:HG23	2.33	0.64
2:B:244:TYR:HB3	2:B:309:ALA:HB2	1.80	0.63
2:E:244:TYR:HB3	2:E:309:ALA:HB2	1.80	0.63
1:A:139:LEU:CD1	1:A:209:GLN:HE21	2.12	0.63
1:D:34:TRP:HE1	1:D:40:LYS:HB3	1.63	0.63
1:A:105:ARG:O	1:A:226:MET:HG3	1.98	0.63
1:A:338:THR:O	1:A:342:PRO:CD	2.45	0.63
1:D:139:LEU:CD1	1:D:209:GLN:HE21	2.12	0.63
1:D:105:ARG:O	1:D:226:MET:HG3	1.98	0.62
3:C:29:LEU:H	3:C:29:LEU:HD12	1.64	0.62
1:A:484:ASN:HA	5:A:1554:TP7:S7	2.40	0.62
1:D:400:GLY:H	1:D:403:GLN:NE2	1.98	0.61
2:E:304:ASN:ND2	2:E:349:ASP:H	1.98	0.61
3:F:29:LEU:HD12	3:F:29:LEU:H	1.64	0.61
1:A:443:ARG:NH1	3:C:101:HIS:CD2	2.68	0.61
1:A:209:GLN:NE2	1:A:234:ILE:HG12	2.16	0.61
2:B:128:PRO:HB2	2:B:180:LEU:HD13	1.82	0.61
1:D:443:ARG:NH1	3:F:101:HIS:CD2	2.68	0.61
2:B:304:ASN:ND2	2:B:349:ASP:H	1.98	0.61
1:A:441:GLN:HE21	1:A:441:GLN:HA	1.66	0.60
1:D:484:ASN:HA	5:D:1554:TP7:S7	2.40	0.60
1:D:209:GLN:NE2	1:D:234:ILE:HG12	2.16	0.60
1:D:441:GLN:HA	1:D:441:GLN:HE21	1.66	0.60
2:E:128:PRO:HB2	2:E:180:LEU:HD13	1.82	0.60
3:F:8:TYR:HB2	3:F:20:ARG:HG2	1.83	0.60
1:D:30:ILE:O	1:D:30:ILE:HG22	2.02	0.60
2:E:292:PHE:C	2:E:292:PHE:CD2	2.75	0.59
2:B:271:ARG:NH2	2:B:342:GLU:OE1	2.35	0.59
1:A:76:GLN:HB2	1:A:83:PHE:CZ	2.38	0.59
1:D:39:ARG:HE	1:D:538:ARG:NH2	2.00	0.59
2:B:257:VAL:HG21	2:B:316:MET:HE1	1.83	0.59
2:B:292:PHE:C	2:B:292:PHE:CD2	2.74	0.59
1:A:472:GLU:HG3	1:A:495:VAL:HG13	1.84	0.59
3:C:8:TYR:HB2	3:C:20:ARG:HG2	1.82	0.59
1:A:362:GLY:HA3	1:A:365:GLU:HG3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:472:GLU:HG3	1:D:495:VAL:HG13	1.84	0.59
1:A:400:GLY:H	1:A:403:GLN:NE2	1.98	0.59
1:A:395:MET:HE2	3:C:95:VAL:HG21	1.84	0.59
1:D:76:GLN:HB2	1:D:83:PHE:CZ	2.38	0.59
1:A:39:ARG:HE	1:A:538:ARG:NH2	2.00	0.58
1:D:395:MET:HE2	3:F:95:VAL:HG21	1.85	0.58
1:D:443:ARG:NH2	3:F:122:SER:O	2.36	0.58
2:E:271:ARG:NH2	2:E:342:GLU:OE1	2.35	0.58
1:A:30:ILE:HG22	1:A:30:ILE:O	2.02	0.58
1:A:443:ARG:NH2	3:C:122:SER:O	2.36	0.58
3:C:19:ARG:HH22	3:C:221:TYR:HE1	1.51	0.58
1:D:362:GLY:HA3	1:D:365:GLU:HG3	1.85	0.58
1:D:196:ALA:O	1:D:200:ILE:HG12	2.04	0.57
1:A:377:ALA:HB3	1:A:432:LEU:HD23	1.86	0.57
2:E:342:GLU:OE1	3:F:7:PHE:HB2	2.04	0.57
1:A:448:GLY:HA2	2:B:358:VAL:HG12	1.85	0.57
2:B:342:GLU:OE1	3:C:7:PHE:HB2	2.04	0.57
3:C:15:VAL:O	3:C:19:ARG:HG3	2.05	0.57
2:B:320:GLY:O	2:B:323:ARG:HD3	2.04	0.57
2:B:257:VAL:HG21	2:B:316:MET:CE	2.35	0.57
1:D:377:ALA:HB3	1:D:432:LEU:HD23	1.86	0.57
2:E:257:VAL:HG21	2:E:316:MET:CE	2.35	0.57
3:F:174:HIS:NE2	3:F:178:ARG:HD2	2.19	0.57
1:D:448:GLY:HA2	2:E:358:VAL:HG12	1.85	0.57
3:C:174:HIS:NE2	3:C:178:ARG:HD2	2.19	0.56
1:A:337:ALA:O	1:A:340:VAL:HG22	2.06	0.56
2:E:320:GLY:O	2:E:323:ARG:HD3	2.04	0.56
1:A:527:GLU:HG2	1:A:530:LYS:HD2	1.88	0.56
1:D:337:ALA:O	1:D:340:VAL:HG22	2.06	0.56
2:E:122:VAL:O	2:E:124:VAL:N	2.38	0.56
3:F:19:ARG:HH22	3:F:221:TYR:HE1	1.52	0.56
3:C:158:THR:HG22	1:D:149:VAL:O	2.05	0.56
1:A:196:ALA:O	1:A:200:ILE:HG12	2.04	0.56
1:A:296:MET:HE2	1:A:306:VAL:HG21	1.88	0.56
1:D:527:GLU:HG2	1:D:530:LYS:HD2	1.88	0.56
3:F:15:VAL:O	3:F:19:ARG:HG3	2.05	0.56
1:A:182:LYS:HD2	1:D:57:PRO:HD3	1.88	0.55
2:B:122:VAL:O	2:B:124:VAL:N	2.38	0.55
1:D:39:ARG:HE	1:D:538:ARG:HH22	1.54	0.55
1:A:39:ARG:HE	1:A:538:ARG:HH22	1.54	0.55
3:C:78:GLU:HB3	3:C:132:ARG:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:309:GLU:CG	1:D:528:PRO:HD2	2.35	0.54
1:D:386:GLU:HG2	3:F:223:ILE:HD13	1.89	0.54
1:A:386:GLU:HG2	3:C:223:ILE:HD13	1.89	0.54
3:F:78:GLU:HB3	3:F:132:ARG:HD2	1.89	0.54
1:D:315:ALA:O	1:D:319:ASP:HB2	2.08	0.54
1:A:57:PRO:HD3	1:D:182:LYS:HD2	1.89	0.54
1:A:464:ARG:HH22	2:B:238:HIS:HE1	1.56	0.54
1:D:464:ARG:HH22	2:E:238:HIS:HE1	1.56	0.54
1:A:246:ALA:HB1	4:A:1553:F43:H9B1	1.90	0.54
1:A:315:ALA:O	1:A:319:ASP:HB2	2.08	0.54
1:A:34:TRP:NE1	1:A:40:LYS:HB3	2.22	0.54
1:A:149:VAL:O	3:F:158:THR:HG22	2.06	0.54
1:A:400:GLY:N	1:A:403:GLN:HE21	2.04	0.54
2:B:336:ASN:ND2	2:B:348:VAL:H	2.06	0.54
1:D:34:TRP:NE1	1:D:40:LYS:HB3	2.22	0.54
1:A:309:GLU:CG	1:A:528:PRO:HD2	2.35	0.54
4:A:1553:F43:O8D	1:D:400:GLY:HA3	2.08	0.53
2:B:253:VAL:O	2:B:257:VAL:HG23	2.09	0.53
2:B:200:HIS:HA	2:B:413:SER:OG	2.09	0.53
3:C:129:VAL:HG22	3:C:130:ARG:N	2.22	0.53
2:E:253:VAL:O	2:E:257:VAL:HG23	2.09	0.53
2:E:336:ASN:ND2	2:E:348:VAL:H	2.06	0.53
2:E:390:ALA:O	2:E:394:VAL:HG23	2.08	0.53
1:A:50:ILE:HG21	1:A:54:ARG:HH21	1.73	0.53
2:B:24:PRO:HB2	2:B:26:GLU:OE2	2.09	0.53
2:B:390:ALA:O	2:B:394:VAL:HG23	2.08	0.53
1:D:34:TRP:CH2	1:D:35:LYS:HE3	2.43	0.53
2:E:257:VAL:HG21	2:E:316:MET:HE1	1.90	0.53
1:A:34:TRP:CH2	1:A:35:LYS:HE3	2.43	0.53
1:D:400:GLY:N	1:D:403:GLN:HE21	2.04	0.53
1:A:464:ARG:HA	2:E:137:MET:CE	2.39	0.53
3:C:29:LEU:HD23	3:C:146:ILE:HD12	1.89	0.53
3:F:129:VAL:HG22	3:F:130:ARG:N	2.22	0.53
1:A:428:ASN:HD21	3:C:246:THR:CG2	2.22	0.53
2:B:137:MET:CE	1:D:464:ARG:HA	2.39	0.53
1:A:167:LYS:HB3	1:A:185:ILE:HG13	1.91	0.53
1:A:338:THR:HB	1:A:342:PRO:CG	2.39	0.53
2:E:194:ARG:HH21	2:E:220:GLN:HE22	1.56	0.53
3:F:213:ASP:O	3:F:217:ARG:HG2	2.09	0.53
3:F:174:HIS:CD2	3:F:178:ARG:HH21	2.27	0.53
1:A:331:VAL:HB	4:D:1553:F43:H9A1	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:338:THR:HB	1:D:342:PRO:CG	2.39	0.52
2:B:210:MET:HB3	2:B:430:MET:HE1	1.92	0.52
3:C:174:HIS:CD2	3:C:178:ARG:HH21	2.27	0.52
2:E:24:PRO:HB2	2:E:26:GLU:OE2	2.09	0.52
2:E:266:VAL:O	2:E:270:VAL:HG23	2.09	0.52
4:A:1553:F43:H9A1	1:D:331:VAL:HB	1.90	0.52
2:B:127:ASP:O	2:B:130:SER:HB2	2.09	0.52
2:E:200:HIS:HA	2:E:413:SER:OG	2.09	0.52
2:E:210:MET:HB3	2:E:430:MET:HE1	1.92	0.52
3:C:213:ASP:O	3:C:217:ARG:HG2	2.09	0.52
2:E:127:ASP:O	2:E:130:SER:HB2	2.09	0.52
1:A:219:HIS:HA	1:D:219:HIS:O	2.10	0.52
2:B:124:VAL:HG11	2:E:30:PRO:HB2	1.92	0.52
1:D:246:ALA:HB1	4:D:1553:F43:H9B1	1.90	0.52
3:F:29:LEU:HD23	3:F:146:ILE:HD12	1.90	0.52
1:A:209:GLN:HE22	1:A:234:ILE:HG12	1.74	0.52
1:D:209:GLN:HE22	1:D:234:ILE:HG12	1.74	0.52
2:B:118:PRO:HG2	2:B:121:ARG:CG	2.33	0.52
2:B:266:VAL:O	2:B:270:VAL:HG23	2.09	0.52
1:D:50:ILE:HG21	1:D:54:ARG:HH21	1.73	0.52
1:D:296:MET:HE2	1:D:306:VAL:HG21	1.90	0.52
2:B:289:LEU:HB3	3:C:233:GLU:OE1	2.11	0.51
3:F:11:GLY:O	3:F:20:ARG:NH2	2.37	0.51
1:A:50:ILE:HG22	1:A:54:ARG:HE	1.76	0.51
1:A:219:HIS:O	1:D:219:HIS:HA	2.09	0.51
2:B:30:PRO:HB2	2:E:124:VAL:HG11	1.92	0.51
1:D:167:LYS:HB3	1:D:185:ILE:HG13	1.91	0.51
1:A:245:ALA:HB2	3:F:88:TYR:CZ	2.46	0.51
2:E:323:ARG:HG3	3:F:60:LEU:HD11	1.93	0.51
2:B:194:ARG:HH21	2:B:220:GLN:HE22	1.56	0.51
2:B:328:VAL:HG21	2:B:394:VAL:HA	1.93	0.51
3:C:88:TYR:CZ	1:D:245:ALA:HB2	2.46	0.51
2:E:289:LEU:HB3	3:F:233:GLU:OE1	2.10	0.51
1:A:135:TYR:OH	1:A:234:ILE:HG23	2.11	0.51
1:A:400:GLY:HA3	4:D:1553:F43:O8D	2.09	0.51
1:A:106:ARG:HD2	1:A:213:MET:O	2.11	0.51
1:A:10:MET:SD	1:A:21:PRO:HG2	2.52	0.50
1:A:232:MET:HG3	2:E:367:ILE:HD12	1.94	0.50
2:B:323:ARG:HG3	3:C:60:LEU:HD11	1.93	0.50
1:D:10:MET:SD	1:D:21:PRO:HG2	2.52	0.50
1:D:135:TYR:OH	1:D:234:ILE:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:THR:HB	1:A:342:PRO:HG3	1.93	0.50
2:E:328:VAL:HG21	2:E:394:VAL:HA	1.93	0.50
3:F:29:LEU:HD12	3:F:29:LEU:N	2.26	0.50
3:C:29:LEU:HD12	3:C:29:LEU:N	2.26	0.50
2:E:118:PRO:HG2	2:E:121:ARG:CG	2.32	0.50
2:E:124:VAL:O	2:E:124:VAL:HG13	2.11	0.50
1:D:50:ILE:HG22	1:D:54:ARG:HE	1.76	0.49
1:D:194:GLU:O	1:D:197:GLU:HG3	2.13	0.49
1:D:338:THR:HB	1:D:342:PRO:HG3	1.93	0.49
1:D:106:ARG:HD2	1:D:213:MET:O	2.11	0.49
1:A:502:ARG:CZ	2:E:72:GLY:HA3	2.43	0.49
1:A:509:HIS:HD2	1:A:511:VAL:H	1.60	0.49
3:C:19:ARG:NH1	3:C:219:THR:O	2.45	0.49
1:A:194:GLU:O	1:A:197:GLU:HG3	2.13	0.49
1:A:296:MET:CE	1:A:306:VAL:HG21	2.43	0.49
3:C:33:ARG:NH1	3:C:149:PRO:HD3	2.25	0.49
1:D:428:ASN:HD21	3:F:246:THR:CG2	2.22	0.49
2:B:210:MET:HG2	3:C:70:ILE:HD11	1.95	0.49
2:B:336:ASN:HD21	2:B:348:VAL:HG23	1.77	0.49
1:D:422:HIS:CD2	1:D:425:ALA:H	2.30	0.49
3:F:19:ARG:NH1	3:F:104:ILE:HD11	2.26	0.49
3:F:19:ARG:NH1	3:F:219:THR:O	2.45	0.49
1:A:422:HIS:CD2	1:A:425:ALA:H	2.30	0.49
2:E:32:ARG:O	2:E:440:GLU:HG2	2.13	0.48
2:E:210:MET:HG2	3:F:70:ILE:HD11	1.95	0.48
2:B:124:VAL:O	2:B:124:VAL:HG13	2.11	0.48
2:B:326:GLN:HE22	3:C:86:ILE:HG21	1.79	0.48
2:E:336:ASN:HD21	2:E:348:VAL:HG23	1.77	0.48
1:A:454:GLN:HE22	3:C:245:ARG:HH22	1.60	0.48
2:B:367:ILE:HD12	1:D:232:MET:HG3	1.94	0.48
3:C:19:ARG:NH1	3:C:104:ILE:HD11	2.26	0.48
2:B:255:ASP:HB3	2:B:276:ARG:HH12	1.79	0.48
1:D:296:MET:CE	1:D:306:VAL:HG21	2.43	0.48
1:D:352:TYR:CD1	1:D:352:TYR:C	2.86	0.48
3:F:33:ARG:NH1	3:F:149:PRO:HD3	2.25	0.48
2:B:32:ARG:O	2:B:440:GLU:HG2	2.13	0.48
3:C:102:PRO:HA	3:C:125:GLN:HG3	1.96	0.48
1:A:316:MET:HA	1:A:320:GLN:HE21	1.79	0.48
2:B:72:GLY:HA3	1:D:502:ARG:CZ	2.43	0.48
2:E:255:ASP:HB3	2:E:276:ARG:HH12	1.79	0.48
1:A:352:TYR:CD1	1:A:352:TYR:C	2.86	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:245:ALA:HB3	1:D:248:GLU:HB2	1.96	0.48
2:E:416:ILE:O	2:E:420:PHE:HB2	2.14	0.48
3:F:102:PRO:HA	3:F:125:GLN:HG3	1.96	0.48
1:A:245:ALA:HB3	1:A:248:GLU:HB2	1.96	0.47
3:F:7:PHE:O	3:F:8:TYR:HB2	2.15	0.47
1:A:227:TYR:CD1	1:A:227:TYR:C	2.88	0.47
1:D:509:HIS:HD2	1:D:511:VAL:H	1.61	0.47
3:C:7:PHE:O	3:C:8:TYR:HB2	2.15	0.47
3:F:237:VAL:O	3:F:241:ILE:HG12	2.15	0.47
2:B:151:MET:HG2	1:D:461:LEU:HD22	1.95	0.47
2:B:416:ILE:O	2:B:420:PHE:HB2	2.14	0.47
1:D:227:TYR:CD1	1:D:227:TYR:C	2.88	0.47
1:A:461:LEU:HD22	2:E:151:MET:HG2	1.96	0.47
2:B:140:VAL:O	2:B:144:ILE:HG12	2.15	0.47
1:D:58:HIS:CG	1:D:59:TYR:H	2.25	0.47
1:A:479:PRO:O	2:B:385:HIS:HE1	1.97	0.47
1:D:392:PRO:HB3	3:F:96:TYR:OH	2.15	0.47
2:E:436:ALA:O	2:E:440:GLU:HB2	2.14	0.47
3:C:25:PRO:HA	3:C:112:ARG:NH2	2.30	0.47
2:B:436:ALA:O	2:B:440:GLU:HB2	2.14	0.47
3:C:11:GLY:O	3:C:20:ARG:NH2	2.37	0.47
1:D:316:MET:HA	1:D:320:GLN:HE21	1.79	0.47
2:E:75:MET:CE	2:E:153:ASP:HB3	2.45	0.47
3:F:25:PRO:HA	3:F:112:ARG:NH2	2.30	0.47
1:A:392:PRO:HB3	3:C:96:TYR:OH	2.15	0.46
2:B:75:MET:CE	2:B:153:ASP:HB3	2.45	0.46
2:E:326:GLN:HE22	3:F:86:ILE:HG21	1.79	0.46
1:D:479:PRO:O	2:E:385:HIS:HE1	1.97	0.46
2:E:221:GLN:HG3	2:E:242:TYR:HB2	1.97	0.46
2:E:328:VAL:O	2:E:332:ILE:HG13	2.16	0.46
3:F:19:ARG:HD2	3:F:104:ILE:HG12	1.97	0.46
2:B:431:LYS:O	2:B:435:GLU:HB2	2.15	0.46
2:E:431:LYS:O	2:E:435:GLU:HB2	2.15	0.46
3:C:237:VAL:O	3:C:241:ILE:HG12	2.15	0.46
1:A:161:THR:HG22	1:A:214:PRO:HB3	1.97	0.46
1:A:443:ARG:HH11	3:C:101:HIS:CD2	2.26	0.46
1:D:294:GLN:O	1:D:297:ARG:HB2	2.16	0.46
1:A:285:PRO:HD2	1:A:288:VAL:HB	1.97	0.46
2:B:42:ASP:HA	2:B:45:ARG:HG2	1.98	0.46
1:D:161:THR:HG22	1:D:214:PRO:HB3	1.97	0.46
2:E:69:LYS:O	2:E:71:PRO:HD3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:MET:HG2	3:C:243:GLU:HG2	1.98	0.46
1:A:58:HIS:CG	1:A:59:TYR:H	2.25	0.46
1:A:294:GLN:O	1:A:297:ARG:HB2	2.16	0.46
3:C:19:ARG:HH11	3:C:104:ILE:CD1	2.27	0.46
3:F:121:LEU:HD13	3:F:124:ARG:HG3	1.97	0.46
2:B:221:GLN:HG3	2:B:242:TYR:HB2	1.98	0.46
2:E:75:MET:HE1	2:E:153:ASP:HB3	1.98	0.46
2:E:140:VAL:O	2:E:144:ILE:HG12	2.15	0.46
1:D:388:TYR:OH	1:D:398:HIS:HD2	1.99	0.45
2:E:414:LYS:HG3	2:E:415:LEU:H	1.80	0.45
2:B:69:LYS:O	2:B:71:PRO:HD3	2.16	0.45
2:B:328:VAL:O	2:B:332:ILE:HG13	2.16	0.45
1:D:8:LEU:HB2	1:D:348:ASP:HB2	1.98	0.45
1:D:339:ALA:HB1	1:D:403:GLN:HB2	1.96	0.45
1:D:370:MET:HG2	3:F:243:GLU:HG2	1.98	0.45
2:E:363:PHE:O	2:E:370:GLY:HA3	2.15	0.45
2:B:169:ILE:HD12	2:B:169:ILE:HA	1.81	0.45
1:D:454:GLN:HE22	3:F:245:ARG:HH22	1.60	0.45
2:E:42:ASP:HA	2:E:45:ARG:HG2	1.98	0.45
3:C:19:ARG:HD2	3:C:104:ILE:HG12	1.97	0.45
3:C:176:LEU:HD12	3:C:178:ARG:NH1	2.31	0.45
1:D:285:PRO:HD2	1:D:288:VAL:HB	1.97	0.45
2:B:316:MET:HB3	2:B:316:MET:HE2	1.92	0.45
1:D:228:ARG:HH11	1:D:228:ARG:HG3	1.82	0.45
2:E:293:LYS:CG	3:F:7:PHE:HZ	2.30	0.45
1:A:8:LEU:HB2	1:A:348:ASP:HB2	1.98	0.45
2:B:414:LYS:HG3	2:B:415:LEU:H	1.80	0.45
1:A:271:PRO:HB3	2:E:169:ILE:HD11	1.98	0.45
1:A:339:ALA:HB1	1:A:403:GLN:HB2	1.96	0.45
2:B:55:GLU:HG3	2:B:79:ILE:HG12	1.99	0.45
2:B:363:PHE:O	2:B:370:GLY:HA3	2.16	0.45
3:F:176:LEU:HD12	3:F:178:ARG:NH1	2.31	0.45
2:B:293:LYS:CG	3:C:7:PHE:HZ	2.30	0.45
1:D:54:ARG:HD3	1:D:56:VAL:HB	1.98	0.45
1:A:228:ARG:HG3	1:A:228:ARG:HH11	1.82	0.45
2:B:194:ARG:NH1	2:B:384:ARG:O	2.50	0.45
3:C:121:LEU:HD13	3:C:124:ARG:HG3	1.97	0.45
1:D:114:THR:HG22	1:D:117:ARG:NH2	2.32	0.45
2:E:356:THR:HG23	2:E:383:THR:HA	1.98	0.45
1:A:388:TYR:OH	1:A:398:HIS:HD2	1.99	0.44
2:B:169:ILE:HD11	1:D:271:PRO:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:155:ARG:HD3	3:C:159:VAL:HA	1.99	0.44
1:D:116:HIS:HA	1:D:119:LEU:HD12	1.99	0.44
3:F:19:ARG:HH11	3:F:104:ILE:CD1	2.27	0.44
2:B:356:THR:HG23	2:B:383:THR:HA	1.98	0.44
1:A:54:ARG:HD3	1:A:56:VAL:HB	1.99	0.44
1:A:339:ALA:HB2	1:A:403:GLN:HB2	1.98	0.44
1:A:116:HIS:HA	1:A:119:LEU:HD12	1.99	0.44
2:B:425:GLU:CD	2:B:425:GLU:H	2.21	0.44
1:D:228:ARG:HH11	1:D:228:ARG:CG	2.30	0.44
1:A:167:LYS:HG3	1:A:183:PHE:C	2.38	0.44
2:B:137:MET:HE2	1:D:464:ARG:HA	1.98	0.44
2:B:181:ASP:O	2:B:182:VAL:O	2.36	0.44
2:B:194:ARG:HH21	2:B:220:GLN:NE2	2.16	0.44
2:E:334:TYR:O	2:E:338:LEU:HG	2.18	0.44
2:E:290:PRO:O	2:E:291:SER:CB	2.66	0.44
2:B:316:MET:HE3	3:C:70:ILE:HD12	1.99	0.44
2:E:55:GLU:HG3	2:E:79:ILE:HG12	1.99	0.44
3:F:192:LYS:HG2	3:F:198:GLU:HA	2.00	0.44
1:A:228:ARG:HH11	1:A:228:ARG:CG	2.30	0.44
2:B:21:GLU:HB3	2:B:22:GLU:H	1.62	0.44
2:B:290:PRO:O	2:B:291:SER:CB	2.66	0.44
1:D:167:LYS:HG3	1:D:183:PHE:C	2.38	0.44
1:A:39:ARG:HA	1:A:39:ARG:HD3	1.39	0.43
1:A:114:THR:HG22	1:A:117:ARG:NH2	2.32	0.43
2:B:46:THR:HG23	2:B:114:LEU:HD11	2.00	0.43
3:C:192:LYS:HG2	3:C:198:GLU:HA	2.00	0.43
1:D:464:ARG:HH22	2:E:238:HIS:CE1	2.35	0.43
2:E:265:THR:O	2:E:269:VAL:HG23	2.18	0.43
3:F:155:ARG:HD3	3:F:159:VAL:HA	1.99	0.43
2:E:425:GLU:CD	2:E:425:GLU:H	2.21	0.43
1:A:464:ARG:HH22	2:B:238:HIS:CE1	2.35	0.43
2:E:194:ARG:NH1	2:E:384:ARG:O	2.50	0.43
1:A:383:TYR:O	1:A:387:GLN:HG2	2.19	0.43
2:B:210:MET:CG	3:C:70:ILE:HD11	2.49	0.43
2:B:265:THR:O	2:B:269:VAL:HG23	2.18	0.43
2:B:290:PRO:O	2:B:291:SER:HB2	2.19	0.43
2:E:194:ARG:HH21	2:E:220:GLN:NE2	2.16	0.43
2:E:221:GLN:O	2:E:225:PHE:HD1	2.02	0.43
2:B:27:VAL:HG11	2:B:437:ALA:O	2.19	0.43
2:B:27:VAL:HG13	2:B:440:GLU:HB3	2.01	0.43
2:B:75:MET:HE1	2:B:153:ASP:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:221:GLN:O	2:B:225:PHE:HD1	2.02	0.43
2:B:334:TYR:O	2:B:338:LEU:HG	2.18	0.43
2:E:27:VAL:HG11	2:E:437:ALA:O	2.19	0.43
1:D:457:ALA:HB1	3:F:246:THR:HG23	2.00	0.43
2:E:273:VAL:HG21	2:E:335:TYR:OH	2.18	0.43
1:A:427:LEU:HD13	1:A:471:LEU:HD13	2.00	0.43
1:A:441:GLN:HE21	1:A:441:GLN:CA	2.32	0.43
2:B:94:ILE:HG23	2:B:118:PRO:HD2	2.01	0.43
1:D:422:HIS:HD2	1:D:424:GLN:N	2.17	0.43
1:D:305:LYS:O	1:D:309:GLU:HB2	2.19	0.42
2:E:46:THR:HG23	2:E:114:LEU:HD11	2.00	0.42
2:E:181:ASP:O	2:E:182:VAL:O	2.36	0.42
2:E:194:ARG:HD2	2:E:220:GLN:OE1	2.19	0.42
1:A:318:TYR:O	1:A:323:LEU:HG	2.19	0.42
1:D:318:TYR:O	1:D:323:LEU:HG	2.19	0.42
1:D:454:GLN:HE21	3:F:245:ARG:HH22	1.65	0.42
2:E:21:GLU:HB3	2:E:22:GLU:H	1.61	0.42
1:A:464:ARG:HA	2:E:137:MET:HE2	2.02	0.42
2:B:194:ARG:HD2	2:B:220:GLN:OE1	2.19	0.42
1:A:173:LEU:HD21	1:A:200:ILE:HD12	2.01	0.42
1:A:457:ALA:HB1	3:C:246:THR:HG23	2.00	0.42
1:A:468:GLY:O	1:A:469:LEU:HB2	2.19	0.42
2:E:187:GLU:HG3	2:E:379:ASN:O	2.20	0.42
2:E:210:MET:CG	3:F:70:ILE:HD11	2.49	0.42
3:F:10:PRO:HG3	3:F:237:VAL:HG11	2.02	0.42
1:A:345:ILE:HD12	1:A:398:HIS:CE1	2.55	0.42
2:E:27:VAL:HG13	2:E:440:GLU:HB3	2.01	0.42
2:B:273:VAL:HG21	2:B:335:TYR:OH	2.18	0.42
1:D:383:TYR:O	1:D:387:GLN:HG2	2.19	0.42
1:A:389:GLU:H	1:A:389:GLU:HG2	1.74	0.42
3:C:10:PRO:HG3	3:C:237:VAL:HG11	2.02	0.42
1:D:114:THR:HG22	1:D:117:ARG:HH22	1.85	0.42
1:D:339:ALA:HB2	1:D:403:GLN:HB2	1.98	0.42
1:D:468:GLY:O	1:D:469:LEU:HB2	2.19	0.42
2:E:316:MET:HE3	3:F:70:ILE:HD12	2.01	0.42
3:F:58:PRO:HA	3:F:59:PRO:HD3	1.90	0.42
1:A:305:LYS:O	1:A:309:GLU:HB2	2.19	0.42
1:A:326:TYR:CZ	1:D:221:CYS:HA	2.55	0.42
2:B:187:GLU:HG3	2:B:379:ASN:O	2.20	0.42
1:D:389:GLU:H	1:D:389:GLU:HG2	1.74	0.42
2:E:94:ILE:HG23	2:E:118:PRO:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:233:PRO:CB	3:F:251:ARG:HD2	2.49	0.42
1:A:250:ALA:O	1:A:253:ASP:HB2	2.20	0.42
4:A:1553:F43:CAD	1:D:335:GLN:H	2.33	0.42
2:B:372:GLY:HA3	2:B:405:GLN:HE22	1.85	0.42
1:D:250:ALA:O	1:D:253:ASP:HB2	2.20	0.42
1:D:443:ARG:HH11	3:F:101:HIS:CD2	2.26	0.42
2:E:293:LYS:HD2	3:F:7:PHE:HZ	1.84	0.42
1:A:457:ALA:O	3:C:246:THR:HG22	2.20	0.41
2:B:289:LEU:O	2:B:291:SER:N	2.47	0.41
3:C:98:ALA:O	3:C:101:HIS:HE1	2.03	0.41
1:D:284:VAL:HA	1:D:285:PRO:HD3	1.83	0.41
2:E:104:VAL:HG13	2:E:115:VAL:HG22	2.01	0.41
2:E:257:VAL:HG21	2:E:316:MET:HE2	2.02	0.41
3:F:90:GLN:HB3	3:F:155:ARG:O	2.20	0.41
3:F:154:VAL:HG12	3:F:194:GLN:HB3	2.02	0.41
1:A:527:GLU:N	1:A:528:PRO:HD3	2.35	0.41
2:B:293:LYS:HD2	3:C:7:PHE:HZ	1.84	0.41
1:D:427:LEU:HD13	1:D:471:LEU:HD13	2.00	0.41
1:A:140:ASN:HB3	1:D:57:PRO:O	2.20	0.41
1:A:495:VAL:O	1:A:499:HIS:HD2	2.03	0.41
2:B:233:PRO:CB	3:C:251:ARG:HD2	2.49	0.41
3:F:201:GLU:HA	3:F:202:PRO:HD3	1.93	0.41
1:A:422:HIS:HD2	1:A:424:GLN:N	2.17	0.41
3:C:90:GLN:HB3	3:C:155:ARG:O	2.20	0.41
3:C:100:ILE:HG13	3:C:101:HIS:N	2.35	0.41
1:D:527:GLU:N	1:D:528:PRO:HD3	2.35	0.41
3:F:37:ASP:O	3:F:41:VAL:HG23	2.20	0.41
1:A:123:LEU:HD13	1:A:244:ILE:HG21	2.02	0.41
1:A:221:CYS:HA	1:D:326:TYR:CZ	2.55	0.41
2:B:104:VAL:HG13	2:B:115:VAL:HG22	2.01	0.41
1:D:495:VAL:O	1:D:499:HIS:HD2	2.03	0.41
2:E:290:PRO:O	2:E:291:SER:HB2	2.19	0.41
3:F:187:GLU:HB3	3:F:206:GLY:HA3	2.03	0.41
1:A:370:MET:HG3	1:A:374:LYS:HD3	2.03	0.41
1:A:478:TYR:HA	1:A:479:PRO:HD3	1.95	0.41
1:D:345:ILE:HD12	1:D:398:HIS:CE1	2.55	0.41
1:A:34:TRP:CZ3	1:A:35:LYS:HG3	2.56	0.41
2:E:372:GLY:HA3	2:E:405:GLN:HE22	1.85	0.41
1:D:227:TYR:C	1:D:227:TYR:HD1	2.24	0.41
1:A:227:TYR:C	1:A:227:TYR:HD1	2.24	0.41
1:A:335:GLN:H	4:D:1553:F43:CAD	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:293:LYS:HD2	3:C:7:PHE:CZ	2.56	0.41
1:D:34:TRP:CE3	1:D:35:LYS:HG3	2.56	0.41
2:E:51:LEU:HD11	2:E:113:LEU:HB2	2.03	0.41
3:F:100:ILE:HG13	3:F:101:HIS:N	2.35	0.41
1:A:316:MET:O	1:A:320:GLN:HB2	2.21	0.41
3:C:37:ASP:O	3:C:41:VAL:HG23	2.20	0.41
1:D:173:LEU:HD21	1:D:200:ILE:HD12	2.01	0.41
1:D:395:MET:HG3	3:F:163:ALA:HB1	2.03	0.41
3:F:98:ALA:O	3:F:101:HIS:HE1	2.03	0.41
1:A:57:PRO:O	1:D:140:ASN:HB3	2.21	0.40
1:D:123:LEU:HD13	1:D:244:ILE:HG21	2.02	0.40
1:D:427:LEU:O	1:D:430:TRP:HB3	2.21	0.40
1:D:457:ALA:O	3:F:246:THR:HG22	2.20	0.40
3:C:123:GLY:N	4:D:1553:F43:O7C	2.41	0.40
3:C:154:VAL:HG12	3:C:194:GLN:HB3	2.02	0.40
1:D:106:ARG:HA	1:D:226:MET:HG3	2.02	0.40
1:D:370:MET:HG3	1:D:374:LYS:HD3	2.03	0.40
2:E:293:LYS:HD2	3:F:7:PHE:CZ	2.56	0.40
2:E:414:LYS:HB2	2:E:414:LYS:HE2	1.87	0.40
1:A:263:VAL:HG22	1:A:264:ILE:N	2.36	0.40
2:B:323:ARG:HA	2:B:323:ARG:HD2	1.95	0.40
1:D:316:MET:O	1:D:320:GLN:HB2	2.21	0.40
1:A:34:TRP:CE3	1:A:35:LYS:HG3	2.56	0.40
1:A:114:THR:HG22	1:A:117:ARG:HH22	1.85	0.40
1:A:509:HIS:HA	1:A:510:PRO:HD3	1.91	0.40
3:C:187:GLU:HB3	3:C:206:GLY:HA3	2.03	0.40
1:D:403:GLN:O	1:D:407:VAL:HG23	2.22	0.40
1:A:232:MET:HE2	1:A:232:MET:HA	2.04	0.40
1:A:395:MET:HG3	3:C:163:ALA:HB1	2.03	0.40
1:A:447:TYR:CE2	2:B:361:SER:HB3	2.56	0.40
1:D:109:ILE:HG12	1:D:210:VAL:HG22	2.04	0.40
1:D:263:VAL:HG22	1:D:264:ILE:N	2.36	0.40
1:D:441:GLN:HE21	1:D:441:GLN:CA	2.32	0.40
1:D:448:GLY:N	2:E:362:PHE:HB2	2.37	0.40
2:E:182:VAL:HG13	2:E:184:MET:SD	2.61	0.40
2:E:260:HIS:HD2	2:E:268:THR:OG1	2.04	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:ASP:OD2	1:D:174:GLU:OE1[1_455]	1.74	0.46
2:E:92:LYS:CE	2:E:262:LYS:NZ[3_655]	1.83	0.37
2:E:92:LYS:CE	2:E:262:LYS:CE[3_655]	2.01	0.19

4.3 Torsion angles [i](#)

4.3.1 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	543/553 (98%)	506 (93%)	34 (6%)	3 (1%)	25	50
1	D	543/553 (98%)	506 (93%)	34 (6%)	3 (1%)	25	50
2	B	434/443 (98%)	406 (94%)	21 (5%)	7 (2%)	9	24
2	E	434/443 (98%)	406 (94%)	21 (5%)	7 (2%)	9	24
3	C	246/258 (95%)	231 (94%)	13 (5%)	2 (1%)	19	43
3	F	246/258 (95%)	231 (94%)	13 (5%)	2 (1%)	19	43
All	All	2446/2508 (98%)	2286 (94%)	136 (6%)	24 (1%)	15	37

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	465	SER
2	B	123	LYS
2	B	182	VAL
2	B	326	GLN
1	D	465	SER
2	E	123	LYS
2	E	182	VAL
2	E	326	GLN
1	A	341	TYR
2	B	21	GLU
2	B	122	VAL
2	B	291	SER
1	D	341	TYR

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Mol	Chain	Res	Type
2	E	21	GLU
2	E	122	VAL
2	E	291	SER
3	C	253	GLU
3	F	253	GLU
1	A	508	VAL
3	C	67	GLU
1	D	508	VAL
3	F	67	GLU
2	B	290	PRO
2	E	290	PRO

4.3.2 Protein sidechains [i](#)

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	432/439 (98%)	401 (93%)	31 (7%)	14	34
1	D	432/439 (98%)	401 (93%)	31 (7%)	14	34
2	B	345/350 (99%)	312 (90%)	33 (10%)	8	19
2	E	345/350 (99%)	312 (90%)	33 (10%)	8	19
3	C	218/224 (97%)	207 (95%)	11 (5%)	24	51
3	F	218/224 (97%)	207 (95%)	11 (5%)	24	51
All	All	1990/2026 (98%)	1840 (92%)	150 (8%)	13	31

All (150) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	SER
1	A	39	ARG
1	A	74	SER
1	A	130	GLU
1	A	139	LEU
1	A	143	LEU
1	A	177	ASP

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Mol	Chain	Res	Type
1	A	180	ASP
1	A	184	LEU
1	A	197	GLU
1	A	201	LYS
1	A	207	THR
1	A	212	ARG
1	A	213	MET
1	A	219	HIS
1	A	227	TYR
1	A	228	ARG
1	A	232	MET
1	A	265	ASN
1	A	296	MET
1	A	298	LYS
1	A	352	TYR
1	A	367	GLU
1	A	379	GLU
1	A	441	GLN
1	A	447	TYR
1	A	473	LEU
1	A	527	GLU
1	A	529	ARG
1	A	534	LYS
1	A	547	ASP
2	B	22	GLU
2	B	28	LEU
2	B	29	SER
2	B	31	MET
2	B	59	GLN
2	B	73	ARG
2	B	76	ASP
2	B	77	VAL
2	B	92	LYS
2	B	112	ARG
2	B	123	LYS
2	B	127	ASP
2	B	130	SER
2	B	164	ARG
2	B	169	ILE
2	B	181	ASP
2	B	184	MET
2	B	218	THR

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Mol	Chain	Res	Type
2	B	237	LEU
2	B	245	GLN
2	B	259	LYS
2	B	278	LEU
2	B	287	GLU
2	B	292	PHE
2	B	300	MET
2	B	302	LEU
2	B	323	ARG
2	B	339	LEU
2	B	387	LYS
2	B	415	LEU
2	B	424	ASP
2	B	435	GLU
2	B	440	GLU
3	C	17	GLU
3	C	52	GLU
3	C	61	GLU
3	C	112	ARG
3	C	124	ARG
3	C	146	ILE
3	C	173	LEU
3	C	174	HIS
3	C	198	GLU
3	C	233	GLU
3	C	234	LEU
1	D	37	SER
1	D	39	ARG
1	D	74	SER
1	D	130	GLU
1	D	139	LEU
1	D	143	LEU
1	D	177	ASP
1	D	180	ASP
1	D	184	LEU
1	D	197	GLU
1	D	201	LYS
1	D	207	THR
1	D	212	ARG
1	D	213	MET
1	D	219	HIS
1	D	227	TYR

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Mol	Chain	Res	Type
1	D	228	ARG
1	D	232	MET
1	D	265	ASN
1	D	296	MET
1	D	298	LYS
1	D	352	TYR
1	D	367	GLU
1	D	379	GLU
1	D	441	GLN
1	D	447	TYR
1	D	473	LEU
1	D	527	GLU
1	D	529	ARG
1	D	534	LYS
1	D	547	ASP
2	E	22	GLU
2	E	28	LEU
2	E	29	SER
2	E	31	MET
2	E	59	GLN
2	E	73	ARG
2	E	76	ASP
2	E	77	VAL
2	E	92	LYS
2	E	112	ARG
2	E	123	LYS
2	E	127	ASP
2	E	130	SER
2	E	164	ARG
2	E	169	ILE
2	E	181	ASP
2	E	184	MET
2	E	218	THR
2	E	237	LEU
2	E	245	GLN
2	E	259	LYS
2	E	278	LEU
2	E	287	GLU
2	E	292	PHE
2	E	300	MET
2	E	302	LEU
2	E	323	ARG

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Mol	Chain	Res	Type
2	E	339	LEU
2	E	387	LYS
2	E	415	LEU
2	E	424	ASP
2	E	435	GLU
2	E	440	GLU
3	F	17	GLU
3	F	52	GLU
3	F	61	GLU
3	F	112	ARG
3	F	124	ARG
3	F	146	ILE
3	F	173	LEU
3	F	174	HIS
3	F	198	GLU
3	F	233	GLU
3	F	234	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	209	GLN
1	A	219	HIS
1	A	320	GLN
1	A	398	HIS
1	A	403	GLN
1	A	422	HIS
1	A	428	ASN
1	A	441	GLN
1	A	452	GLN
1	A	454	GLN
1	A	509	HIS
1	A	520	ASN
2	B	41	ASN
2	B	103	ASN
2	B	186	GLN
2	B	238	HIS
2	B	260	HIS
2	B	304	ASN
2	B	326	GLN
2	B	336	ASN

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Mol	Chain	Res	Type
2	B	405	GLN
3	C	18	ASN
3	C	101	HIS
3	C	182	ASN
1	D	76	GLN
1	D	209	GLN
1	D	219	HIS
1	D	320	GLN
1	D	398	HIS
1	D	403	GLN
1	D	422	HIS
1	D	428	ASN
1	D	441	GLN
1	D	452	GLN
1	D	454	GLN
1	D	509	HIS
1	D	520	ASN
2	E	41	ASN
2	E	103	ASN
2	E	186	GLN
2	E	238	HIS
2	E	260	HIS
2	E	304	ASN
2	E	326	GLN
2	E	336	ASN
2	E	405	GLN
3	F	18	ASN
3	F	101	HIS
3	F	182	ASN

4.3.3 RNA [i](#)

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry

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
4	F43	D	1553	1,6	61,71,71	3.64	20 (32%)	64,118,118	2.04	15 (23%)
4	F43	A	1553	1,6	61,71,71	3.64	20 (32%)	64,118,118	2.04	15 (23%)
5	TP7	D	1554	-	19,20,20	2.00	6 (31%)	24,26,26	1.34	3 (12%)
5	TP7	A	1554	-	19,20,20	2.01	6 (31%)	24,26,26	1.34	3 (12%)
6	COM	D	1555	4	6,6,6	1.68	3 (50%)	7,8,8	1.46	1 (14%)
6	COM	A	1555	4	6,6,6	1.68	3 (50%)	7,8,8	1.46	1 (14%)

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
4	F43	D	1553	1,6	-	10/28/185/185	-
4	F43	A	1553	1,6	-	10/28/185/185	-
5	TP7	D	1554	-	-	0/24/24/24	-
5	TP7	A	1554	-	-	0/24/24/24	-
6	COM	D	1555	4	-	0/4/4/4	-
6	COM	A	1555	4	-	0/4/4/4	-

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1553	F43	CHB-C1B	-15.76	1.42	1.53
4	D	1553	F43	CHB-C1B	-15.71	1.42	1.53
4	D	1553	F43	NI-NA	11.74	2.14	1.89
4	A	1553	F43	NI-NA	11.70	2.14	1.89
4	D	1553	F43	CHD-C1D	9.78	1.56	1.43
4	A	1553	F43	CHD-C1D	9.77	1.56	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1553	F43	NI-NB	9.49	2.10	1.89
4	A	1553	F43	NI-NB	9.49	2.10	1.89
5	A	1554	TP7	C2-C1	-5.31	1.41	1.51
5	D	1554	TP7	C2-C1	-5.29	1.41	1.51
4	A	1553	F43	CAA-C3A	5.26	1.63	1.53
4	D	1553	F43	CAA-C3A	5.25	1.63	1.53
4	A	1553	F43	CHC-C4B	5.15	1.53	1.39
4	D	1553	F43	CHC-C4B	5.12	1.53	1.39
4	D	1553	F43	NI-ND	5.01	2.00	1.89
4	A	1553	F43	NI-ND	5.00	2.00	1.89
4	D	1553	F43	C3D-C4D	-4.02	1.46	1.53
4	A	1553	F43	C3D-C4D	-4.01	1.46	1.53
4	A	1553	F43	C4D-ND	3.70	1.54	1.49
4	D	1553	F43	C4D-ND	3.68	1.54	1.49
5	A	1554	TP7	C1-N	3.43	1.41	1.34
4	A	1553	F43	CAB-C3B	3.40	1.62	1.54
5	D	1554	TP7	C1-N	3.39	1.41	1.34
4	D	1553	F43	CAB-C3B	3.39	1.62	1.54
4	A	1553	F43	CHD-C4C	-3.33	1.31	1.40
4	D	1553	F43	CHD-C4C	-3.33	1.31	1.40
4	D	1553	F43	CHD-C7D	3.25	1.53	1.46
4	A	1553	F43	CHD-C7D	3.24	1.53	1.46
5	A	1554	TP7	P-O4P	3.13	1.65	1.59
5	D	1554	TP7	P-O4P	3.12	1.65	1.59
4	A	1553	F43	C5D-C2D	-3.08	1.47	1.53
4	D	1553	F43	C5D-C2D	-3.05	1.48	1.53
4	A	1553	F43	C6B-N5B	-2.81	1.30	1.34
4	D	1553	F43	C6B-N5B	-2.80	1.30	1.34
5	D	1554	TP7	CA-C	2.58	1.56	1.52
5	A	1554	TP7	CA-C	2.57	1.56	1.52
4	A	1553	F43	C9D-C3D	-2.52	1.48	1.53
5	A	1554	TP7	CA-N	-2.50	1.40	1.45
4	A	1553	F43	C4A-NA	-2.50	1.45	1.49
4	D	1553	F43	C9D-C3D	-2.50	1.48	1.53
5	D	1554	TP7	CA-N	-2.48	1.40	1.45
4	D	1553	F43	C4A-NA	-2.47	1.45	1.49
6	D	1555	COM	O2S-S2	2.41	1.52	1.45
6	A	1555	COM	O2S-S2	2.39	1.52	1.45
4	A	1553	F43	C2D-C1D	-2.37	1.46	1.52
4	A	1553	F43	C4C-NC	2.37	1.38	1.35
4	D	1553	F43	C2D-C1D	-2.36	1.46	1.52
4	D	1553	F43	C4C-NC	2.35	1.38	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1553	F43	C9A-C2A	-2.35	1.50	1.54
4	A	1553	F43	C9A-C2A	-2.35	1.50	1.54
6	A	1555	COM	C2-S2	2.33	1.80	1.77
6	D	1555	COM	C2-S2	2.33	1.80	1.77
5	D	1554	TP7	OXT-C	2.10	1.28	1.22
6	A	1555	COM	O1S-S2	2.09	1.51	1.45
6	D	1555	COM	O1S-S2	2.09	1.51	1.45
5	A	1554	TP7	OXT-C	2.09	1.28	1.22
4	D	1553	F43	C2A-C3A	2.05	1.58	1.54
4	A	1553	F43	C2A-C3A	2.02	1.58	1.54

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1553	F43	CAB-C3B-C2B	-5.65	107.12	119.09
4	D	1553	F43	CAB-C3B-C2B	-5.65	107.14	119.09
4	A	1553	F43	C3B-C4B-CHC	-5.55	111.34	123.32
4	D	1553	F43	C3B-C4B-CHC	-5.54	111.36	123.32
4	A	1553	F43	C4D-ND-C1D	-5.43	101.37	108.51
4	D	1553	F43	C4D-ND-C1D	-5.41	101.40	108.51
4	A	1553	F43	C3D-C4D-ND	5.01	110.13	102.34
4	D	1553	F43	C3D-C4D-ND	5.00	110.12	102.34
4	A	1553	F43	C6D-C7D-CHD	3.97	124.42	116.95
4	D	1553	F43	C6D-C7D-CHD	3.96	124.40	116.95
4	A	1553	F43	O8D-C7D-C6D	-3.94	114.40	120.86
4	D	1553	F43	O8D-C7D-C6D	-3.94	114.41	120.86
5	A	1554	TP7	O1-C1-N	-3.43	117.16	122.95
5	D	1554	TP7	O1-C1-N	-3.41	117.19	122.95
4	A	1553	F43	C3A-C4A-NA	3.09	107.00	102.30
4	D	1553	F43	C3A-C4A-NA	3.09	107.00	102.30
4	D	1553	F43	ODB-CCB-CBB	-2.92	113.70	123.08
4	A	1553	F43	ODB-CCB-CBB	-2.92	113.71	123.08
4	D	1553	F43	C2B-C1B-NB	2.90	106.18	101.84
4	A	1553	F43	C2B-C1B-NB	2.88	106.14	101.84
4	A	1553	F43	OEB-CCB-CBB	2.80	123.03	114.03
4	D	1553	F43	OEB-CCB-CBB	2.80	123.01	114.03
5	A	1554	TP7	C4-C3-C2	2.54	122.34	113.19
5	D	1554	TP7	C4-C3-C2	2.54	122.31	113.19
5	D	1554	TP7	C5-C6-C7	-2.52	108.59	113.09
5	A	1554	TP7	C5-C6-C7	-2.51	108.62	113.09
6	D	1555	COM	O3S-S2-C2	2.47	109.77	105.77
4	D	1553	F43	C2D-C1D-CHD	-2.46	118.70	121.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1553	F43	C2D-C1D-CHD	-2.45	118.71	121.85
6	A	1555	COM	O3S-S2-C2	2.45	109.73	105.77
4	D	1553	F43	C5C-C2C-C3C	-2.32	109.00	114.94
4	A	1553	F43	C5C-C2C-C3C	-2.32	109.00	114.94
4	A	1553	F43	C2B-C3B-C4B	-2.24	99.10	101.63
4	D	1553	F43	C2B-C3B-C4B	-2.21	99.13	101.63
4	A	1553	F43	O7B-C6B-N5B	-2.11	122.39	125.13
4	D	1553	F43	O7B-C6B-N5B	-2.10	122.41	125.13
4	D	1553	F43	C1B-CHB-C4A	2.05	114.26	109.54
4	A	1553	F43	C1B-CHB-C4A	2.05	114.25	109.54

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1553	F43	C3A-CAA-CBA-CCA
4	D	1553	F43	C3A-CAA-CBA-CCA
4	A	1553	F43	C2C-C5C-C6C-O8C
4	D	1553	F43	C2C-C5C-C6C-O8C
4	A	1553	F43	C2C-C5C-C6C-O7C
4	D	1553	F43	C2C-C5C-C6C-O7C
4	A	1553	F43	CAA-CBA-CCA-OEA
4	D	1553	F43	CAA-CBA-CCA-OEA
4	A	1553	F43	C8C-C9C-CAC-OBC
4	D	1553	F43	C8C-C9C-CAC-OBC
4	A	1553	F43	CAA-CBA-CCA-ODA
4	D	1553	F43	CAA-CBA-CCA-ODA
4	A	1553	F43	C8C-C9C-CAC-OCC
4	D	1553	F43	C8C-C9C-CAC-OCC
4	A	1553	F43	CAB-CBB-CCB-OEB
4	D	1553	F43	CAB-CBB-CCB-OEB
4	D	1553	F43	CAB-CBB-CCB-ODB
4	A	1553	F43	CAB-CBB-CCB-ODB
4	A	1553	F43	C3D-C9D-CAD-OB
4	D	1553	F43	C3D-C9D-CAD-OB

There are no ring outliers.

6 monomers are involved in 15 short contacts:

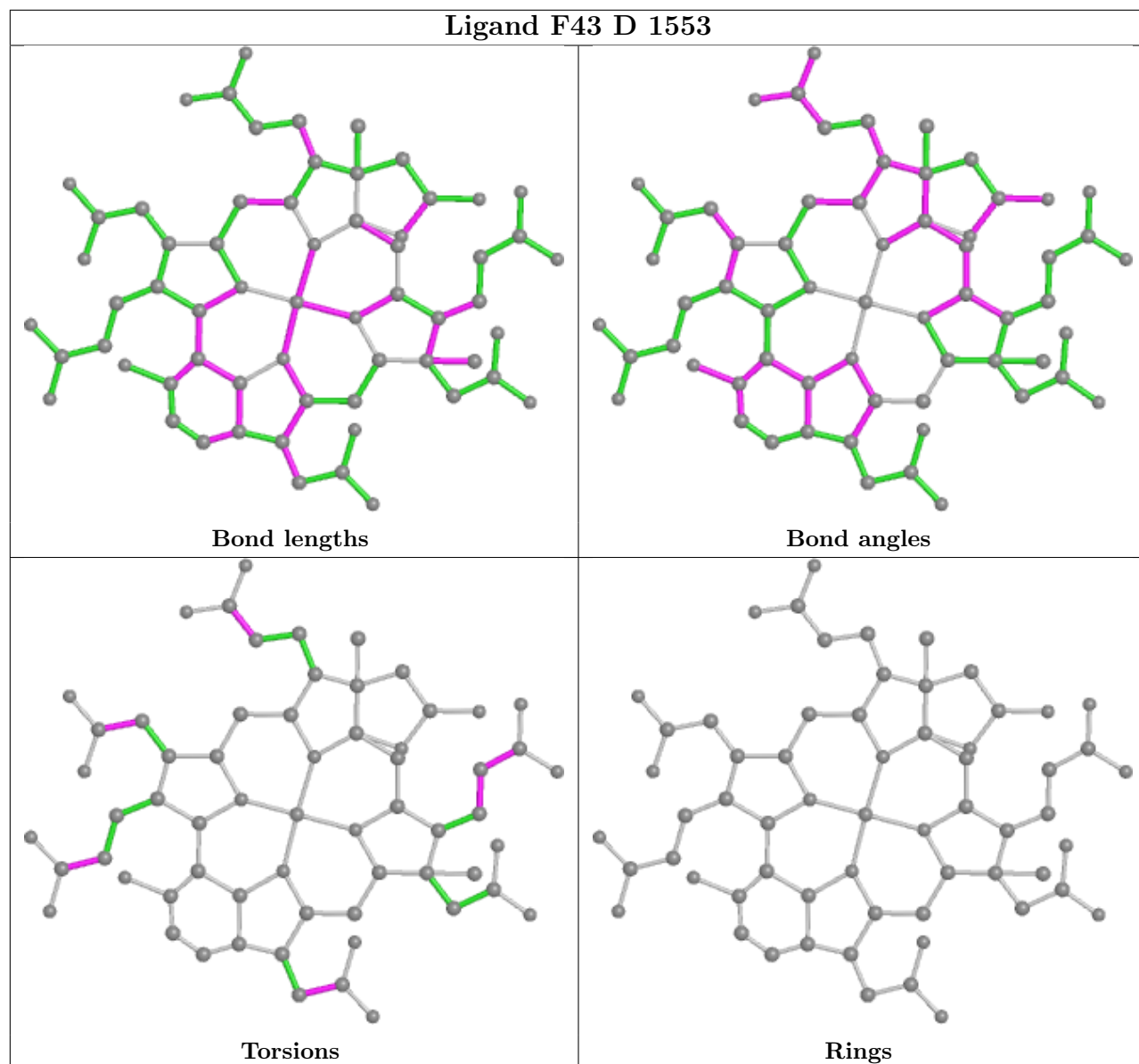
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1553	F43	6	0

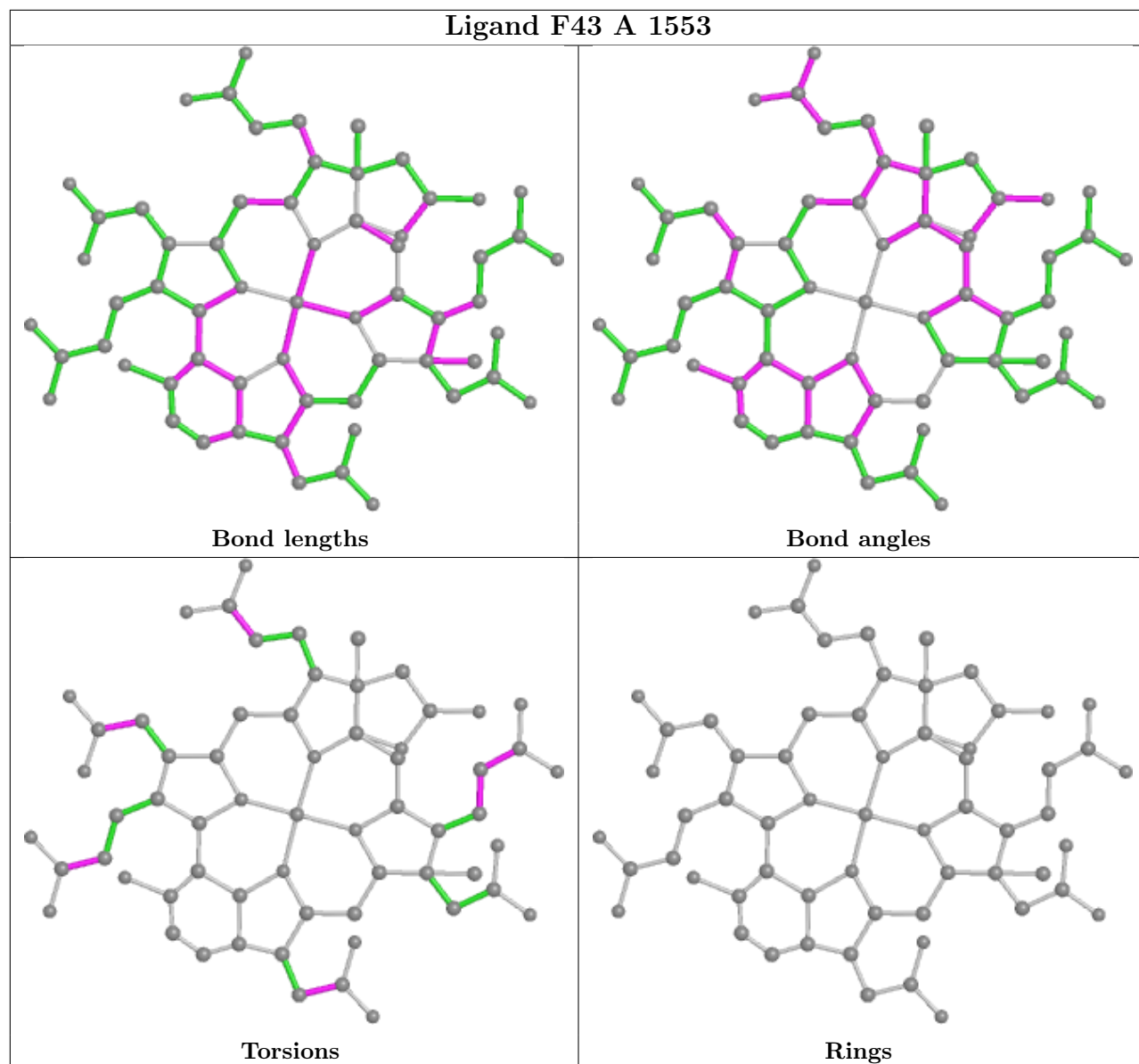
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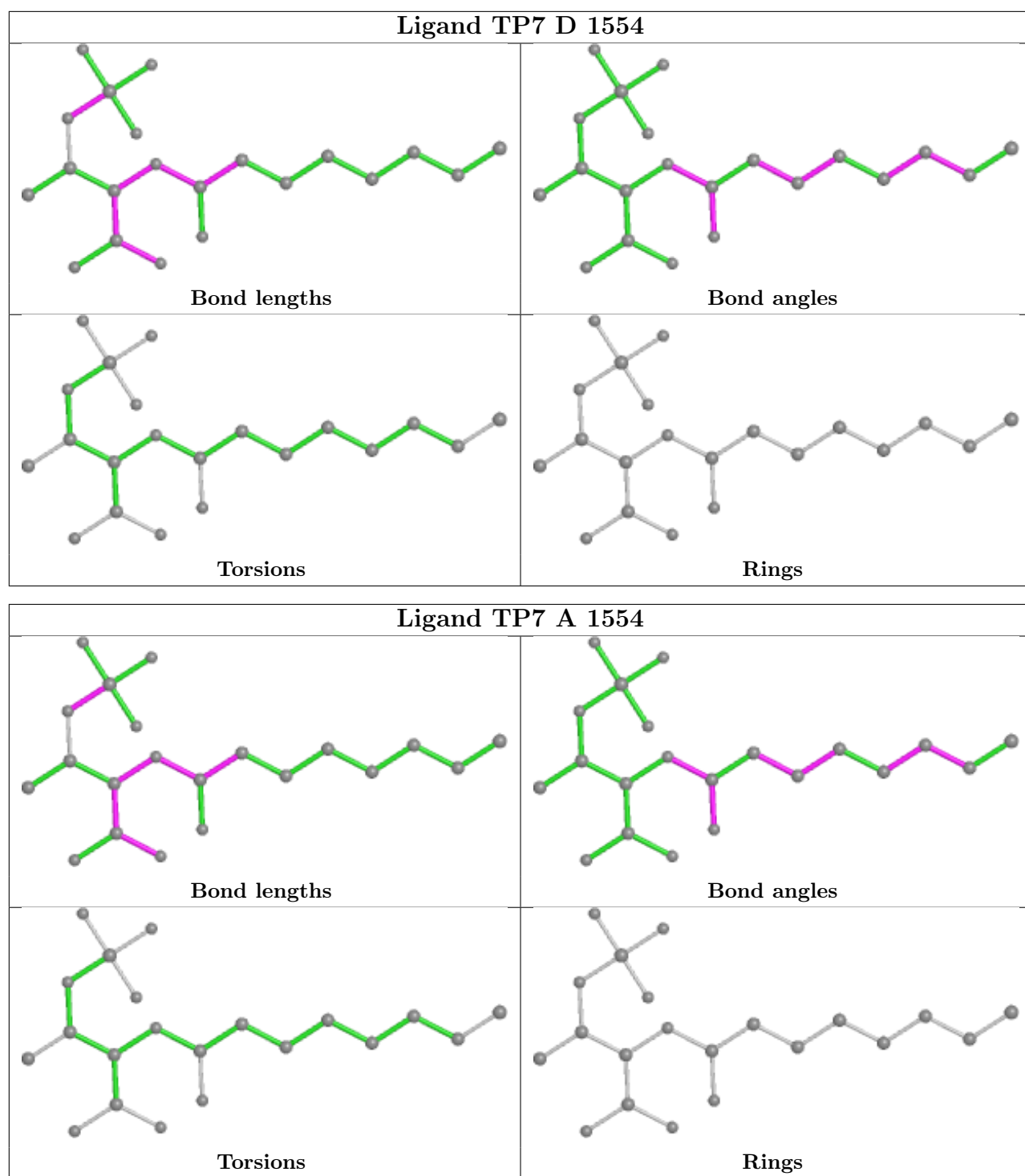
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1553	F43	5	0
5	D	1554	TP7	1	0
5	A	1554	TP7	1	0
6	D	1555	COM	1	0
6	A	1555	COM	1	0

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







4.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

4.8 Polymer linkage issues

There are no chain breaks in this entry.

5 Fit of model and data

5.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	545/553 (98%)	-0.35	2 (0%) 92 93	14, 21, 32, 36	7 (1%)
1	D	545/553 (98%)	-0.34	3 (0%) 89 91	14, 21, 32, 36	7 (1%)
2	B	436/443 (98%)	-0.22	5 (1%) 80 82	15, 25, 35, 42	7 (1%)
2	E	436/443 (98%)	-0.27	1 (0%) 95 96	15, 25, 35, 42	7 (1%)
3	C	248/258 (96%)	0.17	6 (2%) 59 60	20, 32, 40, 46	10 (4%)
3	F	248/258 (96%)	0.06	5 (2%) 65 67	20, 32, 40, 46	10 (4%)
All	All	2458/2508 (98%)	-0.22	22 (0%) 84 85	14, 25, 37, 46	48 (1%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	64	GLU	4.0
1	A	552	ALA	3.9
3	C	67	GLU	3.6
1	D	552	ALA	3.6
3	F	254	GLU	3.5
3	F	64	GLU	3.1
3	C	254	GLU	2.9
3	F	180	ARG	2.7
3	C	201	GLU	2.6
3	C	66	PRO	2.5
3	F	58	PRO	2.5
2	E	442	LYS	2.5
2	B	56	ASN	2.5
3	F	29	LEU	2.4
1	D	367	GLU	2.4
1	D	30	ILE	2.4
2	B	109	ASP	2.3
2	B	21	GLU	2.3
1	A	177	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	20	ALA	2.3
3	C	185	THR	2.2
2	B	76	ASP	2.1

5.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.4 Ligands [i](#)

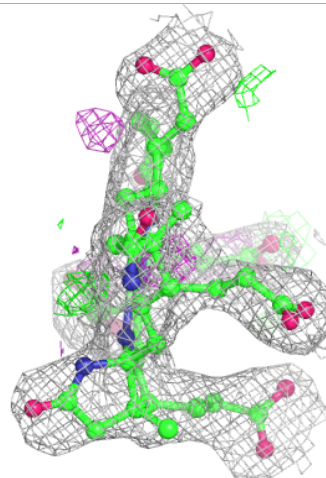
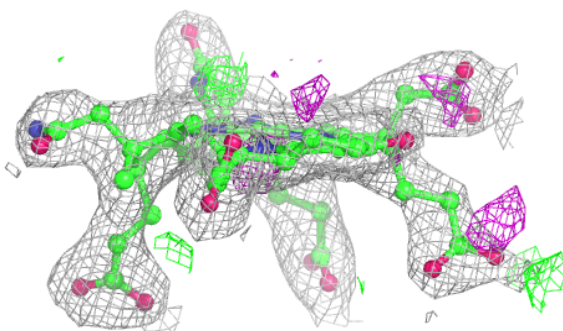
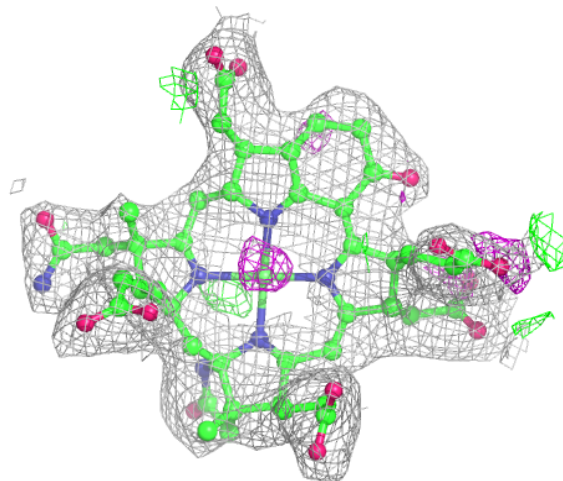
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	F43	A	1553	62/62	0.94	0.17	24,26,26,27	0
4	F43	D	1553	62/62	0.94	0.16	24,26,26,27	0
6	COM	A	1555	7/7	0.94	0.19	16,17,17,17	0
5	TP7	D	1554	21/21	0.96	0.14	23,23,24,24	0
5	TP7	A	1554	21/21	0.96	0.15	23,23,24,24	0
6	COM	D	1555	7/7	0.97	0.12	16,17,17,17	0

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

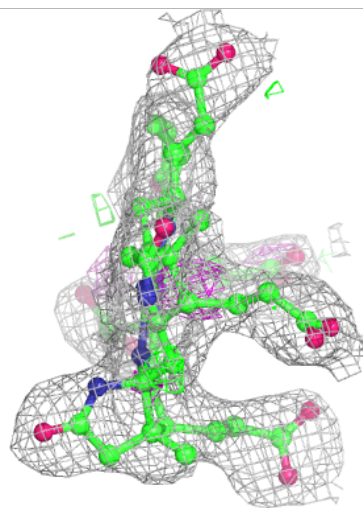
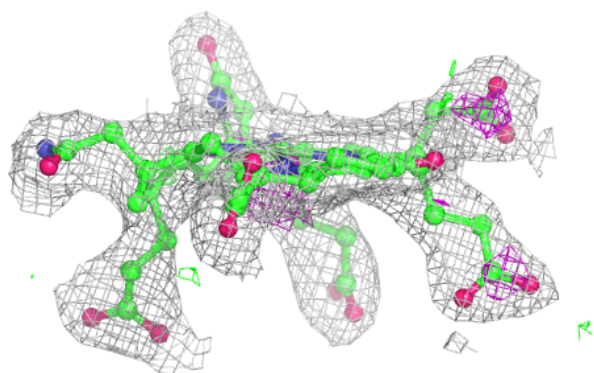
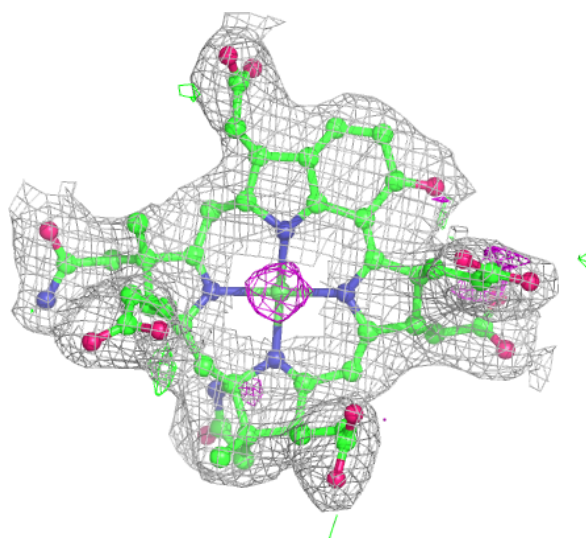
Electron density around F43 A 1553:

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



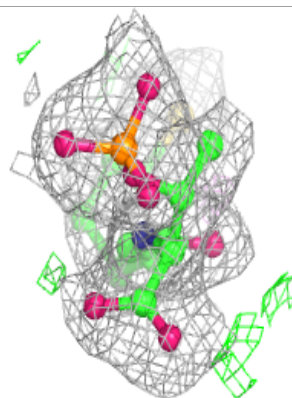
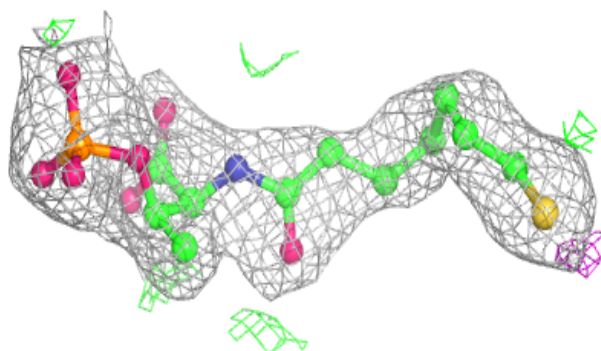
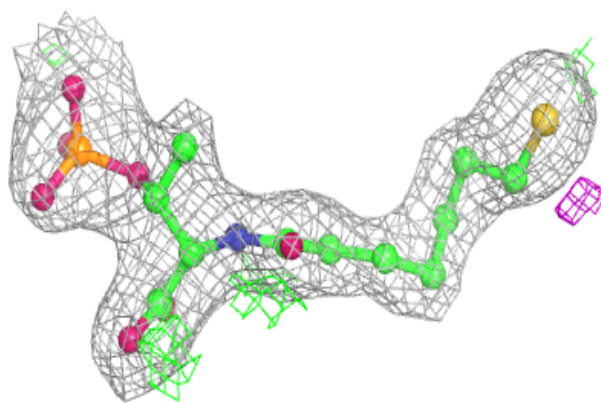
Electron density around F43 D 1553:

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

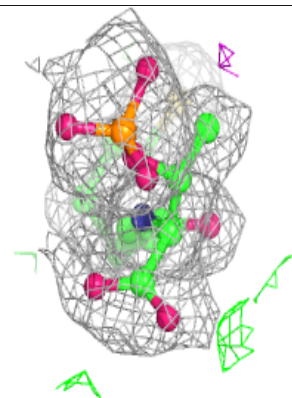
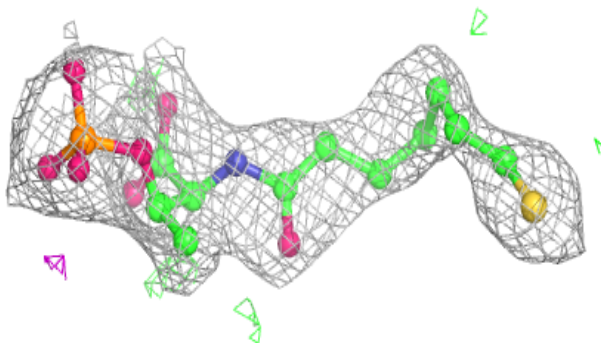
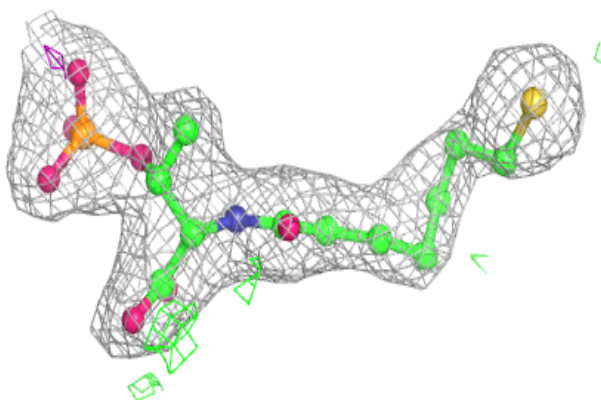


Electron density around TP7 D 1554:

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

**Electron density around TP7 A 1554:**

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



5.5 Other polymers [i](#)

There are no such residues in this entry.