Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 07:55 PM EDT

PDB ID	:	4PDK
Title	:	FadR, Fatty Acid Responsive Transcription Factor from Vibrio cholerae, in Complex with oleoyl-CoA
Authors	:	Shi, W.; Kull, F.J.
Deposited on	:	2014-04-19
Resolution	:	2.80 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 (i) 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 (i)) were used in the production of this report:

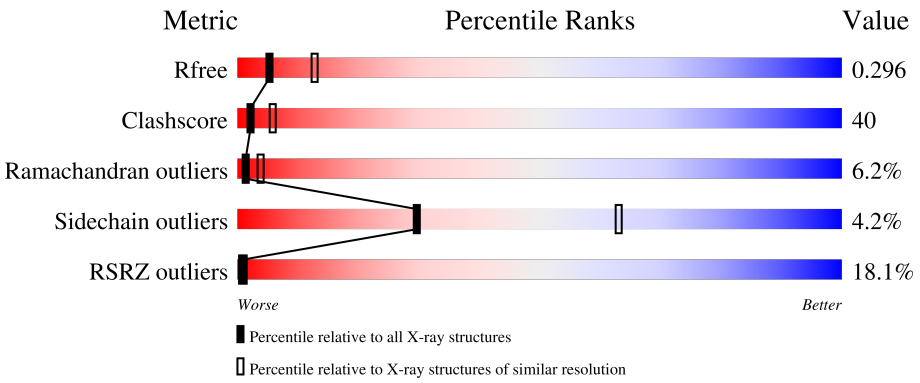
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	279	16% 49% 33% 8% • 7%
1	B	279	18% 51% 35% 6% • 5%

2 Entry composition (i)

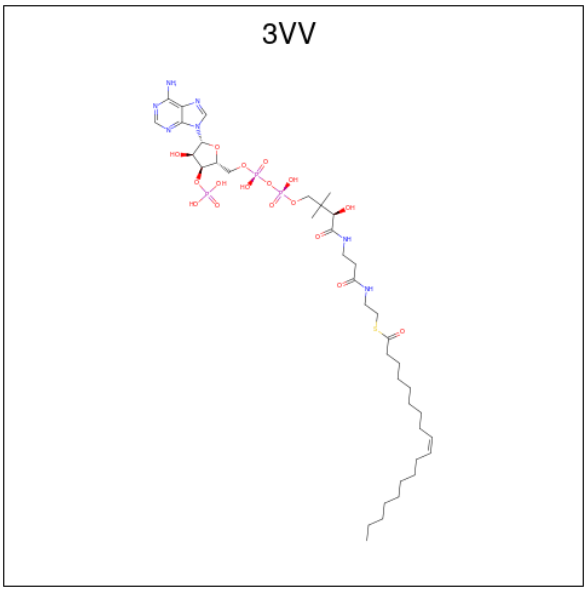
There are 2 unique types of molecules in this entry. The entry contains 4508 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fatty acid metabolism regulator protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	259	2098	1343	361	385	9	0	0	0
1	B	265	2142	1372	368	392	10	0	0	0

- Molecule 2 is S-{(3R,5R,9R)-1-[(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-4-hydroxy-3-(phosphonooxy)tetrahydrofuran-2-yl]-3,5,9-trihydroxy-8,8-dimethyl-3,5-dioxido-10,14-dioxo-2,4,6-trioxa-11,15-diaza-3lambda 5 ,5lambda 5 -diphosphaheptadecan-17-yl} (9Z)-octadec-9-ene thioate (non-preferred name) (three-letter code: 3VV) (formula: $C_{39}H_{68}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	A	1	67	39	7	17	3	1	0	0
2	A	1	67	39	7	17	3	1	0	0
2	B	1	67	39	7	17	3	1	0	0

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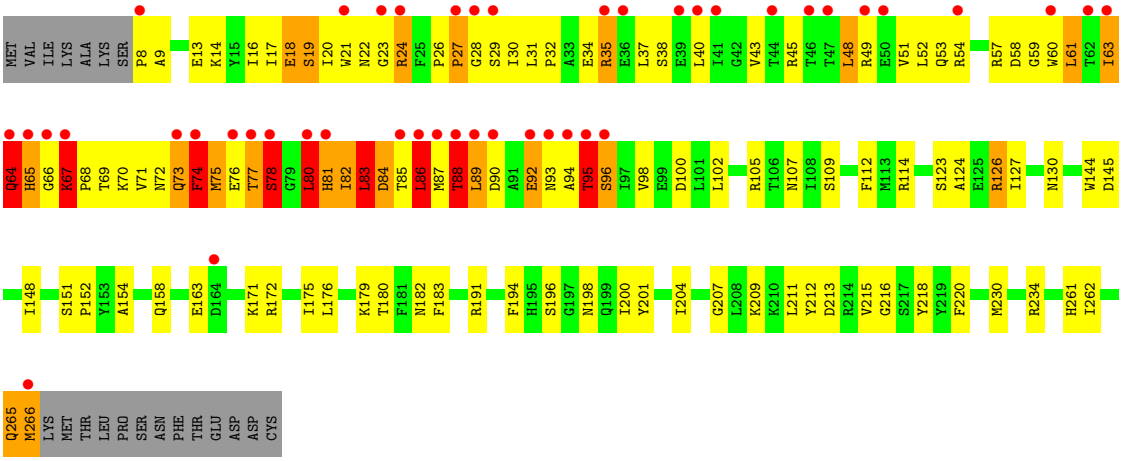
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	B	1	67	39	7	17	3	1	0	0

3 Residue-property plots (i)

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

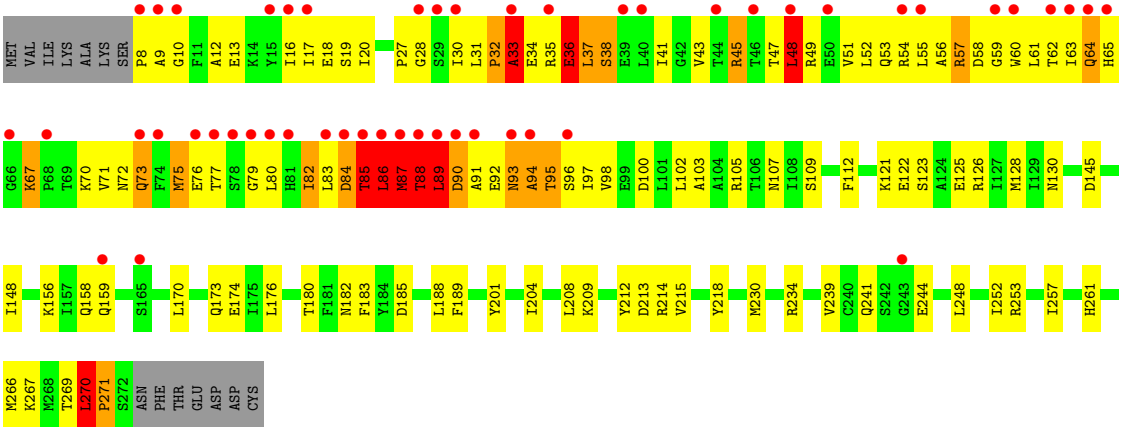
• Molecule 1: Fatty acid metabolism regulator protein

Chain A:



• Molecule 1: Fatty acid metabolism regulator protein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	116.92Å 88.69Å 62.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.73 – 2.80 19.73 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.73-2.80) 99.7 (19.73-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.26 (at 2.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.259 , 0.298 0.258 , 0.296	Depositor DCC
R_{free} test set	1658 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å^2)	34.9	Xtriage
Anisotropy	1.120	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 58.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	4508	wwPDB-VP
Average B, all atoms (Å^2)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.99% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 3VV

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	0.49	0/2145	0.99	14/2899 (0.5%)
1	B	0.50	0/2190	1.03	18/2960 (0.6%)
All	All	0.50	0/4335	1.01	32/5859 (0.5%)

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	8
1	B	0	11
All	All	0	19

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	80	LEU	CA-CB-CG	8.63	135.15	115.30
1	B	48	LEU	CB-CG-CD2	-8.40	96.72	111.00
1	A	266	MET	CG-SD-CE	-8.06	87.31	100.20
1	B	86	LEU	CA-CB-CG	-7.82	97.32	115.30
1	A	81	HIS	N-CA-C	-7.58	90.52	111.00
1	B	85	THR	N-CA-C	7.27	130.62	111.00
1	A	89	LEU	CA-CB-CG	7.20	131.85	115.30
1	B	87	MET	C-N-CA	-7.19	103.72	121.70
1	A	126	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	A	64	GLN	N-CA-C	6.73	129.17	111.00
1	B	87	MET	CA-CB-CG	6.47	124.31	113.30
1	A	67	LYS	N-CA-C	6.34	128.13	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	126	ARG	NE-CZ-NH2	-6.24	117.18	120.30
1	B	90	ASP	CB-CG-OD1	-6.05	112.85	118.30
1	B	33	ALA	C-N-CA	-6.05	106.58	121.70
1	A	61	LEU	CB-CG-CD2	6.01	121.22	111.00
1	B	86	LEU	N-CA-C	5.83	126.75	111.00
1	A	61	LEU	CA-CB-CG	5.83	128.70	115.30
1	B	67	LYS	C-N-CD	5.81	140.60	128.40
1	A	64	GLN	N-CA-CB	-5.81	100.14	110.60
1	B	48	LEU	CB-CG-CD1	5.71	120.71	111.00
1	B	87	MET	CB-CG-SD	5.56	129.09	112.40
1	B	36	GLU	N-CA-C	5.56	126.02	111.00
1	B	86	LEU	CB-CG-CD1	5.49	120.33	111.00
1	A	126	ARG	CD-NE-CZ	5.48	131.27	123.60
1	B	85	THR	CB-CA-C	-5.34	97.18	111.60
1	B	87	MET	N-CA-C	-5.31	96.66	111.00
1	A	86	LEU	CA-CB-CG	-5.26	103.21	115.30
1	B	90	ASP	CB-CG-OD2	5.23	123.01	118.30
1	A	48	LEU	CB-CA-C	-5.17	100.39	110.20
1	B	37	LEU	CB-CG-CD1	5.16	119.78	111.00
1	B	88	THR	N-CA-CB	-5.00	100.80	110.30

There are no chirality outliers.

All (19) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	19	SER	Peptide
1	A	63	ILE	Peptide
1	A	67	LYS	Peptide
1	A	74	PHE	Peptide
1	A	83	LEU	Peptide
1	A	86	LEU	Peptide
1	A	88	THR	Peptide
1	A	95	THR	Peptide
1	B	121	LYS	Peptide
1	B	270	LEU	Peptide
1	B	35	ARG	Peptide
1	B	57	ARG	Peptide
1	B	73	GLN	Peptide
1	B	83	LEU	Peptide
1	B	84	ASP	Peptide
1	B	85	THR	Peptide
1	B	87	MET	Peptide

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Mol	Chain	Res	Type	Group
1	B	88	THR	Peptide
1	B	89	LEU	Peptide

5.2 Too-close contacts (i)

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	2098	0	2084	175	0
1	B	2142	0	2133	177	0
2	A	134	0	128	16	0
2	B	134	0	128	17	0
All	All	4508	0	4473	355	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 40.

All (355) 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:A:14:LYS:HB2	1:A:92:GLU:HB2	1.18	1.13
1:B:93:ASN:HD21	1:B:97:ILE:HD13	1.04	1.08
1:A:45:ARG:HD2	1:A:48:LEU:HB2	1.32	1.06
1:A:29:SER:O	1:A:70:LYS:NZ	1.87	1.05
1:B:34:GLU:O	1:B:45:ARG:NH1	1.90	1.05
1:A:43:VAL:H	1:A:45:ARG:NH1	1.58	1.01
1:B:49:ARG:HG2	1:B:52:LEU:HB2	1.41	1.00
1:A:84:ASP:H	1:A:85:THR:HA	1.28	0.97
1:B:37:LEU:HB3	1:B:45:ARG:HH22	1.30	0.95
1:B:93:ASN:ND2	1:B:97:ILE:HD13	1.80	0.95
1:A:54:ARG:NH1	1:A:58:ASP:OD2	2.00	0.94
1:A:262:ILE:HA	1:A:266:MET:HG2	1.50	0.94
1:B:86:LEU:HD12	1:B:87:MET:N	1.83	0.93
1:A:43:VAL:O	1:A:45:ARG:NH1	2.01	0.92
1:A:64:GLN:HG3	1:A:67:LYS:HB2	1.50	0.92
1:B:84:ASP:H	1:B:86:LEU:HD22	1.33	0.92
1:B:38:SER:OG	1:B:45:ARG:NH1	2.04	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:THR:OG1	1:B:96:SER:N	2.02	0.90
1:A:60:TRP:HE1	1:A:82:ILE:HG22	1.34	0.90
1:A:75:MET:O	1:A:77:THR:N	2.06	0.89
1:A:43:VAL:H	1:A:45:ARG:HH12	1.17	0.89
1:B:86:LEU:O	1:B:214:ARG:NH1	2.06	0.88
1:A:70:LYS:HZ2	1:A:74:PHE:HE2	1.22	0.88
1:B:89:LEU:HG	1:B:90:ASP:H	1.38	0.87
1:A:27:PRO:HG2	2:B:302:3VV:H10	1.57	0.87
1:B:148:ILE:HG12	1:B:158:GLN:HE21	1.38	0.86
1:A:65:HIS:CD2	1:A:66:GLY:H	1.93	0.86
1:A:64:GLN:HG2	1:A:65:HIS:N	1.90	0.85
1:B:84:ASP:H	1:B:86:LEU:CD2	1.89	0.85
1:B:269:THR:HB	1:B:270:LEU:HD13	1.59	0.85
1:A:59:GLY:O	1:A:72:ASN:ND2	2.10	0.84
1:A:95:THR:OG1	1:A:96:SER:N	2.07	0.83
1:A:70:LYS:NZ	1:A:74:PHE:HE2	1.77	0.82
1:B:84:ASP:N	1:B:86:LEU:HD22	1.94	0.81
1:B:86:LEU:HG	1:B:87:MET:HG2	1.63	0.81
1:B:93:ASN:O	1:B:95:THR:N	2.13	0.80
1:A:191:ARG:HH21	2:A:302:3VV:H57	1.46	0.80
1:A:89:LEU:HD12	1:A:90:ASP:H	1.47	0.80
1:A:93:ASN:HB3	1:A:95:THR:HG22	1.64	0.79
1:A:90:ASP:OD2	1:A:215:VAL:HG22	1.82	0.79
1:B:88:THR:OG1	1:B:90:ASP:HA	1.82	0.79
1:A:61:LEU:HA	1:A:71:VAL:HA	1.64	0.78
1:B:12:ALA:HA	1:B:41:ILE:HD11	1.66	0.78
1:B:56:ALA:HA	1:B:61:LEU:HB2	1.66	0.78
1:B:84:ASP:HB2	1:B:85:THR:OG1	1.84	0.77
1:B:93:ASN:HD21	1:B:97:ILE:CD1	1.92	0.77
1:B:96:SER:O	1:B:100:ASP:N	2.16	0.77
1:B:57:ARG:HA	1:B:59:GLY:H	1.49	0.77
1:A:14:LYS:HB2	1:A:92:GLU:CB	2.10	0.76
1:A:163:GLU:HB2	1:A:179:LYS:NZ	2.01	0.76
1:B:60:TRP:NE1	1:B:84:ASP:OD2	2.18	0.76
1:A:38:SER:HB3	1:A:45:ARG:CG	2.16	0.75
1:A:81:HIS:O	1:A:84:ASP:HB3	1.87	0.74
1:A:212:TYR:HA	2:A:301:3VV:H34	1.69	0.74
1:A:86:LEU:HG	1:A:92:GLU:OE2	1.87	0.73
1:A:92:GLU:HA	1:A:93:ASN:HB2	1.69	0.73
1:B:38:SER:HA	1:B:41:ILE:HG22	1.71	0.72
1:A:49:ARG:O	1:A:53:GLN:HG3	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:GLY:HA2	1:A:26:PRO:HA	1.73	0.71
1:B:34:GLU:HG2	1:B:49:ARG:HD3	1.73	0.70
1:B:79:GLY:HA2	1:B:82:ILE:HD11	1.72	0.70
1:B:103:ALA:O	1:B:107:ASN:ND2	2.24	0.70
1:A:37:LEU:O	1:A:40:LEU:HB3	1.92	0.70
1:B:85:THR:OG1	1:B:86:LEU:HD13	1.92	0.69
1:A:83:LEU:HD11	1:A:87:MET:HG2	1.74	0.69
1:B:90:ASP:CG	1:B:94:ALA:HB3	2.12	0.69
1:A:30:ILE:HG13	1:A:70:LYS:HD2	1.72	0.69
1:A:64:GLN:HG3	1:A:67:LYS:HD3	1.74	0.69
1:B:73:GLN:HB3	1:B:75:MET:HA	1.73	0.68
1:A:65:HIS:CG	1:A:66:GLY:H	2.07	0.68
1:B:62:THR:HG22	1:B:70:LYS:H	1.58	0.68
1:B:212:TYR:HA	2:B:301:3VV:H34	1.76	0.68
1:B:266:MET:C	1:B:269:THR:O	2.32	0.68
1:B:88:THR:HG23	1:B:89:LEU:H	1.59	0.68
1:B:156:LYS:HA	1:B:159:GLN:HG2	1.74	0.67
1:A:18:GLU:HB2	1:A:95:THR:HB	1.75	0.67
1:A:96:SER:O	1:A:100:ASP:N	2.24	0.67
1:A:71:VAL:CG2	1:A:74:PHE:HZ	2.08	0.67
1:A:218:TYR:OH	1:A:265:GLN:NE2	2.27	0.67
1:A:93:ASN:CG	1:A:95:THR:H	1.99	0.66
1:A:21:TRP:CE3	1:A:22:ASN:HB3	2.31	0.66
1:A:75:MET:C	1:A:77:THR:H	1.98	0.66
1:A:86:LEU:C	1:A:88:THR:H	1.99	0.66
1:A:90:ASP:CG	1:A:98:VAL:HG21	2.15	0.66
1:B:105:ARG:NH1	1:B:189:PHE:CZ	2.64	0.66
1:A:38:SER:HB3	1:A:45:ARG:HG2	1.77	0.65
1:B:30:ILE:HD12	1:B:70:LYS:HG2	1.78	0.65
1:A:38:SER:HB3	1:A:45:ARG:HG3	1.79	0.65
1:A:29:SER:C	1:A:70:LYS:HZ3	1.95	0.65
1:A:114:ARG:HB3	2:B:301:3VV:H68	1.78	0.64
1:B:92:GLU:CG	1:B:93:ASN:H	2.11	0.64
1:B:182:ASN:HD21	1:B:213:ASP:HA	1.62	0.64
1:A:93:ASN:OD1	1:A:94:ALA:N	2.29	0.64
1:A:43:VAL:N	1:A:45:ARG:NH1	2.40	0.64
1:A:60:TRP:NE1	1:A:82:ILE:HG22	2.10	0.64
1:A:105:ARG:HB2	2:A:301:3VV:H28	1.80	0.64
1:B:93:ASN:OD1	1:B:95:THR:HG23	1.97	0.64
1:A:71:VAL:HG22	1:A:74:PHE:HZ	1.64	0.63
1:B:86:LEU:HD12	1:B:87:MET:H	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:THR:HG21	1:A:92:GLU:OE1	1.98	0.63
1:A:92:GLU:O	1:A:92:GLU:HG2	1.97	0.63
1:B:257:ILE:HG13	2:B:301:3VV:H49	1.80	0.63
1:B:94:ALA:HB1	1:B:98:VAL:HB	1.82	0.62
1:B:230:MET:O	1:B:234:ARG:HG3	2.00	0.62
1:B:45:ARG:HA	1:B:48:LEU:HB3	1.81	0.61
1:B:58:ASP:CG	1:B:85:THR:HG23	2.20	0.61
1:A:64:GLN:CG	1:A:67:LYS:HB2	2.28	0.61
1:B:10:GLY:O	1:B:13:GLU:HB3	2.00	0.61
1:B:93:ASN:OD1	1:B:94:ALA:N	2.34	0.61
1:A:64:GLN:HG3	1:A:67:LYS:CB	2.29	0.60
1:A:84:ASP:H	1:A:85:THR:CA	2.09	0.60
2:A:302:3VV:H17	2:B:302:3VV:H27	1.82	0.60
1:B:16:ILE:HD11	1:B:55:LEU:HD13	1.83	0.60
1:B:56:ALA:HB1	1:B:63:ILE:HG22	1.82	0.60
1:B:63:ILE:O	1:B:64:GLN:HB3	2.01	0.60
1:A:18:GLU:O	1:A:21:TRP:HB3	2.01	0.60
1:B:38:SER:N	1:B:45:ARG:HH12	1.99	0.60
1:A:88:THR:OG1	1:A:89:LEU:N	2.34	0.60
1:B:34:GLU:HB3	1:B:45:ARG:CZ	2.32	0.59
1:B:94:ALA:CB	1:B:98:VAL:HB	2.33	0.59
1:B:257:ILE:O	1:B:261:HIS:ND1	2.18	0.59
1:B:38:SER:H	1:B:45:ARG:HH12	1.51	0.59
1:B:54:ARG:O	1:B:58:ASP:OD2	2.20	0.59
1:B:8:PRO:HD2	1:B:9:ALA:H	1.68	0.59
1:B:28:GLY:H	1:B:71:VAL:HG13	1.68	0.59
1:B:90:ASP:OD1	1:B:94:ALA:HB3	2.02	0.59
1:B:253:ARG:NH2	2:B:301:3VV:O40	2.36	0.58
1:B:38:SER:HG	1:B:45:ARG:HH11	1.51	0.58
1:B:18:GLU:CG	1:B:95:THR:HB	2.33	0.58
1:A:14:LYS:HG3	1:A:93:ASN:HB2	1.84	0.58
1:B:257:ILE:HG23	1:B:261:HIS:CE1	2.38	0.58
1:A:130:ASN:OD1	2:A:302:3VV:N63	2.37	0.58
1:B:56:ALA:O	1:B:57:ARG:HG2	2.03	0.58
1:B:88:THR:HG23	1:B:89:LEU:N	2.18	0.57
1:B:18:GLU:HG2	1:B:95:THR:HB	1.84	0.57
1:A:154:ALA:O	1:A:158:GLN:HG3	2.04	0.57
1:A:43:VAL:N	1:A:45:ARG:HH12	1.95	0.57
1:A:58:ASP:OD2	1:A:86:LEU:HD22	2.04	0.57
1:A:90:ASP:OD1	1:A:98:VAL:HG21	2.06	0.56
1:B:92:GLU:HG3	1:B:93:ASN:H	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:ARG:NH1	1:A:58:ASP:CG	2.58	0.56
1:A:23:GLY:HA2	1:A:26:PRO:CA	2.36	0.56
1:B:148:ILE:HG12	1:B:158:GLN:NE2	2.15	0.56
1:A:200:ILE:HG23	2:A:302:3VV:H29	1.88	0.55
1:B:123:SER:HB3	1:B:126:ARG:NH2	2.21	0.55
1:B:32:PRO:HB3	1:B:36:GLU:HG2	1.88	0.55
1:B:73:GLN:NE2	1:B:75:MET:O	2.39	0.55
1:B:87:MET:CA	1:B:88:THR:HB	2.37	0.55
1:A:17:ILE:HG12	1:A:60:TRP:CZ3	2.41	0.54
1:B:51:VAL:HG12	1:B:54:ARG:HD3	1.89	0.54
1:B:58:ASP:CB	1:B:85:THR:HG23	2.37	0.54
1:A:65:HIS:CG	1:A:66:GLY:N	2.75	0.54
1:B:17:ILE:HG21	1:B:93:ASN:HB2	1.88	0.54
1:A:262:ILE:HA	1:A:266:MET:CG	2.32	0.54
1:A:171:LYS:O	1:A:175:ILE:HG22	2.07	0.54
1:B:253:ARG:HH21	2:B:301:3VV:H53	1.72	0.54
1:A:17:ILE:HG23	1:A:60:TRP:HZ3	1.73	0.54
1:A:261:HIS:HB3	1:A:266:MET:CE	2.38	0.54
1:A:163:GLU:HG2	1:A:175:ILE:HD13	1.88	0.53
1:A:13:GLU:OE2	1:A:86:LEU:HD21	2.07	0.53
1:A:71:VAL:HG23	1:A:74:PHE:CZ	2.44	0.53
1:B:77:THR:O	1:B:77:THR:OG1	2.16	0.53
1:B:130:ASN:OD1	2:B:302:3VV:N63	2.40	0.53
1:A:16:ILE:O	1:A:20:ILE:HG13	2.08	0.53
1:A:17:ILE:HG23	1:A:60:TRP:CZ3	2.44	0.53
1:B:38:SER:N	1:B:45:ARG:NH1	2.57	0.53
1:A:38:SER:HA	1:A:45:ARG:CZ	2.40	0.52
1:A:87:MET:HB3	1:A:211:LEU:HD13	1.90	0.52
1:A:29:SER:C	1:A:70:LYS:NZ	2.58	0.52
1:B:180:THR:O	1:B:183:PHE:HB3	2.09	0.52
1:A:29:SER:N	1:A:70:LYS:HZ1	2.08	0.52
1:A:60:TRP:O	1:A:71:VAL:HB	2.10	0.52
1:B:86:LEU:HD11	1:B:87:MET:HE2	1.91	0.52
1:B:84:ASP:N	1:B:86:LEU:CD2	2.64	0.51
1:B:185:ASP:OD1	1:B:212:TYR:OH	2.19	0.51
1:B:62:THR:CG2	1:B:70:LYS:H	2.22	0.51
1:B:94:ALA:O	1:B:96:SER:N	2.44	0.51
1:A:73:GLN:O	1:A:74:PHE:CG	2.64	0.51
1:B:17:ILE:CD1	1:B:60:TRP:HZ3	2.24	0.51
1:B:79:GLY:CA	1:B:82:ILE:HD11	2.39	0.51
1:A:204:ILE:HD11	2:B:301:3VV:H14	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:SER:HA	1:B:112:PHE:CE2	2.45	0.51
1:B:8:PRO:HB2	1:B:47:THR:HG21	1.93	0.51
1:B:67:LYS:HD3	1:B:67:LYS:N	2.26	0.51
1:B:102:LEU:HD21	2:B:301:3VV:H3	1.93	0.51
1:A:67:LYS:HG2	1:A:69:THR:HA	1.92	0.50
1:A:34:GLU:HG2	1:A:49:ARG:HD3	1.91	0.50
1:B:87:MET:HB3	1:B:88:THR:HG22	1.92	0.50
1:B:13:GLU:O	1:B:17:ILE:HG13	2.12	0.50
1:A:32:PRO:HB2	1:A:37:LEU:CD1	2.42	0.50
1:B:38:SER:H	1:B:45:ARG:NH1	2.09	0.50
1:A:60:TRP:O	1:A:61:LEU:HD23	2.12	0.50
1:B:63:ILE:HD12	1:B:64:GLN:HB2	1.94	0.50
1:B:239:VAL:HG11	1:B:248:LEU:HD13	1.94	0.50
1:A:123:SER:O	1:A:126:ARG:HB3	2.12	0.50
1:B:32:PRO:C	1:B:33:ALA:O	2.50	0.49
1:A:21:TRP:HE3	1:A:22:ASN:HB3	1.78	0.49
1:B:57:ARG:CA	1:B:59:GLY:H	2.22	0.49
1:B:267:LYS:N	1:B:269:THR:O	2.44	0.49
1:A:52:LEU:O	1:A:61:LEU:HD12	2.13	0.49
1:A:84:ASP:N	1:A:85:THR:HA	2.08	0.49
1:A:191:ARG:NH2	2:A:302:3VV:H57	2.23	0.49
1:A:71:VAL:CG2	1:A:74:PHE:CZ	2.92	0.49
1:B:89:LEU:HD12	1:B:218:TYR:CD1	2.48	0.49
1:A:17:ILE:HD12	1:A:92:GLU:HG2	1.95	0.49
1:B:87:MET:CA	1:B:92:GLU:HG2	2.43	0.49
1:A:26:PRO:O	1:A:28:GLY:N	2.46	0.48
1:B:92:GLU:CG	1:B:93:ASN:N	2.76	0.48
1:B:17:ILE:CG2	1:B:93:ASN:HD22	2.26	0.48
1:B:89:LEU:HD12	1:B:218:TYR:CG	2.48	0.48
1:B:34:GLU:O	1:B:37:LEU:N	2.44	0.48
1:A:45:ARG:HA	1:A:48:LEU:H	1.79	0.48
1:B:32:PRO:HB2	1:B:37:LEU:HD22	1.95	0.48
1:A:28:GLY:N	1:A:74:PHE:CD2	2.81	0.48
1:A:65:HIS:CE1	1:A:67:LYS:HD2	2.48	0.48
1:B:61:LEU:HD23	1:B:71:VAL:HA	1.95	0.48
1:B:86:LEU:HD21	1:B:87:MET:HE3	1.96	0.48
1:A:8:PRO:HD2	1:A:9:ALA:H	1.78	0.48
1:A:163:GLU:HB2	1:A:179:LYS:HZ1	1.77	0.48
1:A:220:PHE:CE2	2:A:301:3VV:H7	2.48	0.48
1:A:84:ASP:N	1:A:85:THR:OG1	2.47	0.48
1:B:60:TRP:HA	1:B:72:ASN:HD22	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:PRO:HG2	1:B:37:LEU:HD13	1.96	0.47
1:B:73:GLN:HB3	1:B:75:MET:CA	2.42	0.47
1:B:32:PRO:HB3	1:B:36:GLU:CG	2.43	0.47
1:B:49:ARG:O	1:B:53:GLN:NE2	2.47	0.47
1:B:270:LEU:HB3	1:B:271:PRO:HD3	1.97	0.47
1:A:194:PHE:CD1	2:A:302:3VV:H42	2.48	0.47
1:A:64:GLN:HG2	1:A:65:HIS:C	2.34	0.47
1:A:84:ASP:N	1:A:85:THR:CA	2.75	0.47
1:B:16:ILE:O	1:B:20:ILE:HG13	2.15	0.47
1:B:204:ILE:HG23	2:B:302:3VV:H17	1.97	0.47
1:B:253:ARG:NH2	2:B:301:3VV:H53	2.30	0.47
1:A:27:PRO:HB3	1:A:74:PHE:CD1	2.50	0.47
1:B:93:ASN:CG	1:B:95:THR:HG23	2.35	0.47
1:A:21:TRP:CZ3	1:A:22:ASN:HB3	2.50	0.47
1:B:31:LEU:HD12	1:B:32:PRO:HD2	1.97	0.47
1:B:86:LEU:O	1:B:87:MET:HB2	2.14	0.47
1:A:261:HIS:HB3	1:A:266:MET:HE1	1.96	0.47
1:B:170:LEU:O	1:B:174:GLU:HG2	2.15	0.47
1:A:92:GLU:N	1:A:93:ASN:HA	2.30	0.46
1:A:17:ILE:HG12	1:A:60:TRP:HZ3	1.80	0.46
1:A:86:LEU:HA	1:A:86:LEU:HD12	1.23	0.46
1:B:182:ASN:ND2	1:B:213:ASP:HA	2.28	0.46
1:A:180:THR:O	1:A:183:PHE:HB3	2.16	0.46
1:B:17:ILE:HD13	1:B:60:TRP:HZ3	1.80	0.46
1:A:102:LEU:HD21	2:A:301:3VV:H3	1.98	0.46
2:B:302:3VV:H31	2:B:302:3VV:H26	1.80	0.46
1:A:163:GLU:HB2	1:A:179:LYS:HZ2	1.75	0.46
1:A:163:GLU:O	1:A:172:ARG:NH1	2.49	0.46
1:A:112:PHE:HB2	1:A:196:SER:HB3	1.97	0.46
1:A:211:LEU:O	1:A:215:VAL:HG23	2.16	0.46
1:A:109:SER:HA	1:A:112:PHE:CE2	2.50	0.45
1:B:60:TRP:CA	1:B:72:ASN:HD22	2.29	0.45
1:B:89:LEU:HG	1:B:90:ASP:N	2.19	0.45
1:A:98:VAL:O	1:A:102:LEU:HG	2.16	0.45
1:A:144:TRP:CE3	1:A:176:LEU:HD13	2.52	0.45
1:B:49:ARG:HA	1:B:52:LEU:H	1.80	0.45
1:A:86:LEU:C	1:A:88:THR:N	2.65	0.45
2:A:302:3VV:H8	1:B:27:PRO:HG2	1.99	0.45
1:B:86:LEU:HD11	1:B:87:MET:CE	2.47	0.45
1:A:58:ASP:O	1:A:84:ASP:OD2	2.34	0.45
1:A:64:GLN:HG3	1:A:67:LYS:CD	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:HIS:O	1:A:81:HIS:ND1	2.50	0.45
1:B:30:ILE:CD1	1:B:70:LYS:HG2	2.46	0.45
1:A:34:GLU:HG2	1:A:49:ARG:HH11	1.80	0.45
1:A:73:GLN:NE2	1:A:78:SER:O	2.50	0.45
1:B:12:ALA:HB3	1:B:51:VAL:HG21	1.98	0.45
1:A:34:GLU:HB3	1:A:48:LEU:HD23	1.98	0.44
1:A:78:SER:C	1:A:80:LEU:H	2.18	0.44
1:A:87:MET:HG3	1:A:211:LEU:HD22	2.00	0.44
2:B:301:3VV:H2	2:B:301:3VV:H29	1.99	0.44
1:A:27:PRO:HD2	2:B:302:3VV:H12	1.99	0.44
1:B:94:ALA:O	1:B:97:ILE:N	2.50	0.44
1:B:145:ASP:O	1:B:148:ILE:HG22	2.18	0.44
1:B:208:LEU:HD11	2:B:302:3VV:H18	2.00	0.44
1:A:64:GLN:NE2	1:A:65:HIS:O	2.50	0.44
1:A:85:THR:HG23	1:A:87:MET:HB2	2.00	0.44
1:A:182:ASN:ND2	1:A:216:GLY:HA3	2.33	0.44
1:A:63:ILE:HG22	1:A:64:GLN:H	1.82	0.44
1:A:124:ALA:HA	1:A:127:ILE:HD12	2.00	0.44
1:A:53:GLN:O	1:A:57:ARG:HG3	2.18	0.43
1:A:102:LEU:HD13	2:A:301:3VV:O25	2.18	0.43
1:A:123:SER:O	1:A:127:ILE:HG13	2.18	0.43
1:B:17:ILE:CD1	1:B:92:GLU:OE2	2.66	0.43
1:A:18:GLU:C	1:A:21:TRP:HB3	2.39	0.43
1:A:88:THR:HG21	1:A:92:GLU:CD	2.38	0.43
2:A:302:3VV:H30	2:A:302:3VV:H33	1.73	0.43
1:B:59:GLY:C	1:B:60:TRP:CD1	2.92	0.43
1:A:80:LEU:CG	1:A:81:HIS:H	2.31	0.43
1:B:125:GLU:OE2	1:B:241:GLN:HA	2.18	0.43
1:A:262:ILE:N	1:A:266:MET:HE2	2.33	0.43
1:B:48:LEU:HD12	1:B:48:LEU:O	2.19	0.43
1:B:95:THR:OG1	1:B:97:ILE:HD12	2.18	0.43
1:B:93:ASN:C	1:B:95:THR:N	2.72	0.43
1:B:16:ILE:HA	1:B:19:SER:HB2	2.01	0.43
1:B:270:LEU:HB3	1:B:271:PRO:CD	2.48	0.43
1:A:70:LYS:NZ	1:A:74:PHE:CE2	2.70	0.43
1:A:77:THR:HA	1:A:78:SER:C	2.39	0.43
1:A:207:GLY:HA3	2:A:302:3VV:H19	2.00	0.43
1:A:51:VAL:HA	1:A:54:ARG:HB3	2.01	0.42
1:A:64:GLN:CG	1:A:67:LYS:HD3	2.47	0.42
1:A:64:GLN:HG2	1:A:65:HIS:CA	2.49	0.42
1:B:84:ASP:C	1:B:86:LEU:HB2	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ARG:O	1:A:38:SER:OG	2.27	0.42
1:A:85:THR:CG2	1:A:87:MET:HB2	2.49	0.42
1:A:107:ASN:ND2	1:B:201:TYR:OH	2.51	0.42
1:B:38:SER:OG	1:B:45:ARG:CG	2.68	0.42
1:B:38:SER:HB3	1:B:43:VAL:O	2.19	0.42
1:B:84:ASP:H	1:B:86:LEU:HB2	1.85	0.42
1:A:64:GLN:HE21	1:A:66:GLY:C	2.23	0.42
1:B:18:GLU:HG3	1:B:95:THR:HB	2.01	0.42
1:A:265:GLN:OE1	1:A:265:GLN:HA	2.19	0.42
1:A:151:SER:HA	1:A:152:PRO:HD3	1.91	0.42
1:B:31:LEU:H	1:B:70:LYS:HA	1.84	0.42
1:B:88:THR:CG2	1:B:89:LEU:N	2.81	0.42
1:B:88:THR:HG23	1:B:215:VAL:HG22	2.01	0.42
2:B:302:3VV:H33	2:B:302:3VV:H29	1.71	0.42
2:A:302:3VV:H8	1:B:27:PRO:CG	2.49	0.41
1:B:86:LEU:CD1	1:B:87:MET:N	2.69	0.41
1:B:145:ASP:OD1	1:B:173:GLN:NE2	2.53	0.41
1:A:198:ASN:OD1	1:B:100:ASP:HB3	2.20	0.41
1:A:201:TYR:HA	1:A:204:ILE:HD12	2.02	0.41
1:B:34:GLU:C	1:B:36:GLU:HB3	2.41	0.41
1:B:37:LEU:HB3	1:B:45:ARG:NH2	2.14	0.41
1:B:60:TRP:CE2	1:B:84:ASP:OD2	2.73	0.41
1:B:87:MET:HA	1:B:92:GLU:HG2	2.00	0.41
1:B:88:THR:HG23	1:B:89:LEU:O	2.21	0.41
1:A:145:ASP:O	1:A:148:ILE:HG22	2.20	0.41
1:B:87:MET:HA	1:B:92:GLU:CG	2.50	0.41
1:A:31:LEU:HA	1:A:32:PRO:HD2	1.83	0.41
1:A:209:LYS:NZ	1:A:213:ASP:OD2	2.47	0.41
1:B:109:SER:HB3	1:B:252:ILE:HD13	2.03	0.41
1:B:17:ILE:HD13	1:B:92:GLU:OE2	2.21	0.41
1:B:49:ARG:O	1:B:53:GLN:CD	2.58	0.41
1:B:63:ILE:HD12	1:B:64:GLN:CB	2.50	0.41
1:B:92:GLU:HG3	1:B:93:ASN:N	2.36	0.41
1:B:128:MET:HG2	1:B:188:LEU:HD21	2.02	0.41
1:B:257:ILE:HG23	1:B:261:HIS:HE1	1.83	0.41
1:A:105:ARG:NH1	2:A:301:3VV:H4	2.36	0.41
1:A:230:MET:O	1:A:234:ARG:HG3	2.21	0.41
1:B:176:LEU:HD23	1:B:176:LEU:HA	1.89	0.41
1:A:80:LEU:HB3	1:A:82:ILE:HG12	2.04	0.40
1:B:34:GLU:HB3	1:B:45:ARG:NE	2.36	0.40
1:B:75:MET:O	1:B:76:GLU:C	2.60	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ILE:CG2	1:A:60:TRP:HZ3	2.32	0.40
1:A:34:GLU:HB3	1:A:48:LEU:CD2	2.51	0.40
1:A:176:LEU:HD23	1:A:176:LEU:HA	1.94	0.40
1:A:19:SER:HB3	1:A:24:ARG:HB2	2.03	0.40
1:A:28:GLY:C	1:A:70:LYS:HZ1	2.24	0.40
1:B:38:SER:OG	1:B:45:ARG:HG2	2.19	0.40
1:B:89:LEU:HD23	1:B:91:ALA:H	1.86	0.40
1:A:93:ASN:ND2	1:A:95:THR:HA	2.37	0.40
1:B:87:MET:HB3	1:B:88:THR:CG2	2.51	0.40
1:B:209:LYS:NZ	1:B:213:ASP:OD2	2.54	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	257/279 (92%)	225 (88%)	17 (7%)	15 (6%)	1 4
1	B	263/279 (94%)	224 (85%)	22 (8%)	17 (6%)	1 3
All	All	520/558 (93%)	449 (86%)	39 (8%)	32 (6%)	1 4

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	LYS
1	A	74	PHE
1	A	76	GLU
1	A	95	THR
1	B	36	GLU
1	B	75	MET
1	B	80	LEU
1	B	85	THR

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Mol	Chain	Res	Type
1	B	89	LEU
1	B	94	ALA
1	B	122	GLU
1	B	271	PRO
1	A	65	HIS
1	A	73	GLN
1	A	75	MET
1	A	80	LEU
1	A	88	THR
1	B	64	GLN
1	B	65	HIS
1	B	87	MET
1	B	93	ASN
1	B	95	THR
1	B	270	LEU
1	A	68	PRO
1	A	78	SER
1	A	84	ASP
1	A	265	GLN
1	B	32	PRO
1	B	88	THR
1	A	77	THR
1	A	27	PRO
1	B	33	ALA

5.3.2 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	223/242 (92%)	212 (95%)	11 (5%)	25	57
1	B	228/242 (94%)	220 (96%)	8 (4%)	36	70
All	All	451/484 (93%)	432 (96%)	19 (4%)	30	63

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	GLU
1	A	24	ARG
1	A	35	ARG
1	A	64	GLN
1	A	78	SER
1	A	80	LEU
1	A	82	ILE
1	A	83	LEU
1	A	86	LEU
1	A	92	GLU
1	A	96	SER
1	B	38	SER
1	B	45	ARG
1	B	48	LEU
1	B	82	ILE
1	B	85	THR
1	B	86	LEU
1	B	88	THR
1	B	244	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	65	HIS
1	A	73	GLN
1	A	107	ASN
1	B	158	GLN
1	B	182	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	3VV	B	302	-	61,69,69	2.83	14 (22%)	72,95,95	1.51	12 (16%)
2	3VV	B	301	-	61,69,69	2.80	15 (24%)	72,95,95	1.39	12 (16%)
2	3VV	A	302	-	61,69,69	2.80	14 (22%)	72,95,95	1.37	10 (13%)
2	3VV	A	301	-	61,69,69	2.78	14 (22%)	72,95,95	1.40	10 (13%)

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	3VV	B	302	-	-	29/64/84/84	0/3/3/3
2	3VV	B	301	-	-	15/64/84/84	0/3/3/3
2	3VV	A	302	-	-	26/64/84/84	0/3/3/3
2	3VV	A	301	-	-	17/64/84/84	0/3/3/3

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	3VV	C50-C52	-12.60	1.24	1.52
2	A	301	3VV	C50-C52	-12.38	1.25	1.52
2	B	302	3VV	C50-C52	-12.27	1.25	1.52
2	B	301	3VV	C50-C52	-12.16	1.25	1.52
2	B	301	3VV	O48-C49	7.95	1.52	1.41
2	A	301	3VV	O48-C49	7.83	1.52	1.41
2	A	302	3VV	O48-C49	7.77	1.51	1.41
2	B	302	3VV	O48-C49	7.52	1.51	1.41
2	B	301	3VV	O48-C47	-7.45	1.28	1.45
2	B	302	3VV	O48-C47	-7.34	1.28	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	3VV	O48-C47	-7.21	1.28	1.45
2	A	301	3VV	O48-C47	-7.16	1.29	1.45
2	B	302	3VV	C29-N28	6.40	1.47	1.33
2	B	302	3VV	C24-N23	6.28	1.47	1.33
2	A	301	3VV	C29-N28	6.00	1.46	1.33
2	A	302	3VV	C29-N28	5.97	1.46	1.33
2	B	301	3VV	C29-N28	5.97	1.46	1.33
2	B	301	3VV	C24-N23	5.81	1.46	1.33
2	A	302	3VV	C24-N23	5.81	1.46	1.33
2	A	301	3VV	C24-N23	5.80	1.46	1.33
2	B	302	3VV	C17-C18	4.64	1.55	1.50
2	B	302	3VV	C52-C47	4.51	1.65	1.52
2	A	301	3VV	C52-C47	4.48	1.64	1.52
2	B	301	3VV	C52-C47	4.40	1.64	1.52
2	A	302	3VV	C52-C47	4.33	1.64	1.52
2	B	301	3VV	C17-C18	4.32	1.55	1.50
2	A	302	3VV	C17-C18	4.31	1.55	1.50
2	B	302	3VV	O51-C50	4.20	1.52	1.43
2	B	301	3VV	O51-C50	4.14	1.52	1.43
2	A	301	3VV	C17-C18	4.13	1.55	1.50
2	A	302	3VV	O51-C50	4.11	1.52	1.43
2	A	301	3VV	O51-C50	4.11	1.52	1.43
2	B	302	3VV	C18-S20	3.81	1.85	1.76
2	B	301	3VV	C18-S20	3.58	1.84	1.76
2	A	301	3VV	C18-S20	3.56	1.84	1.76
2	A	302	3VV	C18-S20	3.44	1.84	1.76
2	A	302	3VV	C62-N63	3.41	1.46	1.34
2	B	301	3VV	C62-N63	3.40	1.46	1.34
2	B	302	3VV	C62-N63	3.39	1.46	1.34
2	A	301	3VV	C62-N63	3.37	1.46	1.34
2	B	302	3VV	P54-O53	2.89	1.64	1.59
2	A	302	3VV	P54-O53	2.89	1.64	1.59
2	A	301	3VV	P54-O53	2.77	1.64	1.59
2	B	301	3VV	P54-O53	2.67	1.64	1.59
2	A	301	3VV	O30-C29	-2.45	1.18	1.23
2	B	302	3VV	O30-C29	-2.40	1.18	1.23
2	B	301	3VV	C50-C49	2.39	1.57	1.53
2	B	301	3VV	O30-C29	-2.35	1.18	1.23
2	A	302	3VV	O30-C29	-2.35	1.18	1.23
2	B	301	3VV	O25-C24	-2.28	1.18	1.23
2	A	301	3VV	O25-C24	-2.22	1.18	1.23
2	A	301	3VV	C50-C49	2.17	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	302	3VV	O25-C24	-2.14	1.18	1.23
2	A	302	3VV	O25-C24	-2.13	1.18	1.23
2	A	302	3VV	C50-C49	2.08	1.56	1.53
2	B	302	3VV	C50-C49	2.03	1.56	1.53
2	B	301	3VV	P42-O45	2.03	1.67	1.59

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	302	3VV	C17-C18-S20	6.10	120.55	113.46
2	A	301	3VV	C17-C18-S20	5.76	120.16	113.46
2	B	301	3VV	C17-C18-S20	5.59	119.96	113.46
2	A	302	3VV	C17-C18-S20	4.85	119.11	113.46
2	B	302	3VV	N66-C65-N64	-4.33	121.91	128.68
2	A	301	3VV	N66-C65-N64	-4.31	121.94	128.68
2	A	302	3VV	N66-C65-N64	-4.21	122.09	128.68
2	B	301	3VV	N66-C65-N64	-3.92	122.55	128.68
2	B	302	3VV	C26-C24-N23	3.87	122.94	116.42
2	B	302	3VV	O19-C18-C17	-3.76	119.55	123.99
2	A	301	3VV	O19-C18-C17	-3.19	120.22	123.99
2	A	301	3VV	P42-O41-P38	-3.15	122.02	132.83
2	A	302	3VV	C35-C33-C36	3.15	113.37	108.23
2	B	302	3VV	P42-O41-P38	-3.11	122.16	132.83
2	A	302	3VV	O19-C18-C17	-3.02	120.42	123.99
2	B	301	3VV	O19-C18-C17	-2.92	120.55	123.99
2	A	301	3VV	C27-N28-C29	-2.87	117.46	122.59
2	A	302	3VV	P42-O41-P38	-2.78	123.28	132.83
2	B	301	3VV	P42-O41-P38	-2.73	123.46	132.83
2	B	302	3VV	C21-S20-C18	2.66	110.14	101.87
2	A	301	3VV	O37-C36-C33	-2.60	106.37	110.55
2	B	301	3VV	C27-N28-C29	-2.57	118.00	122.59
2	A	302	3VV	O30-C29-N28	-2.50	117.63	122.99
2	B	301	3VV	C67-C61-N60	-2.47	106.82	109.40
2	B	301	3VV	O19-C18-S20	-2.40	119.50	122.61
2	B	302	3VV	O25-C24-N23	-2.37	118.53	123.01
2	A	301	3VV	O19-C18-S20	-2.31	119.62	122.61
2	B	301	3VV	C52-C50-C49	2.29	104.95	99.89
2	B	302	3VV	C52-C50-C49	2.25	104.87	99.89
2	B	302	3VV	C67-C61-N60	-2.24	107.07	109.40
2	B	301	3VV	O48-C49-C50	-2.23	103.66	106.93
2	B	302	3VV	C50-C52-C47	2.20	107.13	103.22
2	B	301	3VV	C21-C22-N23	-2.16	107.88	112.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	302	3VV	C27-N28-C29	-2.14	118.77	122.59
2	A	302	3VV	C26-C24-N23	2.11	119.97	116.42
2	A	301	3VV	C52-C50-C49	2.11	104.56	99.89
2	B	302	3VV	O19-C18-S20	-2.09	119.90	122.61
2	A	302	3VV	C22-N23-C24	-2.09	118.96	122.84
2	A	302	3VV	C46-C47-C52	-2.09	107.48	114.40
2	B	301	3VV	C22-N23-C24	-2.05	119.02	122.84
2	B	302	3VV	O48-C49-C50	-2.04	103.94	106.93
2	A	301	3VV	C67-C61-N60	-2.03	107.29	109.40
2	B	301	3VV	C26-C24-N23	2.02	119.82	116.42
2	A	301	3VV	C35-C33-C31	2.01	112.31	108.82

There are no chirality outliers.

All (87) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	3VV	C36-O37-P38-O41
2	A	301	3VV	C46-O45-P42-O41
2	A	301	3VV	C46-O45-P42-O43
2	A	301	3VV	C46-O45-P42-O44
2	A	301	3VV	O45-C46-C47-O48
2	A	302	3VV	C22-C21-S20-C18
2	A	302	3VV	C36-O37-P38-O39
2	A	302	3VV	C36-O37-P38-O40
2	A	302	3VV	C36-O37-P38-O41
2	A	302	3VV	C17-C18-S20-C21
2	A	302	3VV	O19-C18-S20-C21
2	A	302	3VV	N28-C29-C31-O32
2	A	302	3VV	N28-C29-C31-C33
2	A	302	3VV	O30-C29-C31-C33
2	A	302	3VV	C31-C33-C36-O37
2	A	302	3VV	C34-C33-C36-O37
2	B	301	3VV	C22-C21-S20-C18
2	B	301	3VV	C36-O37-P38-O40
2	B	302	3VV	S20-C21-C22-N23
2	B	302	3VV	C22-C21-S20-C18
2	B	302	3VV	C46-O45-P42-O44
2	B	302	3VV	C52-O53-P54-O56
2	B	302	3VV	C17-C18-S20-C21
2	B	302	3VV	O19-C18-S20-C21
2	B	302	3VV	N28-C29-C31-O32
2	B	302	3VV	N28-C29-C31-C33

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Mol	Chain	Res	Type	Atoms
2	B	302	3VV	O30-C29-C31-O32
2	B	302	3VV	O30-C29-C31-C33
2	B	302	3VV	C31-C33-C36-O37
2	B	302	3VV	C34-C33-C36-O37
2	B	302	3VV	C26-C24-N23-C22
2	B	302	3VV	C06-C07-C08-C09
2	A	301	3VV	O45-C46-C47-C52
2	B	302	3VV	O25-C24-N23-C22
2	A	302	3VV	C06-C07-C08-C09
2	A	302	3VV	C35-C33-C36-O37
2	B	302	3VV	C35-C33-C36-O37
2	A	302	3VV	C12-C13-C14-C15
2	A	302	3VV	C04-C05-C06-C07
2	B	301	3VV	C03-C04-C05-C06
2	A	301	3VV	C14-C15-C16-C17
2	B	302	3VV	C02-C03-C04-C05
2	A	302	3VV	C05-C06-C07-C08
2	B	301	3VV	C14-C15-C16-C17
2	B	302	3VV	C05-C06-C07-C08
2	B	302	3VV	C12-C13-C14-C15
2	A	302	3VV	C03-C04-C05-C06
2	A	301	3VV	C10-C11-C12-C13
2	B	301	3VV	C10-C11-C12-C13
2	B	302	3VV	C10-C11-C12-C13
2	A	302	3VV	O30-C29-C31-O32
2	B	301	3VV	C05-C06-C07-C08
2	B	302	3VV	C01-C02-C03-C04
2	A	302	3VV	C01-C02-C03-C04
2	B	302	3VV	P42-O41-P38-O40
2	A	302	3VV	C14-C15-C16-C17
2	A	301	3VV	S20-C21-C22-N23
2	B	301	3VV	O45-C46-C47-O48
2	A	301	3VV	C24-C26-C27-N28
2	B	301	3VV	C47-C46-O45-P42
2	B	302	3VV	C46-O45-P42-O41
2	A	301	3VV	P38-O41-P42-O43
2	A	302	3VV	C33-C36-O37-P38
2	B	301	3VV	C11-C12-C13-C14
2	A	301	3VV	C36-O37-P38-O39
2	B	302	3VV	C46-O45-P42-O43
2	B	301	3VV	C13-C14-C15-C16
2	B	301	3VV	C04-C05-C06-C07

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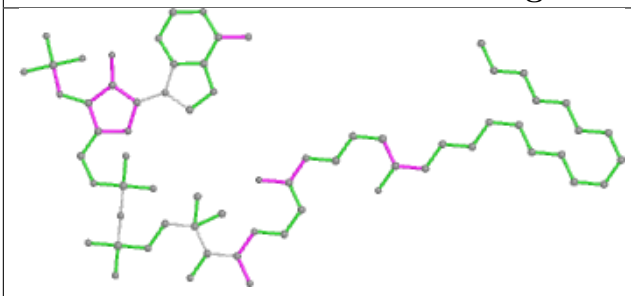
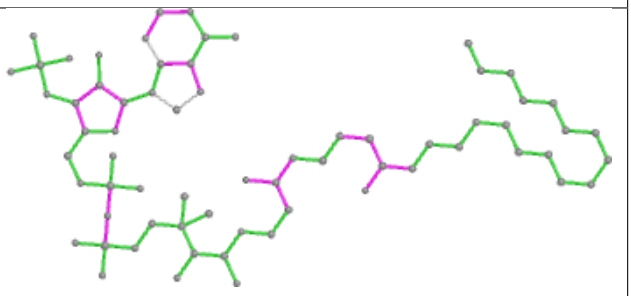
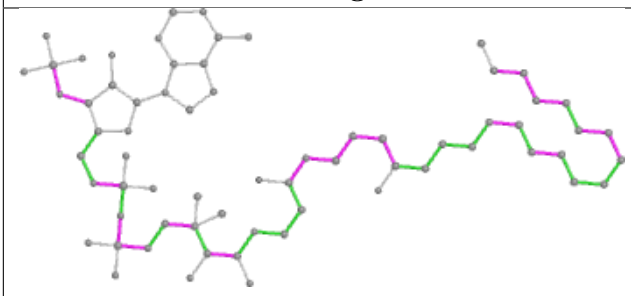
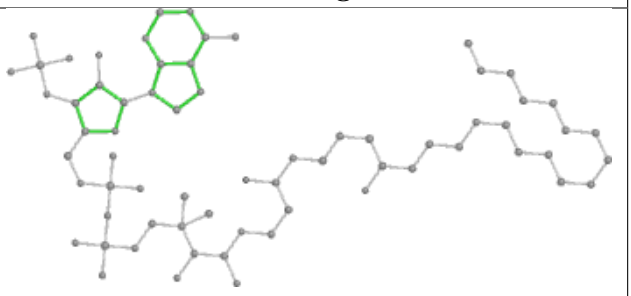
Mol	Chain	Res	Type	Atoms
2	A	302	3VV	O45-C46-C47-O48
2	A	302	3VV	P42-O41-P38-O39
2	B	301	3VV	S20-C21-C22-N23
2	B	302	3VV	C21-C22-N23-C24
2	A	301	3VV	C22-C21-S20-C18
2	A	302	3VV	O45-C46-C47-C52
2	B	302	3VV	C47-C52-O53-P54
2	A	301	3VV	C15-C16-C17-C18
2	A	302	3VV	C15-C16-C17-C18
2	B	301	3VV	C07-C08-C09-C10
2	B	302	3VV	C52-O53-P54-O55
2	A	301	3VV	P42-O41-P38-O39
2	A	301	3VV	P38-O41-P42-O44
2	A	302	3VV	P42-O41-P38-O40
2	A	301	3VV	C36-O37-P38-O40
2	B	301	3VV	C46-O45-P42-O44
2	B	302	3VV	C36-O37-P38-O40
2	B	301	3VV	O45-C46-C47-C52
2	B	302	3VV	C03-C04-C05-C06

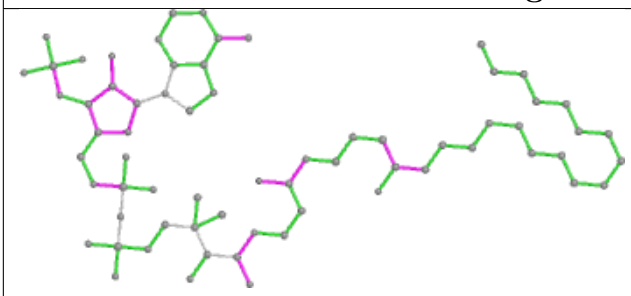
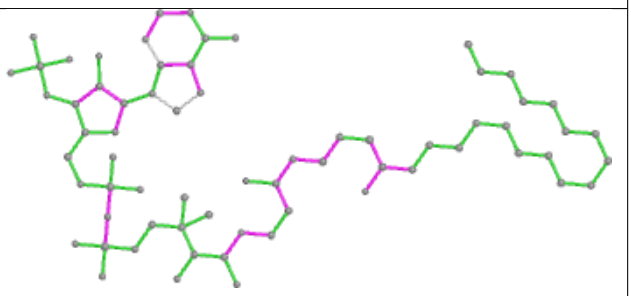
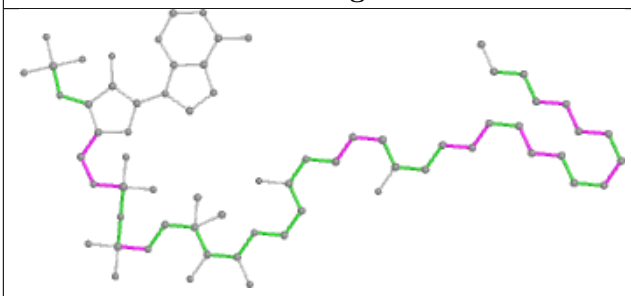
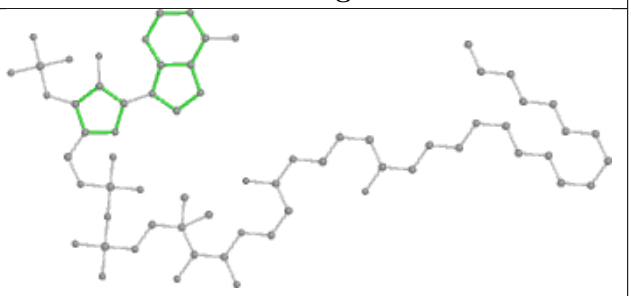
There are no ring outliers.

4 monomers are involved in 32 short contacts:

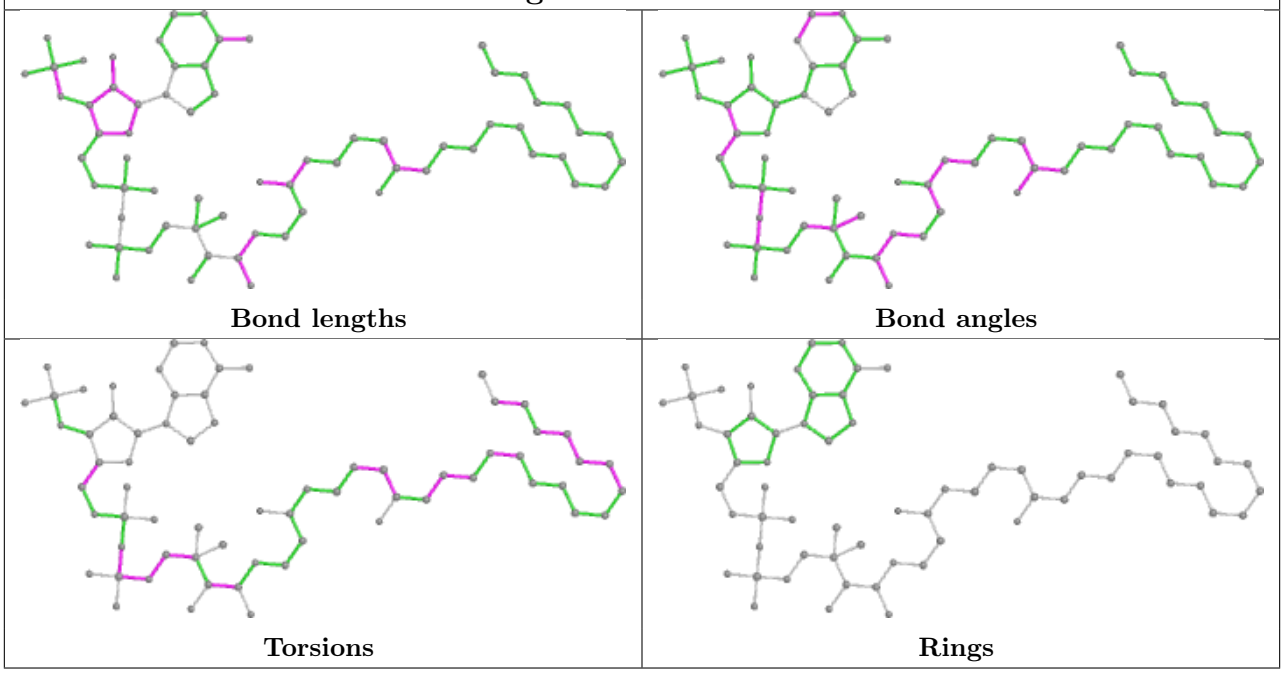
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	3VV	8	0
2	B	301	3VV	9	0
2	A	302	3VV	10	0
2	A	301	3VV	6	0

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 then 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.

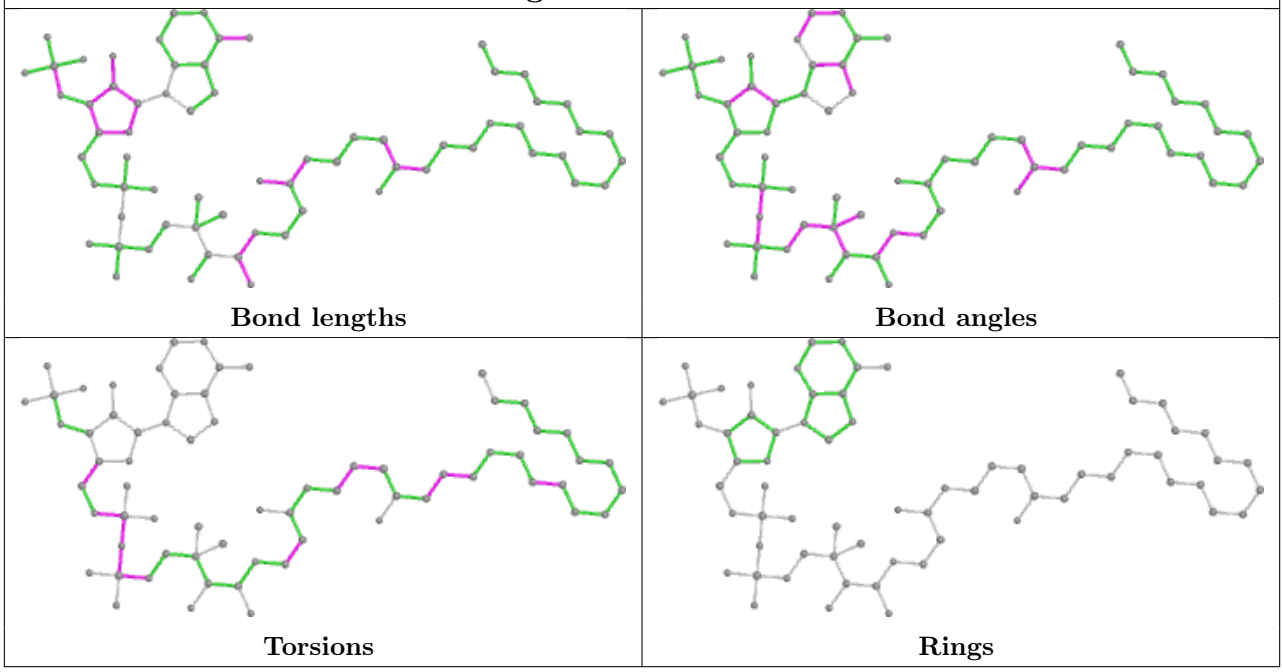
Ligand 3VV B 302	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand 3VV B 301	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand 3VV A 302



Ligand 3VV A 301



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	259/279 (92%)	0.84	45 (17%)	1	1	19, 53, 159, 199	0
1	B	265/279 (94%)	0.99	50 (18%)	1	1	23, 63, 142, 169	0
All	All	524/558 (93%)	0.91	95 (18%)	1	1	19, 57, 150, 199	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	28	GLY	10.7
1	A	39	GLU	8.1
1	B	79	GLY	7.7
1	B	91	ALA	7.4
1	A	93	ASN	6.6
1	B	54	ARG	6.1
1	A	66	GLY	5.9
1	B	89	LEU	5.7
1	B	85	THR	5.7
1	B	90	ASP	5.6
1	A	47	THR	5.6
1	B	8	PRO	5.3
1	A	80	LEU	5.1
1	A	88	THR	5.1
1	A	73	GLN	4.9
1	A	90	ASP	4.9
1	B	44	THR	4.9
1	B	64	GLN	4.8
1	A	77	THR	4.7
1	A	41	ILE	4.7
1	B	60	TRP	4.6
1	A	50	GLU	4.5
1	B	65	HIS	4.4
1	B	76	GLU	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	86	LEU	4.4
1	B	63	ILE	4.4
1	B	10	GLY	4.2
1	A	40	LEU	4.2
1	A	36	GLU	3.9
1	B	77	THR	3.9
1	B	74	PHE	3.8
1	A	87	MET	3.8
1	A	54	ARG	3.7
1	B	165	SER	3.6
1	B	66	GLY	3.6
1	B	96	SER	3.5
1	A	164	ASP	3.5
1	A	27	PRO	3.5
1	B	78	SER	3.4
1	A	266	MET	3.4
1	B	28	GLY	3.4
1	A	81	HIS	3.4
1	B	50	GLU	3.3
1	B	62	THR	3.3
1	A	95	THR	3.2
1	B	88	THR	3.2
1	B	84	ASP	3.2
1	A	65	HIS	3.1
1	B	30	ILE	3.1
1	B	68	PRO	3.1
1	A	89	LEU	3.1
1	A	60	TRP	3.1
1	B	48	LEU	3.0
1	A	78	SER	3.0
1	A	35	ARG	3.0
1	B	29	SER	3.0
1	A	46	THR	2.9
1	B	159	GLN	2.9
1	A	85	THR	2.9
1	A	21	TRP	2.9
1	A	29	SER	2.9
1	A	63	ILE	2.9
1	B	87	MET	2.8
1	A	44	THR	2.8
1	A	23	GLY	2.8
1	A	24	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	93	ASN	2.7
1	A	94	ALA	2.7
1	A	8	PRO	2.7
1	A	67	LYS	2.7
1	B	17	ILE	2.7
1	B	39	GLU	2.7
1	B	40	LEU	2.7
1	B	80	LEU	2.5
1	A	64	GLN	2.5
1	B	35	ARG	2.5
1	A	74	PHE	2.5
1	A	62	THR	2.5
1	B	55	LEU	2.5
1	B	86	LEU	2.4
1	A	49	ARG	2.3
1	A	92	GLU	2.3
1	B	9	ALA	2.3
1	B	16	ILE	2.2
1	B	33	ALA	2.2
1	B	243	GLY	2.2
1	A	76	GLU	2.2
1	B	15	TYR	2.2
1	B	94	ALA	2.2
1	B	59	GLY	2.2
1	A	96	SER	2.1
1	B	83	LEU	2.1
1	B	46	THR	2.1
1	B	81	HIS	2.1
1	B	73	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

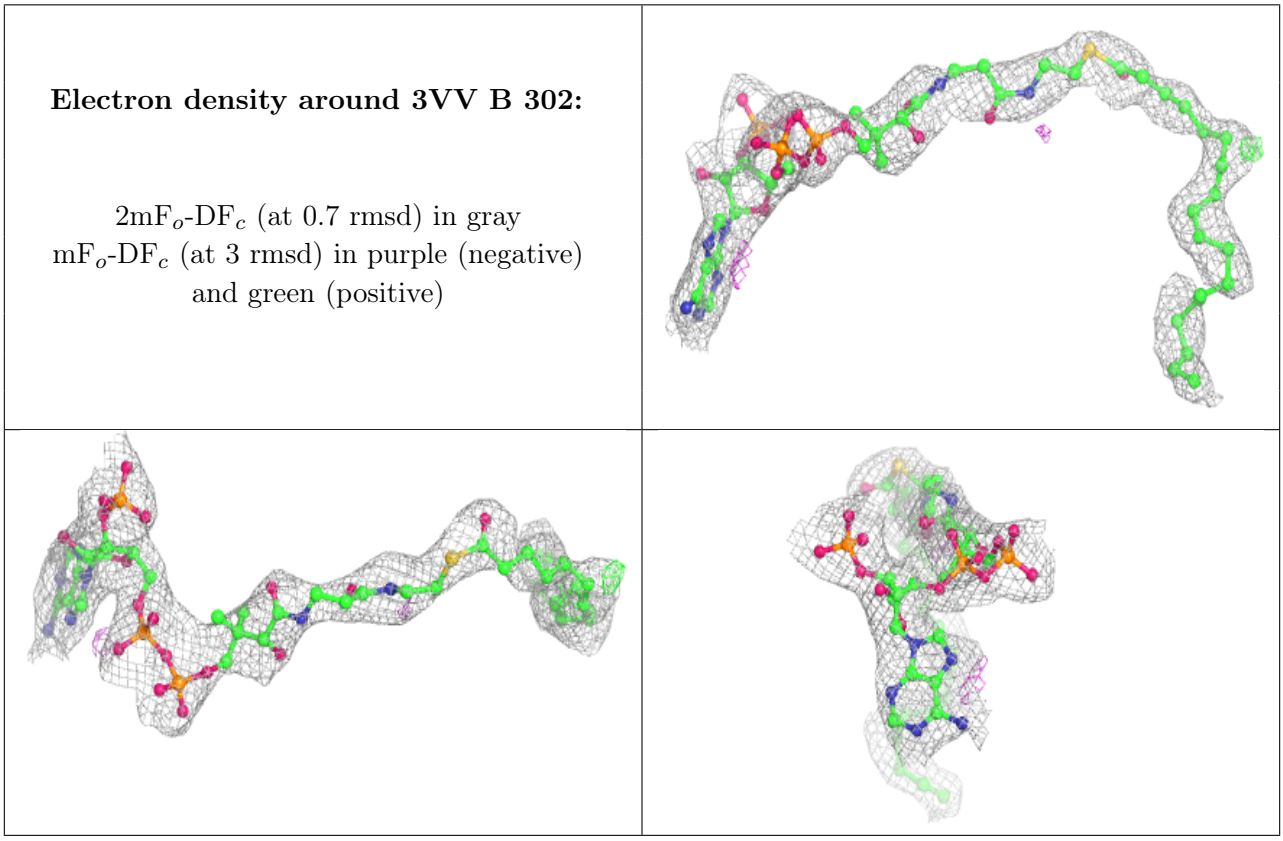
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	3VV	B	302	67/67	0.81	0.24	42,77,112,125	0
2	3VV	A	302	67/67	0.85	0.22	19,67,94,124	0
2	3VV	B	301	67/67	0.86	0.26	18,56,87,105	22
2	3VV	A	301	67/67	0.92	0.16	11,31,80,107	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.

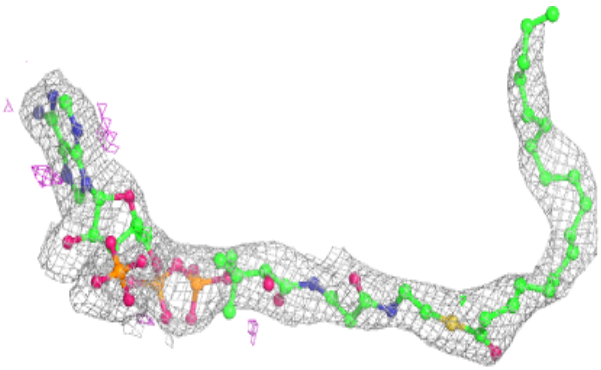
Electron density around 3VV B 302:

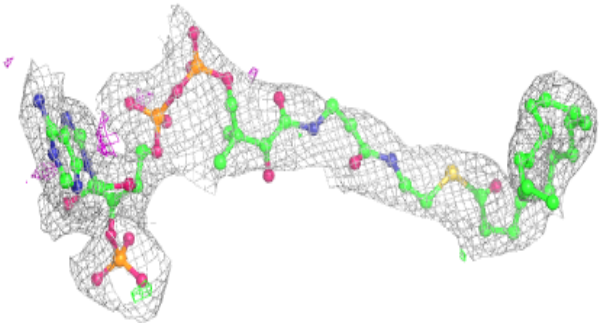
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

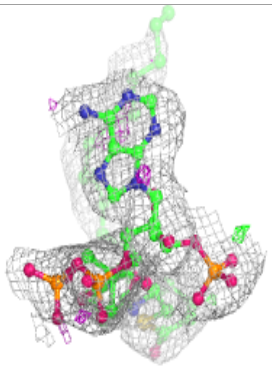


Electron density around 3VV A 302:

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

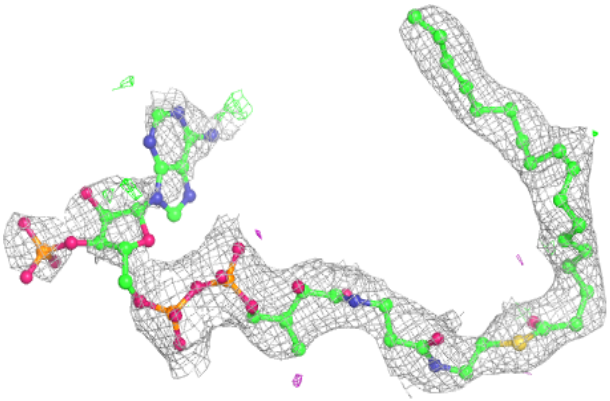


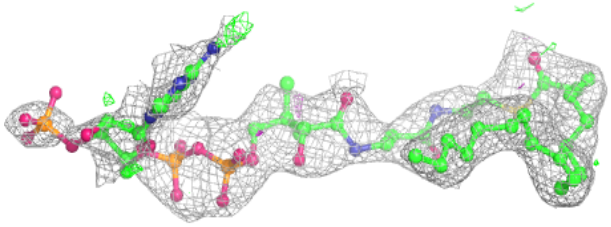


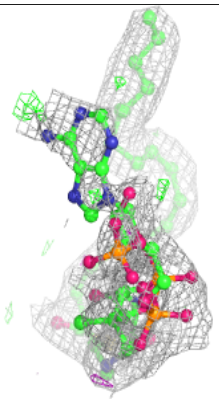


Electron density around 3VV B 301:

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

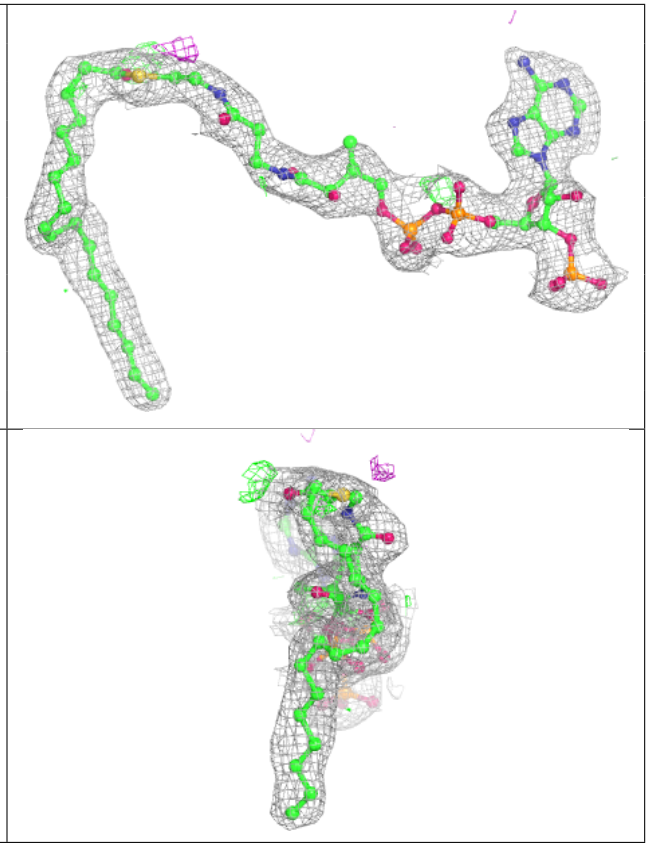






Electron density around 3VV A 301:

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



6.5 Other polymers (i)

There are no such residues in this entry.