



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:31 AM UTC

PDB ID : 3KRP / pdb_00003krp
Title : Mint heterotetrameric geranyl pyrophosphate synthase in complex with magnesium and GPP
Authors : Chang, T.-H.; Hsieh, F.-L.; Ko, T.-P.; Wang, A.H.-J.
Deposited on : 2009-11-19
Resolution : 2.42 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

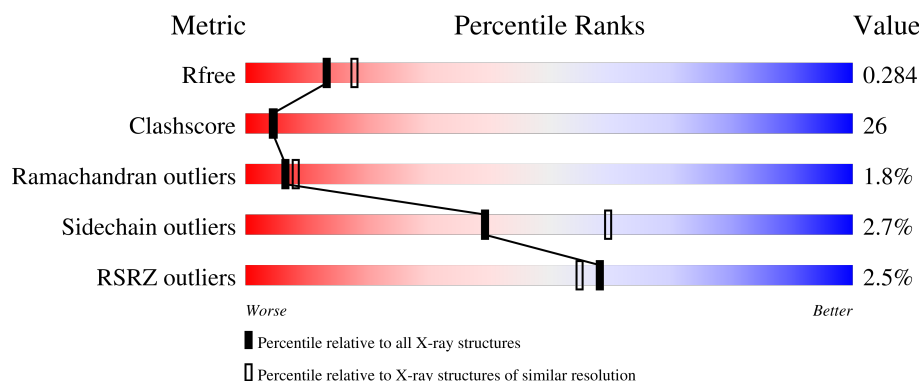
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	295	 59% 34% • 5%
1	D	295	 3% 49% 44% • •
2	B	274	 5% 50% 40% • 7%
2	C	274	 % 52% 37% 5% 5%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9010 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 Geranyl diphosphate synthase large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	0	0	0
			2120	1338	372	392	18			
1	D	284	Total	C	N	O	S	0	0	0
			2155	1359	378	400	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q9SBR3
D	1	MET	-	expression tag	UNP Q9SBR3

- Molecule 2 is a protein called Geranyl diphosphate synthase small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	256	Total	C	N	O	S	0	0	0
			1929	1216	334	365	14			
2	C	259	Total	C	N	O	S	0	0	0
			1954	1233	339	368	14			

There are 18 discrepancies between the modelled and reference sequences:

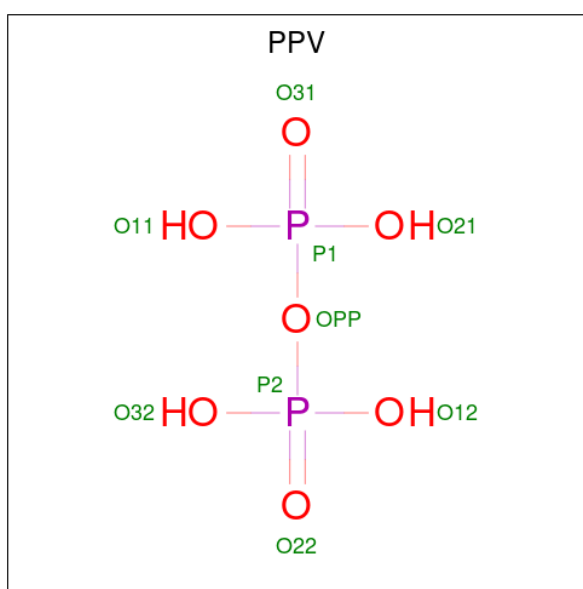
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	expression tag	UNP Q9SBR4
B	267	HIS	-	expression tag	UNP Q9SBR4
B	268	HIS	-	expression tag	UNP Q9SBR4
B	269	HIS	-	expression tag	UNP Q9SBR4
B	270	HIS	-	expression tag	UNP Q9SBR4
B	271	HIS	-	expression tag	UNP Q9SBR4
B	272	HIS	-	expression tag	UNP Q9SBR4
B	273	HIS	-	expression tag	UNP Q9SBR4
B	274	HIS	-	expression tag	UNP Q9SBR4
C	1	MET	-	expression tag	UNP Q9SBR4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	267	HIS	-	expression tag	UNP Q9SBR4
C	268	HIS	-	expression tag	UNP Q9SBR4
C	269	HIS	-	expression tag	UNP Q9SBR4
C	270	HIS	-	expression tag	UNP Q9SBR4
C	271	HIS	-	expression tag	UNP Q9SBR4
C	272	HIS	-	expression tag	UNP Q9SBR4
C	273	HIS	-	expression tag	UNP Q9SBR4
C	274	HIS	-	expression tag	UNP Q9SBR4

- Molecule 3 is PYROPHOSPHATE (CCD ID: PPV) (formula: $\text{H}_4\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 9 7 2	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

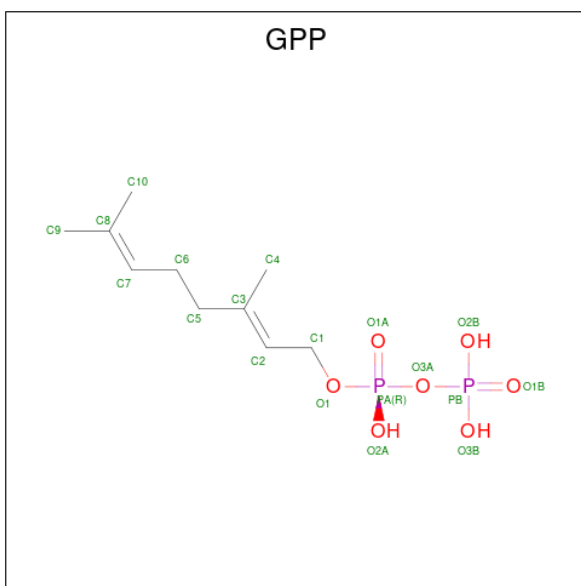
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mg 2 2	0	0
4	D	2	Total Mg 2 2	0	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is GERANYL DIPHOSPHATE (CCD ID: GPP) (formula: $C_{10}H_{20}O_7P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	1	Total	C	O	P	0	0
			19	10	7	2		

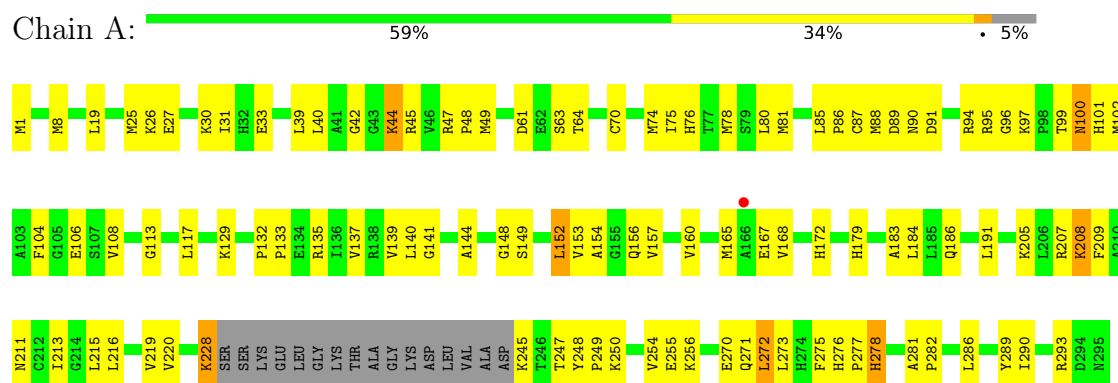
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	224	Total 224	O 224	0	0
7	B	177	Total 177	O 177	0	0
7	C	221	Total 221	O 221	0	0
7	D	194	Total 194	O 194	0	0

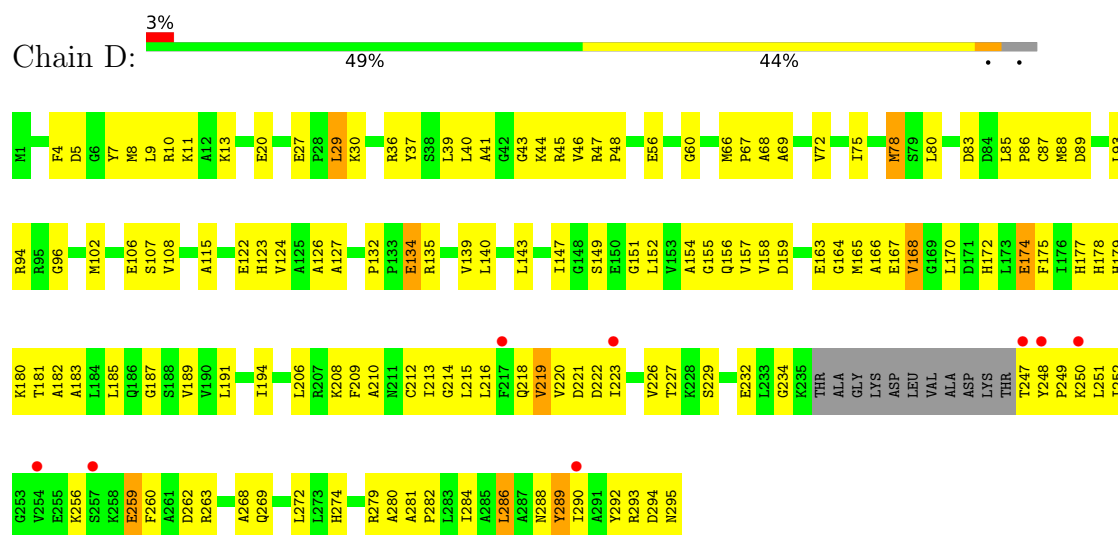
3 Residue-property plots [i](#)

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.

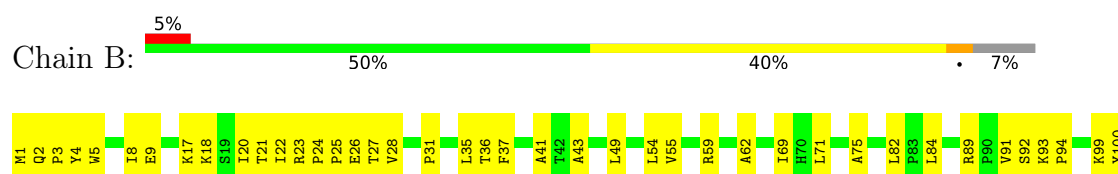
• Molecule 1: Geranyl diphosphate synthase large subunit

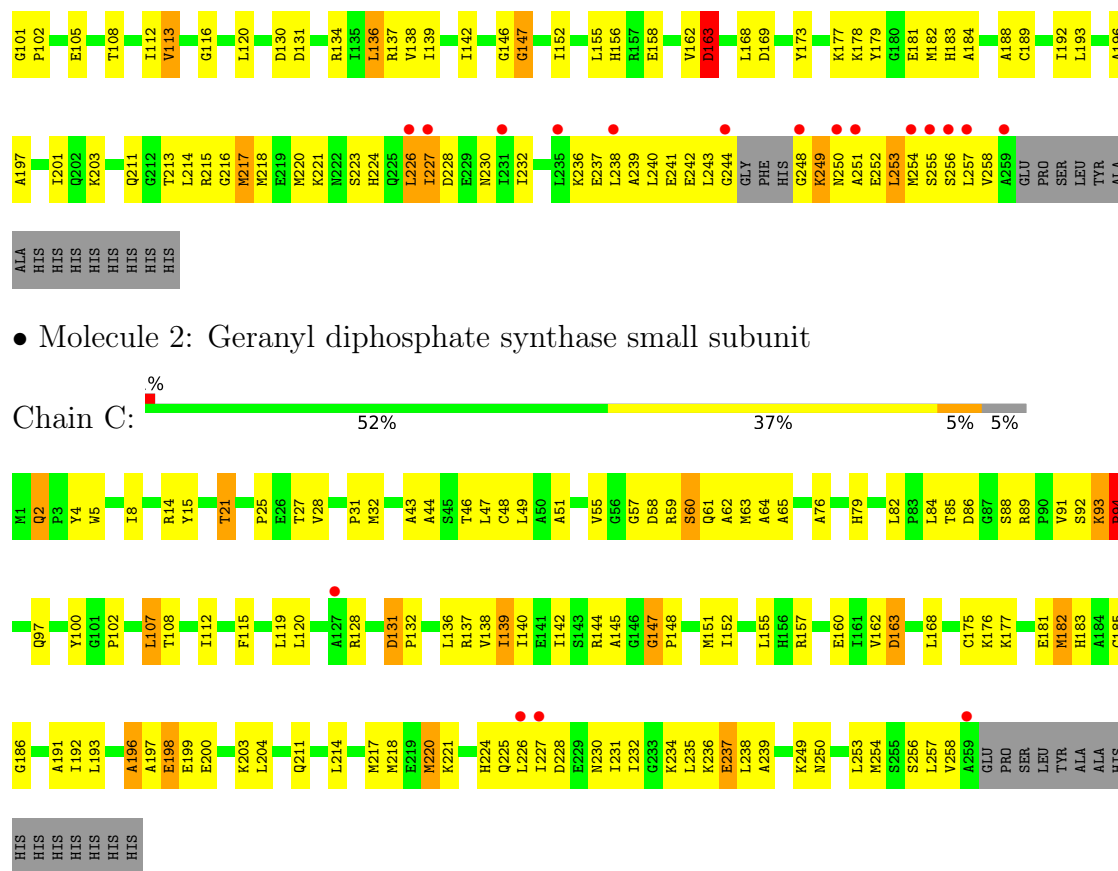


• Molecule 1: Geranyl diphosphate synthase large subunit



• Molecule 2: Geranyl diphosphate synthase small subunit





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.26Å 109.02Å 182.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.42 30.00 – 2.42	Depositor EDS
% Data completeness (in resolution range)	85.9 (30.00-2.42) 85.9 (30.00-2.42)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.82 (at 2.24Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.223 , 0.284 0.224 , 0.284	Depositor DCC
R_{free} test set	2171 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	1.003	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 59.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9010	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.52	0/2155	0.95	4/2903 (0.1%)
1	D	0.48	0/2190	0.96	9/2948 (0.3%)
2	B	0.45	0/1965	0.97	9/2656 (0.3%)
2	C	0.48	0/1993	0.94	7/2695 (0.3%)
All	All	0.48	0/8303	0.96	29/11202 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	226	LEU	N-CA-C	-7.91	101.82	112.26
2	B	254	MET	N-CA-C	-6.98	103.86	112.38
1	D	289	TYR	N-CA-C	-6.83	104.43	112.89
2	B	147	GLY	CA-C-N	6.64	126.33	119.56
2	B	147	GLY	C-N-CA	6.64	126.33	119.56
2	B	253	LEU	N-CA-C	-6.56	105.82	113.88
2	C	147	GLY	CA-C-N	6.46	127.01	120.04
2	C	147	GLY	C-N-CA	6.46	127.01	120.04
2	B	217	MET	N-CA-C	-6.08	104.77	111.82
1	D	29	LEU	N-CA-C	6.08	118.43	111.02
2	C	220	MET	N-CA-C	-6.05	105.48	112.92
1	D	85	LEU	CA-C-N	6.03	125.75	119.05
1	D	85	LEU	C-N-CA	6.03	125.75	119.05
2	C	85	THR	N-CA-C	6.03	117.85	111.28
1	A	100	ASN	N-CA-C	5.98	118.28	111.11
2	C	2	GLN	CA-C-N	5.83	125.30	119.24
2	C	2	GLN	C-N-CA	5.83	125.30	119.24
1	D	39	LEU	N-CA-C	5.70	117.29	111.14
1	D	259	GLU	N-CA-C	-5.62	105.53	112.90
1	D	172	HIS	N-CA-C	-5.62	105.07	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	113	GLY	N-CA-C	-5.61	105.59	112.77
2	B	184	ALA	N-CA-C	-5.44	105.04	110.97
1	A	85	LEU	CA-C-N	5.40	125.47	119.32
1	A	85	LEU	C-N-CA	5.40	125.47	119.32
2	B	113	VAL	CB-CA-C	-5.32	108.64	113.70
2	C	107	LEU	N-CA-C	5.30	116.74	111.07
2	B	99	LYS	N-CA-C	-5.28	105.17	111.03
1	D	219	VAL	N-CA-C	-5.11	104.66	111.05
1	D	43	GLY	N-CA-C	-5.06	104.17	111.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2120	0	2146	86	0
1	D	2155	0	2182	139	0
2	B	1929	0	1928	126	0
2	C	1954	0	1948	104	0
3	A	9	0	0	0	0
4	A	2	0	0	0	0
4	D	2	0	0	0	0
5	C	4	0	6	1	0
6	D	19	0	17	2	0
7	A	224	0	0	13	0
7	B	177	0	0	17	0
7	C	221	0	0	11	0
7	D	194	0	0	12	0
All	All	9010	0	8227	432	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 26.

All (432) 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:23:ARG:HB2	2:B:24:PRO:HD2	1.37	1.06
2:B:249:LYS:HE2	2:B:253:LEU:HD13	1.46	0.98
2:B:8:ILE:HG12	2:B:59:ARG:HH12	1.28	0.97
2:C:163:ASP:HA	7:C:702:HOH:O	1.65	0.95
1:A:168:VAL:HG13	1:A:172:HIS:HB3	1.48	0.94
2:B:249:LYS:HD3	2:B:253:LEU:HB2	1.49	0.93
2:C:221:LYS:HD3	2:C:227:ILE:HG22	1.50	0.90
2:C:43:ALA:O	2:C:47:LEU:HB2	1.70	0.90
2:C:14:ARG:HG2	7:C:535:HOH:O	1.73	0.88
1:A:139:VAL:HG13	1:A:191:LEU:HD22	1.57	0.85
2:B:248:GLY:N	2:B:251:ALA:HB3	1.91	0.84
1:D:260:PHE:HD1	1:D:263:ARG:HH21	1.25	0.84
2:B:55:VAL:HG11	2:B:196:ALA:HB2	1.60	0.82
2:C:8:ILE:HG23	2:C:49:LEU:HD12	1.63	0.81
2:C:28:VAL:HG22	1:D:157:VAL:CG2	2.11	0.81
2:B:2:GLN:HB2	2:B:3:PRO:HD3	1.63	0.80
1:A:289:TYR:O	1:A:293:ARG:HG2	1.82	0.79
1:D:135:ARG:HG3	1:D:135:ARG:HH11	1.46	0.79
1:A:277:PRO:HD2	1:A:278:HIS:CE1	2.17	0.78
1:D:10:ARG:HD2	7:D:394:HOH:O	1.84	0.78
1:D:139:VAL:HG13	1:D:191:LEU:HD22	1.64	0.77
2:C:15:TYR:HB2	2:C:63:MET:HE3	1.66	0.77
1:A:76:HIS:CE1	1:A:80:LEU:HD11	2.20	0.77
2:C:214:LEU:O	2:C:218:MET:HG2	1.86	0.76
2:C:76:ALA:HB2	2:C:151:MET:HE3	1.68	0.76
1:D:9:LEU:O	1:D:13:LYS:HD3	1.86	0.75
2:C:199:GLU:O	2:C:203:LYS:HG2	1.87	0.74
1:A:256:LYS:HE2	7:A:378:HOH:O	1.87	0.73
1:D:216:LEU:O	1:D:220:VAL:HG23	1.88	0.73
2:B:2:GLN:HB2	2:B:3:PRO:CD	2.19	0.72
1:A:183:ALA:HB3	7:A:341:HOH:O	1.89	0.71
2:C:97:GLN:NE2	7:C:773:HOH:O	2.22	0.71
1:A:47:ARG:HB2	1:A:48:PRO:HD3	1.73	0.71
2:B:138:VAL:O	2:B:142:ILE:HG13	1.91	0.71
2:B:249:LYS:O	2:B:253:LEU:N	2.24	0.70
2:C:137:ARG:HD3	2:C:192:ILE:HD13	1.72	0.70
2:B:28:VAL:O	2:B:31:PRO:HD2	1.91	0.70
2:C:230:ASN:O	2:C:234:LYS:HG3	1.92	0.70
1:A:96:GLY:C	1:A:97:LYS:HD2	2.15	0.70
2:B:221:LYS:HG3	2:B:227:ILE:HG21	1.72	0.70
2:C:221:LYS:HD3	2:C:227:ILE:CG2	2.20	0.69
1:A:256:LYS:HE3	7:A:379:HOH:O	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:248:GLY:O	2:B:252:GLU:HB2	1.91	0.69
1:A:286:LEU:O	1:A:290:ILE:HG13	1.93	0.69
2:B:23:ARG:HB2	2:B:24:PRO:CD	2.19	0.69
2:B:92:SER:C	2:B:94:PRO:HD2	2.18	0.69
2:B:137:ARG:HD3	2:B:192:ILE:HD13	1.75	0.69
2:B:253:LEU:HD12	2:B:256:SER:OG	1.93	0.68
1:D:20:GLU:HA	1:D:40:LEU:HD11	1.75	0.68
2:B:1:MET:HG3	2:B:2:GLN:H	1.57	0.68
1:D:135:ARG:NH1	7:D:368:HOH:O	2.27	0.68
2:B:25:PRO:HB2	2:B:27:THR:HG22	1.76	0.68
2:C:51:ALA:HA	2:C:254:MET:HE1	1.74	0.67
2:C:62:ALA:HA	2:C:193:LEU:HD13	1.74	0.67
1:A:19:LEU:HD22	1:A:39:LEU:HD21	1.75	0.67
1:D:168:VAL:HG23	1:D:251:LEU:HD21	1.75	0.67
1:A:64:THR:HG23	1:A:129:LYS:O	1.94	0.67
1:A:149:SER:HB2	7:A:493:HOH:O	1.92	0.67
1:D:252:ILE:O	1:D:256:LYS:HB2	1.94	0.67
2:B:227:ILE:HG13	2:B:227:ILE:O	1.93	0.67
1:A:1:MET:N	7:A:352:HOH:O	2.23	0.67
2:B:220:MET:HB3	2:B:224:HIS:HE1	1.59	0.67
2:B:8:ILE:HG12	2:B:59:ARG:NH1	2.06	0.67
2:C:58:ASP:OD1	2:C:61:GLN:HG3	1.95	0.66
2:B:168:LEU:HD21	2:B:226:LEU:HB2	1.77	0.66
1:A:101:HIS:CD2	1:A:102:MET:HE2	2.31	0.66
2:C:86:ASP:OD2	1:D:106:GLU:HG3	1.96	0.66
1:A:88:MET:HE3	1:A:160:VAL:HG21	1.77	0.66
2:B:177:LYS:O	7:B:508:HOH:O	2.13	0.66
1:D:208:LYS:HD2	7:D:694:HOH:O	1.95	0.66
1:D:260:PHE:HD1	1:D:263:ARG:NH2	1.93	0.66
2:C:145:ALA:HB1	2:C:182:MET:HE3	1.78	0.65
2:B:249:LYS:CD	2:B:253:LEU:HB2	2.23	0.65
1:D:170:LEU:O	1:D:174:GLU:HB2	1.96	0.65
1:A:167:GLU:HG2	7:A:546:HOH:O	1.96	0.65
1:D:80:LEU:HD23	6:D:900:GPP:H92	1.77	0.65
1:A:90:ASN:O	7:A:412:HOH:O	2.13	0.65
2:B:43:ALA:HB3	2:B:183:HIS:HE1	1.62	0.65
2:B:43:ALA:HB3	2:B:183:HIS:CE1	2.32	0.64
1:D:151:GLY:O	1:D:179:HIS:HB3	1.97	0.64
1:A:245:LYS:HD2	1:A:245:LYS:O	1.97	0.64
2:B:188:ALA:O	2:B:192:ILE:HG13	1.97	0.64
1:A:81:MET:HE3	1:A:100:ASN:ND2	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:LYS:HE2	7:D:529:HOH:O	1.96	0.64
1:A:276:HIS:HB3	1:A:278:HIS:ND1	2.12	0.64
1:D:251:LEU:HG	7:D:665:HOH:O	1.96	0.64
2:B:177:LYS:NZ	2:B:181:GLU:OE2	2.32	0.63
1:A:153:VAL:O	1:A:157:VAL:HG23	1.97	0.63
2:B:214:LEU:O	2:B:218:MET:HG2	1.98	0.63
2:B:236:LYS:O	2:B:240:LEU:HG	1.97	0.63
2:C:214:LEU:HD21	2:C:232:ILE:HG23	1.81	0.63
1:A:149:SER:CB	7:A:493:HOH:O	2.47	0.63
2:C:175:CYS:SG	2:C:220:MET:HG3	2.38	0.63
2:C:191:ALA:HB2	2:C:204:LEU:HD12	1.80	0.63
1:D:256:LYS:HA	1:D:259:GLU:CD	2.24	0.63
1:A:165:MET:HE1	7:A:757:HOH:O	1.99	0.62
2:C:64:ALA:HB3	7:C:350:HOH:O	1.98	0.62
2:B:136:LEU:O	2:B:136:LEU:HD12	1.99	0.62
2:B:71:LEU:HD22	2:B:112:ILE:HG23	1.82	0.62
2:C:92:SER:C	2:C:94:PRO:HD2	2.25	0.62
2:C:145:ALA:CB	2:C:182:MET:HE3	2.29	0.62
2:C:136:LEU:O	2:C:140:ILE:HG13	2.00	0.62
2:B:1:MET:HG3	2:B:2:GLN:N	2.14	0.62
2:C:28:VAL:HG22	1:D:157:VAL:HG21	1.82	0.62
1:D:187:GLY:O	1:D:191:LEU:HG	1.99	0.62
1:A:30:LYS:HE2	1:A:108:VAL:HG21	1.81	0.62
1:D:132:PRO:HB2	1:D:134:GLU:OE2	2.00	0.61
1:D:220:VAL:HG22	7:D:576:HOH:O	2.01	0.61
1:A:86:PRO:HG2	1:A:106:GLU:OE2	2.01	0.61
1:D:139:VAL:CG1	1:D:191:LEU:HD22	2.31	0.61
1:D:93:LEU:HG	1:D:234:GLY:O	2.00	0.61
2:C:65:ALA:HA	2:C:120:LEU:HD21	1.83	0.60
2:C:25:PRO:HB2	2:C:27:THR:HG22	1.83	0.60
1:D:93:LEU:HD11	1:D:96:GLY:HA2	1.83	0.60
1:A:270:GLU:HA	1:A:273:LEU:HG	1.84	0.60
1:A:76:HIS:NE2	1:A:80:LEU:HD11	2.17	0.60
1:A:61:ASP:OD2	1:A:63:SER:HB3	2.03	0.59
2:C:227:ILE:HG23	2:C:227:ILE:O	2.01	0.59
2:B:179:TYR:OH	7:B:306:HOH:O	2.16	0.59
1:D:30:LYS:HE2	1:D:108:VAL:HG21	1.84	0.59
2:B:221:LYS:HG3	2:B:227:ILE:CG2	2.32	0.59
1:D:7:TYR:HA	1:D:10:ARG:HH11	1.67	0.59
2:C:237:GLU:O	2:C:239:ALA:N	2.36	0.59
2:B:31:PRO:HA	2:B:100:TYR:CE2	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:MET:HE3	1:A:117:LEU:HB2	1.85	0.58
2:C:162:VAL:HG12	2:C:163:ASP:OD1	2.04	0.58
1:D:20:GLU:HA	1:D:40:LEU:CD1	2.33	0.58
2:C:181:GLU:OE2	2:C:181:GLU:HA	2.04	0.58
1:D:177:HIS:HD1	1:D:218:GLN:HG2	1.66	0.58
2:B:240:LEU:HD21	2:B:258:VAL:HG21	1.86	0.58
2:B:257:LEU:HD21	7:B:505:HOH:O	2.03	0.58
1:A:228:LYS:HB2	1:A:228:LYS:NZ	2.19	0.57
1:D:177:HIS:ND1	1:D:218:GLN:HG2	2.19	0.57
2:C:93:LYS:N	2:C:94:PRO:HD2	2.20	0.57
2:B:147:GLY:HA3	7:B:296:HOH:O	2.04	0.57
2:C:79:HIS:HA	2:C:82:LEU:HG	1.87	0.57
1:D:83:ASP:O	1:D:89:ASP:HB2	2.05	0.57
2:B:137:ARG:NH1	7:B:499:HOH:O	2.37	0.57
2:B:23:ARG:CB	2:B:24:PRO:HD2	2.24	0.56
2:B:213:THR:O	2:B:217:MET:HB2	2.05	0.56
1:D:208:LYS:HG2	7:D:529:HOH:O	2.05	0.56
1:D:222:ASP:HB3	1:D:249:PRO:HD2	1.87	0.56
1:D:281:ALA:HB3	1:D:282:PRO:HD3	1.87	0.56
2:B:120:LEU:HD11	2:B:142:ILE:HD12	1.87	0.56
2:C:214:LEU:C	2:C:214:LEU:HD13	2.30	0.56
2:B:54:LEU:HB2	2:B:250:ASN:OD1	2.05	0.56
2:C:58:ASP:HB2	7:C:634:HOH:O	2.05	0.56
2:C:217:MET:HE2	2:C:231:ILE:HG21	1.88	0.56
1:A:133:PRO:O	1:A:137:VAL:HG23	2.05	0.56
2:C:102:PRO:HB2	1:D:87:CYS:HB2	1.87	0.56
1:D:180:LYS:HG2	6:D:900:GPP:H62	1.86	0.56
2:B:55:VAL:CG1	2:B:196:ALA:HB2	2.33	0.56
2:C:253:LEU:O	2:C:256:SER:OG	2.23	0.56
1:D:252:ILE:O	1:D:256:LYS:HE3	2.06	0.56
1:D:7:TYR:HA	1:D:10:ARG:NH1	2.21	0.55
2:B:9:GLU:OE1	7:B:503:HOH:O	2.18	0.55
2:B:214:LEU:HD23	2:B:258:VAL:HG12	1.88	0.55
1:D:263:ARG:HB2	1:D:263:ARG:NH1	2.21	0.55
1:A:205:LYS:HD3	1:A:271:GLN:O	2.07	0.55
2:B:168:LEU:HD21	2:B:226:LEU:CB	2.37	0.55
2:B:220:MET:HB3	2:B:224:HIS:CE1	2.40	0.55
2:C:43:ALA:HB3	2:C:183:HIS:HE1	1.72	0.55
1:A:75:ILE:O	1:A:78:MET:HG3	2.07	0.55
2:C:4:TYR:CE2	2:C:8:ILE:HD11	2.42	0.55
2:C:221:LYS:HD2	2:C:225:GLN:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:222:ASP:CG	1:D:247:THR:HB	2.30	0.55
1:D:289:TYR:O	1:D:293:ARG:HB3	2.07	0.55
2:C:44:ALA:CB	2:C:183:HIS:HA	2.37	0.55
1:A:95:ARG:NH1	1:A:95:ARG:HG2	2.22	0.54
1:D:68:ALA:HA	1:D:124:VAL:HG22	1.89	0.54
1:A:87:CYS:HB2	2:B:102:PRO:HB2	1.89	0.54
1:A:95:ARG:HG2	1:A:95:ARG:HH11	1.73	0.54
2:B:137:ARG:HD2	7:B:499:HOH:O	2.07	0.54
2:C:51:ALA:CA	2:C:254:MET:HE1	2.38	0.54
1:D:294:ASP:O	1:D:295:ASN:HB3	2.07	0.54
1:D:229:SER:OG	1:D:232:GLU:HG3	2.08	0.53
1:A:140:LEU:HD13	2:B:139:ILE:HD13	1.91	0.53
1:D:37:TYR:O	1:D:41:ALA:HB2	2.08	0.53
1:D:280:ALA:O	1:D:284:ILE:HG13	2.08	0.53
2:B:93:LYS:N	2:B:94:PRO:HD2	2.22	0.53
2:C:76:ALA:CB	2:C:151:MET:HE3	2.39	0.53
2:C:92:SER:CB	2:C:94:PRO:HD2	2.38	0.53
1:D:212:CYS:C	1:D:214:GLY:H	2.14	0.53
1:A:31:ILE:HD11	2:B:156:HIS:CD2	2.43	0.53
2:C:108:THR:O	2:C:112:ILE:HG12	2.08	0.53
1:D:177:HIS:ND1	1:D:218:GLN:CG	2.72	0.53
1:A:281:ALA:HB3	1:A:282:PRO:HD3	1.91	0.53
2:B:91:VAL:HG23	2:B:91:VAL:O	2.09	0.53
1:D:149:SER:O	1:D:154:ALA:HB2	2.08	0.53
1:D:123:HIS:O	1:D:127:ALA:HB3	2.09	0.53
1:A:208:LYS:HE3	1:A:271:GLN:CD	2.34	0.53
2:B:155:LEU:HD23	2:B:178:LYS:HE3	1.91	0.53
2:B:69:ILE:HG23	2:B:182:MET:HE2	1.90	0.52
1:D:223:ILE:O	1:D:227:THR:HG23	2.10	0.52
2:C:84:LEU:HD12	2:C:89:ARG:HG2	1.92	0.52
1:D:66:MET:HB3	1:D:67:PRO:HD3	1.92	0.52
1:D:135:ARG:HG3	1:D:135:ARG:NH1	2.21	0.52
1:A:186:GLN:NE2	1:A:211:ASN:OD1	2.42	0.52
1:D:75:ILE:HD13	1:D:147:ILE:HD13	1.90	0.52
2:B:41:ALA:HB1	7:B:300:HOH:O	2.10	0.52
2:C:197:ALA:O	2:C:198:GLU:C	2.52	0.52
2:B:1:MET:HE1	7:B:430:HOH:O	2.09	0.52
2:B:62:ALA:HA	2:B:193:LEU:HD13	1.91	0.52
1:D:134:GLU:H	1:D:134:GLU:CD	2.19	0.51
1:A:94:ARG:HG2	1:A:95:ARG:HG3	1.93	0.51
2:B:258:VAL:HG22	7:B:616:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:157:ARG:O	2:C:160:GLU:HB2	2.10	0.51
1:D:178:HIS:O	1:D:182:ALA:HB3	2.11	0.51
1:A:89:ASP:OD1	7:A:733:HOH:O	2.18	0.51
2:C:43:ALA:HB3	2:C:183:HIS:CE1	2.46	0.51
2:C:93:LYS:O	2:C:94:PRO:C	2.54	0.51
2:C:8:ILE:CG2	2:C:49:LEU:HD12	2.37	0.51
1:D:56:GLU:HA	1:D:60:GLY:O	2.11	0.50
1:D:215:LEU:O	1:D:219:VAL:HG23	2.10	0.50
2:B:17:LYS:HG3	2:B:37:PHE:CE2	2.46	0.50
1:D:279:ARG:O	1:D:282:PRO:HD2	2.10	0.50
2:B:4:TYR:O	2:B:8:ILE:HG13	2.11	0.50
2:B:249:LYS:HD2	2:B:253:LEU:HD22	1.93	0.50
1:D:93:LEU:HD11	1:D:96:GLY:CA	2.41	0.50
2:B:5:TRP:HZ2	2:B:253:LEU:HG	1.77	0.50
2:B:146:GLY:O	2:B:152:ILE:HG12	2.11	0.50
2:C:88:SER:OG	1:D:107:SER:HB3	2.12	0.50
2:C:177:LYS:HA	2:C:181:GLU:HG2	1.93	0.50
1:D:214:GLY:C	1:D:216:LEU:N	2.68	0.50
2:B:242:GLU:O	2:B:243:LEU:HD23	2.12	0.49
1:D:183:ALA:HB3	7:D:313:HOH:O	2.12	0.49
1:D:250:LYS:HD3	7:D:665:HOH:O	2.10	0.49
1:A:148:GLY:O	1:A:153:VAL:HB	2.12	0.49
2:B:75:ALA:CB	2:B:113:VAL:HG23	2.42	0.49
2:C:211:GLN:HG2	2:C:258:VAL:HG12	1.95	0.49
1:D:248:TYR:HB2	1:D:249:PRO:HD3	1.95	0.49
1:A:149:SER:C	1:A:154:ALA:HB2	2.37	0.49
1:A:209:PHE:CE1	1:A:272:LEU:HD11	2.48	0.49
2:B:162:VAL:HG12	2:B:162:VAL:O	2.13	0.49
2:B:253:LEU:C	2:B:255:SER:N	2.67	0.49
2:C:221:LYS:CD	2:C:227:ILE:HG22	2.33	0.49
1:D:288:ASN:O	1:D:292:TYR:HD1	1.96	0.49
2:B:221:LYS:HD3	2:B:227:ILE:HB	1.95	0.49
2:B:239:ALA:O	2:B:242:GLU:HB2	2.13	0.49
1:D:69:ALA:O	1:D:72:VAL:HG22	2.12	0.49
2:B:101:GLY:O	2:B:105:GLU:HG3	2.13	0.48
1:A:8:MET:SD	1:A:49:MET:HG3	2.53	0.48
2:B:179:TYR:CD1	2:B:216:GLY:HA3	2.48	0.48
2:C:59:ARG:HG3	2:C:60:SER:N	2.28	0.48
2:C:8:ILE:HG22	2:C:46:THR:HG22	1.95	0.48
2:C:93:LYS:N	2:C:94:PRO:CD	2.76	0.48
1:A:148:GLY:HA3	7:B:335:HOH:O	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:230:ASN:HB3	7:B:279:HOH:O	2.13	0.48
2:B:21:THR:O	2:B:21:THR:HG22	2.13	0.48
2:B:211:GLN:CD	2:B:257:LEU:HD23	2.39	0.48
1:A:152:LEU:HA	1:A:179:HIS:O	2.12	0.48
1:A:247:THR:HG21	7:A:736:HOH:O	2.13	0.48
2:B:71:LEU:HD12	2:B:116:GLY:HA2	1.95	0.48
2:C:55:VAL:HG21	2:C:196:ALA:HB2	1.95	0.48
2:C:176:LYS:HE3	7:C:286:HOH:O	2.12	0.48
1:D:288:ASN:ND2	7:D:310:HOH:O	2.47	0.48
1:D:47:ARG:HG2	1:D:185:LEU:HD23	1.95	0.48
7:C:291:HOH:O	1:D:149:SER:HB2	2.12	0.48
1:D:212:CYS:C	1:D:214:GLY:N	2.70	0.48
1:A:70:CYS:O	1:A:74:MET:HG3	2.13	0.48
2:C:51:ALA:HA	2:C:254:MET:CE	2.41	0.48
1:D:214:GLY:C	1:D:216:LEU:H	2.21	0.47
1:A:168:VAL:CG1	1:A:172:HIS:HB3	2.33	0.47
2:B:20:ILE:HD12	2:B:36:THR:HG21	1.95	0.47
2:C:168:LEU:HD13	2:C:231:ILE:HG13	1.95	0.47
1:A:19:LEU:HB3	1:A:40:LEU:HD21	1.96	0.47
2:B:173:TYR:CE2	1:D:29:LEU:HD21	2.49	0.47
2:C:249:LYS:O	2:C:253:LEU:HG	2.15	0.47
2:B:203:LYS:HD2	2:B:244:GLY:H	1.79	0.47
2:B:228:ASP:OD2	2:B:230:ASN:HB2	2.15	0.47
2:B:258:VAL:HG13	7:B:616:HOH:O	2.14	0.47
2:C:139:ILE:HG21	1:D:140:LEU:HD13	1.96	0.47
2:C:228:ASP:OD1	2:C:230:ASN:HB2	2.15	0.47
1:D:164:GLY:O	1:D:166:ALA:N	2.48	0.47
1:A:89:ASP:HB3	1:A:91:ASP:OD1	2.14	0.47
2:B:248:GLY:N	7:B:714:HOH:O	2.48	0.47
2:C:48:CYS:HA	2:C:186:GLY:O	2.15	0.47
2:C:131:ASP:N	2:C:132:PRO:HD2	2.30	0.47
1:D:248:TYR:C	1:D:250:LYS:H	2.23	0.47
2:C:128:ARG:NH1	2:C:128:ARG:HB3	2.30	0.47
1:D:152:LEU:C	1:D:152:LEU:HD23	2.40	0.47
2:B:197:ALA:O	2:B:201:ILE:HG13	2.14	0.46
1:D:4:PHE:HB2	1:D:8:MET:CE	2.45	0.46
1:D:27:GLU:HB2	7:D:578:HOH:O	2.14	0.46
1:D:222:ASP:OD2	1:D:247:THR:HB	2.15	0.46
1:A:254:VAL:O	1:A:255:GLU:C	2.58	0.46
2:B:216:GLY:O	2:B:220:MET:HG2	2.15	0.46
2:B:226:LEU:O	2:B:228:ASP:N	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:181:THR:HG21	1:D:218:GLN:HB2	1.97	0.46
2:B:1:MET:O	2:B:5:TRP:CD1	2.68	0.46
2:B:218:MET:HE1	2:B:232:ILE:CD1	2.45	0.46
2:C:46:THR:HG23	5:C:902:EDO:O2	2.16	0.46
1:D:93:LEU:HD12	1:D:94:ARG:H	1.81	0.46
1:A:19:LEU:HD22	1:A:39:LEU:CD2	2.45	0.46
1:A:101:HIS:HD2	1:A:102:MET:HE2	1.76	0.46
2:B:25:PRO:C	2:B:27:THR:H	2.24	0.46
2:B:169:ASP:CG	1:D:36:ARG:HE	2.23	0.46
2:C:32:MET:HA	2:C:108:THR:CG2	2.46	0.46
1:D:223:ILE:HA	1:D:226:VAL:HG12	1.97	0.46
1:D:286:LEU:O	1:D:290:ILE:HG13	2.15	0.46
1:A:25:MET:O	1:A:26:LYS:HB3	2.16	0.46
2:C:151:MET:O	2:C:155:LEU:HG	2.16	0.46
1:D:157:VAL:HG23	1:D:158:VAL:N	2.31	0.46
1:A:186:GLN:HG3	1:A:207:ARG:HG3	1.96	0.46
2:C:28:VAL:HG22	1:D:157:VAL:HG23	1.93	0.46
2:C:115:PHE:CE1	2:C:119:LEU:HD11	2.51	0.45
2:C:199:GLU:HG3	2:C:200:GLU:N	2.31	0.45
2:B:257:LEU:HD11	7:B:505:HOH:O	2.16	0.45
1:D:5:ASP:O	1:D:9:LEU:HB2	2.15	0.45
1:A:97:LYS:HD2	1:A:97:LYS:N	2.31	0.45
1:A:167:GLU:CG	7:A:546:HOH:O	2.60	0.45
2:C:107:LEU:HD21	1:D:88:MET:HE2	1.98	0.45
1:A:27:GLU:OE1	2:C:157:ARG:NH2	2.50	0.45
2:B:20:ILE:CD1	2:B:36:THR:HG21	2.47	0.45
2:B:1:MET:HG3	2:B:2:GLN:HG2	1.98	0.45
2:B:130:ASP:O	2:B:134:ARG:HG3	2.17	0.45
2:B:162:VAL:O	2:B:163:ASP:HB2	2.17	0.45
2:B:253:LEU:HD12	2:B:256:SER:HG	1.81	0.45
2:C:21:THR:HG22	7:C:766:HOH:O	2.17	0.45
1:D:178:HIS:HE1	7:D:362:HOH:O	1.99	0.45
1:D:135:ARG:O	1:D:139:VAL:HG23	2.17	0.44
2:B:218:MET:HE1	2:B:232:ILE:HD13	1.99	0.44
2:B:137:ARG:HD3	2:B:192:ILE:HG21	1.98	0.44
1:A:132:PRO:O	1:A:135:ARG:N	2.46	0.44
1:D:222:ASP:HB3	1:D:249:PRO:CD	2.47	0.44
2:B:4:TYR:CE2	2:B:8:ILE:HD11	2.52	0.44
2:B:237:GLU:HG2	2:B:241:GLU:OE2	2.16	0.44
1:D:7:TYR:O	1:D:11:LYS:HG2	2.18	0.44
1:D:263:ARG:HB2	1:D:263:ARG:HH11	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:272:LEU:C	1:D:274:HIS:N	2.74	0.44
1:A:132:PRO:HG2	1:A:135:ARG:HB2	1.99	0.44
2:C:25:PRO:HG2	2:C:28:VAL:CG2	2.48	0.44
1:D:135:ARG:HH11	1:D:135:ARG:CG	2.20	0.44
1:A:27:GLU:CD	2:C:157:ARG:HH21	2.25	0.44
2:C:44:ALA:HB2	2:C:183:HIS:CD2	2.52	0.44
1:D:167:GLU:N	1:D:167:GLU:CD	2.76	0.44
2:C:86:ASP:OD2	1:D:106:GLU:CG	2.63	0.44
1:D:135:ARG:NH1	1:D:135:ARG:CG	2.81	0.44
2:C:5:TRP:CZ2	2:C:257:LEU:HB2	2.53	0.43
2:C:31:PRO:HG3	2:C:100:TYR:CE1	2.53	0.43
1:D:212:CYS:HB3	1:D:268:ALA:HB2	1.99	0.43
1:D:269:GLN:HE22	1:D:288:ASN:HD21	1.64	0.43
1:A:45:ARG:C	1:A:48:PRO:HD2	2.43	0.43
2:B:203:LYS:HE3	2:B:243:LEU:HA	2.00	0.43
1:A:33:GLU:HG2	1:A:104:PHE:CE2	2.53	0.43
1:A:44:LYS:HB2	1:A:44:LYS:NZ	2.33	0.43
1:A:213:ILE:HA	1:A:216:LEU:HB3	2.00	0.43
2:B:21:THR:O	2:B:22:ILE:C	2.62	0.43
1:A:245:LYS:HD3	1:A:250:LYS:NZ	2.33	0.43
2:B:8:ILE:HG23	2:B:49:LEU:HD12	1.99	0.43
2:B:93:LYS:N	2:B:94:PRO:CD	2.82	0.43
2:C:59:ARG:HG2	7:C:635:HOH:O	2.19	0.43
1:A:27:GLU:HG3	1:D:27:GLU:HB3	2.01	0.43
1:A:219:VAL:HG22	1:A:248:TYR:CE2	2.53	0.43
1:A:215:LEU:HD12	1:A:215:LEU:HA	1.93	0.43
1:A:276:HIS:HA	1:A:277:PRO:HD3	1.91	0.43
1:D:222:ASP:O	1:D:226:VAL:HG12	2.18	0.43
1:D:272:LEU:C	1:D:274:HIS:H	2.26	0.43
1:D:209:PHE:CZ	1:D:213:ILE:HD13	2.54	0.43
1:D:168:VAL:HG23	1:D:168:VAL:O	2.19	0.42
1:A:247:THR:HB	1:A:249:PRO:HD2	2.00	0.42
2:B:18:LYS:O	2:B:18:LYS:HG2	2.19	0.42
1:D:175:PHE:CD1	1:D:175:PHE:C	2.98	0.42
2:B:220:MET:O	2:B:221:LYS:C	2.62	0.42
1:D:189:VAL:HG21	1:D:210:ALA:HB2	2.00	0.42
1:A:96:GLY:N	7:A:730:HOH:O	2.38	0.42
2:B:82:LEU:HD21	2:B:105:GLU:HB2	2.01	0.42
1:A:216:LEU:O	1:A:220:VAL:HG23	2.20	0.42
1:A:272:LEU:HA	1:A:275:PHE:CE1	2.54	0.42
1:D:179:HIS:HA	1:D:183:ALA:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:137:ARG:CD	2:B:192:ILE:HD13	2.47	0.42
1:A:47:ARG:HE	1:A:184:LEU:HD23	1.85	0.42
2:C:88:SER:CB	1:D:107:SER:H	2.33	0.42
2:B:35:LEU:HD12	2:B:108:THR:HG21	2.02	0.42
2:C:148:PRO:HA	2:C:152:ILE:HB	2.01	0.42
1:A:156:GLN:HG3	1:A:156:GLN:O	2.20	0.42
2:C:28:VAL:O	2:C:31:PRO:HD2	2.20	0.42
2:B:1:MET:CG	2:B:2:GLN:N	2.77	0.41
1:D:4:PHE:HB2	1:D:8:MET:HE3	2.02	0.41
1:D:45:ARG:HB3	1:D:48:PRO:HG2	2.01	0.41
1:D:226:VAL:O	1:D:226:VAL:HG13	2.18	0.41
2:B:2:GLN:H	2:B:2:GLN:HG2	1.67	0.41
2:C:144:ARG:HG3	1:D:122:GLU:CD	2.45	0.41
1:D:256:LYS:HA	1:D:259:GLU:OE1	2.20	0.41
1:D:262:ASP:O	1:D:263:ARG:C	2.63	0.41
1:A:141:GLY:O	1:A:144:ALA:HB3	2.20	0.41
2:C:182:MET:O	2:C:185:CYS:HB3	2.20	0.41
1:A:88:MET:CE	1:A:160:VAL:HG21	2.48	0.41
2:B:1:MET:HE2	2:B:2:GLN:OE1	2.20	0.41
2:B:158:GLU:CD	2:B:178:LYS:HZ2	2.25	0.41
1:D:135:ARG:HD3	1:D:194:ILE:O	2.21	0.41
1:D:168:VAL:CG2	1:D:251:LEU:HD21	2.45	0.41
1:A:140:LEU:HD13	2:B:139:ILE:HG21	2.01	0.41
1:D:189:VAL:CG1	1:D:206:LEU:HB3	2.51	0.41
2:B:84:LEU:HD12	2:B:89:ARG:HB3	2.03	0.41
2:B:238:LEU:O	2:B:242:GLU:HG2	2.21	0.41
2:C:235:LEU:O	2:C:236:LYS:C	2.64	0.41
1:D:20:GLU:HG2	1:D:40:LEU:HD13	2.01	0.41
1:D:167:GLU:O	1:D:168:VAL:C	2.63	0.41
1:D:220:VAL:O	1:D:221:ASP:C	2.64	0.41
2:C:86:ASP:OD1	2:C:86:ASP:C	2.64	0.41
1:D:47:ARG:N	1:D:48:PRO:HD2	2.36	0.41
1:D:149:SER:C	1:D:154:ALA:HB2	2.46	0.41
1:D:156:GLN:O	1:D:159:ASP:HB3	2.21	0.41
1:D:293:ARG:HH21	1:D:293:ARG:HG3	1.86	0.41
2:B:5:TRP:CZ2	2:B:253:LEU:HG	2.56	0.41
2:B:138:VAL:HG22	2:B:189:CYS:SG	2.61	0.41
2:B:181:GLU:HB2	7:B:508:HOH:O	2.21	0.41
2:C:138:VAL:O	2:C:142:ILE:HG13	2.21	0.41
1:D:86:PRO:HG2	1:D:106:GLU:CD	2.46	0.41
1:D:155:GLY:O	1:D:159:ASP:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:2:GLN:HG2	7:C:476:HOH:O	2.21	0.41
2:C:92:SER:HB2	2:C:94:PRO:HD2	2.02	0.41
1:D:44:LYS:HB2	1:D:46:VAL:HG23	2.03	0.41
1:A:248:TYR:N	1:A:249:PRO:CD	2.84	0.40
2:C:139:ILE:HD13	1:D:140:LEU:HD13	2.03	0.40
2:C:148:PRO:HB3	1:D:115:ALA:HB1	2.04	0.40
2:C:59:ARG:O	2:C:60:SER:C	2.63	0.40
2:C:2:GLN:HB2	2:C:5:TRP:HD1	1.86	0.40
2:C:140:ILE:HD13	1:D:126:ALA:HB2	2.03	0.40
2:C:147:GLY:HA3	7:C:341:HOH:O	2.21	0.40
1:D:4:PHE:CE1	1:D:5:ASP:OD1	2.75	0.40
1:D:75:ILE:O	1:D:78:MET:HG3	2.21	0.40
1:D:269:GLN:HE22	1:D:288:ASN:ND2	2.19	0.40
2:B:25:PRO:C	2:B:27:THR:N	2.79	0.40
2:B:54:LEU:HD13	2:B:250:ASN:OD1	2.22	0.40
2:C:93:LYS:HA	2:C:93:LYS:HE2	2.03	0.40
1:D:143:LEU:HD13	1:D:191:LEU:HD11	2.03	0.40
1:D:214:GLY:O	1:D:216:LEU:N	2.55	0.40
2:B:49:LEU:HD22	2:B:59:ARG:HG3	2.04	0.40
2:B:214:LEU:HD11	2:B:232:ILE:HG23	2.03	0.40
2:B:215:ARG:NH2	7:B:325:HOH:O	2.50	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	275/295 (93%)	252 (92%)	22 (8%)	1 (0%)	30	42
1	D	280/295 (95%)	252 (90%)	25 (9%)	3 (1%)	11	16
2	B	252/274 (92%)	218 (86%)	30 (12%)	4 (2%)	7	10
2	C	257/274 (94%)	224 (87%)	22 (9%)	11 (4%)	2	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1064/1138 (94%)	946 (89%)	99 (9%)	19 (2%)	6 8

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	223	SER
2	B	227	ILE
2	C	94	PRO
2	C	196	ALA
2	C	198	GLU
2	C	238	LEU
1	D	165	MET
1	D	168	VAL
1	A	42	GLY
2	C	57	GLY
2	C	224	HIS
2	C	226	LEU
2	C	237	GLU
2	B	26	GLU
2	C	91	VAL
2	B	163	ASP
2	C	60	SER
2	C	163	ASP
1	D	286	LEU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	222/234 (95%)	215 (97%)	7 (3%)	34 54
1	D	226/234 (97%)	221 (98%)	5 (2%)	45 66
2	B	199/214 (93%)	195 (98%)	4 (2%)	48 68
2	C	201/214 (94%)	194 (96%)	7 (4%)	32 51
All	All	848/896 (95%)	825 (97%)	23 (3%)	39 60

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	LYS
1	A	99	THR
1	A	152	LEU
1	A	208	LYS
1	A	228	LYS
1	A	272	LEU
1	A	278	HIS
2	B	131	ASP
2	B	136	LEU
2	B	163	ASP
2	B	249	LYS
2	C	21	THR
2	C	93	LYS
2	C	94	PRO
2	C	131	ASP
2	C	139	ILE
2	C	182	MET
2	C	250	ASN
1	D	78	MET
1	D	102	MET
1	D	134	GLU
1	D	163	GLU
1	D	174	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	ASN
1	A	179	HIS
1	A	186	GLN
1	A	211	ASN
2	B	34	HIS
2	B	156	HIS
2	B	206	ASN
2	B	224	HIS
2	C	33	HIS
1	D	32	HIS
1	D	265	ASN
1	D	288	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	EDO	C	902	-	3,3,3	0.42	0	2,2,2	0.40	0
3	PPV	A	901	4	6,8,8	1.38	0	12,13,13	0.90	0
6	GPP	D	900	-	17,18,18	2.14	6 (35%)	21,25,25	1.25	4 (19%)

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
5	EDO	C	902	-	-	0/1/1/1	-
3	PPV	A	901	4	-	0/6/6/6	-
6	GPP	D	900	-	-	2/19/19/19	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	900	GPP	C6-C7	-3.66	1.39	1.50
6	D	900	GPP	PB-O1B	3.56	1.61	1.50
6	D	900	GPP	C1-C2	-3.42	1.39	1.49
6	D	900	GPP	PA-O1A	3.08	1.61	1.50
6	D	900	GPP	C2-C3	2.74	1.39	1.33
6	D	900	GPP	C7-C8	2.39	1.39	1.32

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	900	GPP	C4-C3-C5	2.55	119.66	115.23
6	D	900	GPP	C10-C8-C9	2.27	119.81	114.59
6	D	900	GPP	O3B-PB-O3A	2.22	112.08	104.64
6	D	900	GPP	O2A-PA-O3A	2.09	112.93	107.27

There are no chirality outliers.

All (2) torsion outliers are listed below:

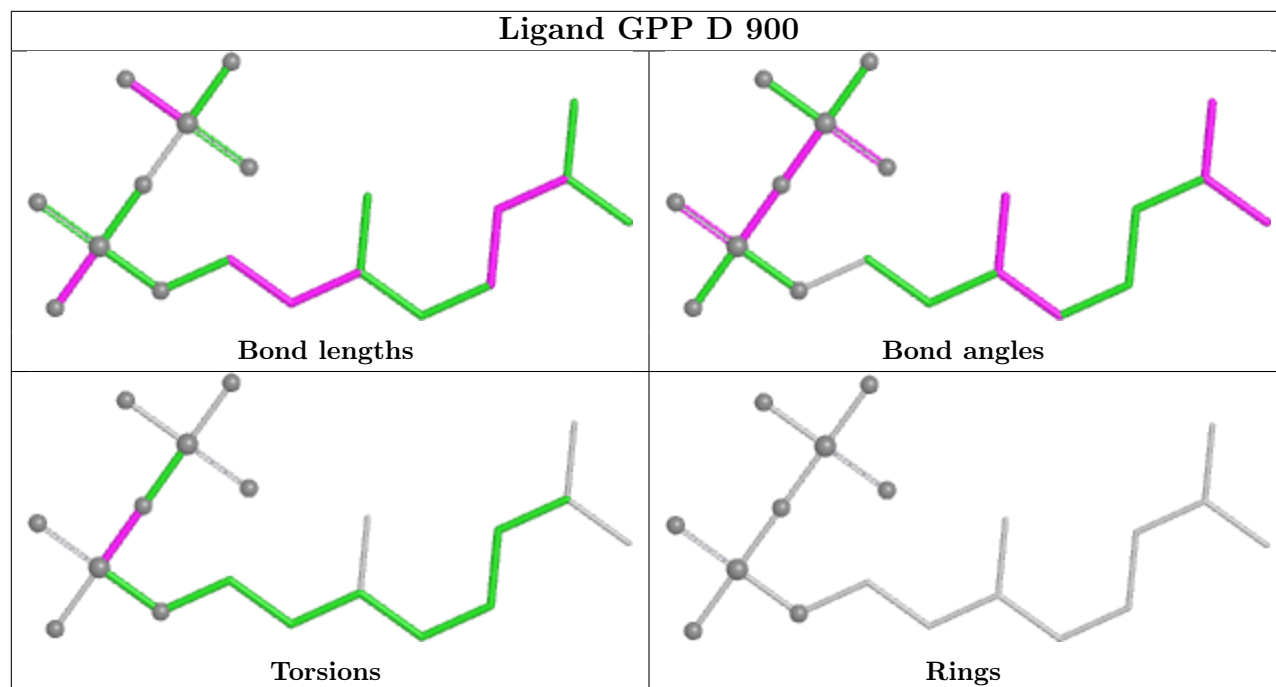
Mol	Chain	Res	Type	Atoms
6	D	900	GPP	PB-O3A-PA-O2A
6	D	900	GPP	PB-O3A-PA-O1A

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	902	EDO	1	0
6	D	900	GPP	2	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	279/295 (94%)	-0.31	1 (0%) 88 87	28, 42, 65, 95	0
1	D	284/295 (96%)	0.04	8 (2%) 55 51	30, 47, 109, 125	0
2	B	256/274 (93%)	0.14	14 (5%) 30 27	32, 50, 128, 145	0
2	C	259/274 (94%)	0.04	4 (1%) 72 69	30, 50, 97, 106	0
All	All	1078/1138 (94%)	-0.02	27 (2%) 58 55	28, 46, 104, 145	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	251	ALA	7.3
1	D	217	PHE	4.4
2	B	254	MET	4.1
2	B	250	ASN	3.6
1	D	223	ILE	3.3
2	B	259	ALA	3.3
2	B	256	SER	3.2
2	C	127	ALA	2.9
2	B	257	LEU	2.8
2	B	244	GLY	2.8
1	D	290	ILE	2.8
2	C	226	LEU	2.8
1	D	257	SER	2.6
2	C	259	ALA	2.5
2	B	231	ILE	2.4
1	D	248	TYR	2.3
2	B	226	LEU	2.3
2	B	227	ILE	2.3
1	D	254	VAL	2.3
1	D	250	LYS	2.3
2	B	238	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	248	GLY	2.2
1	A	166	ALA	2.2
2	B	235	LEU	2.2
2	C	227	ILE	2.1
2	B	255	SER	2.1
1	D	247	THR	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](#)

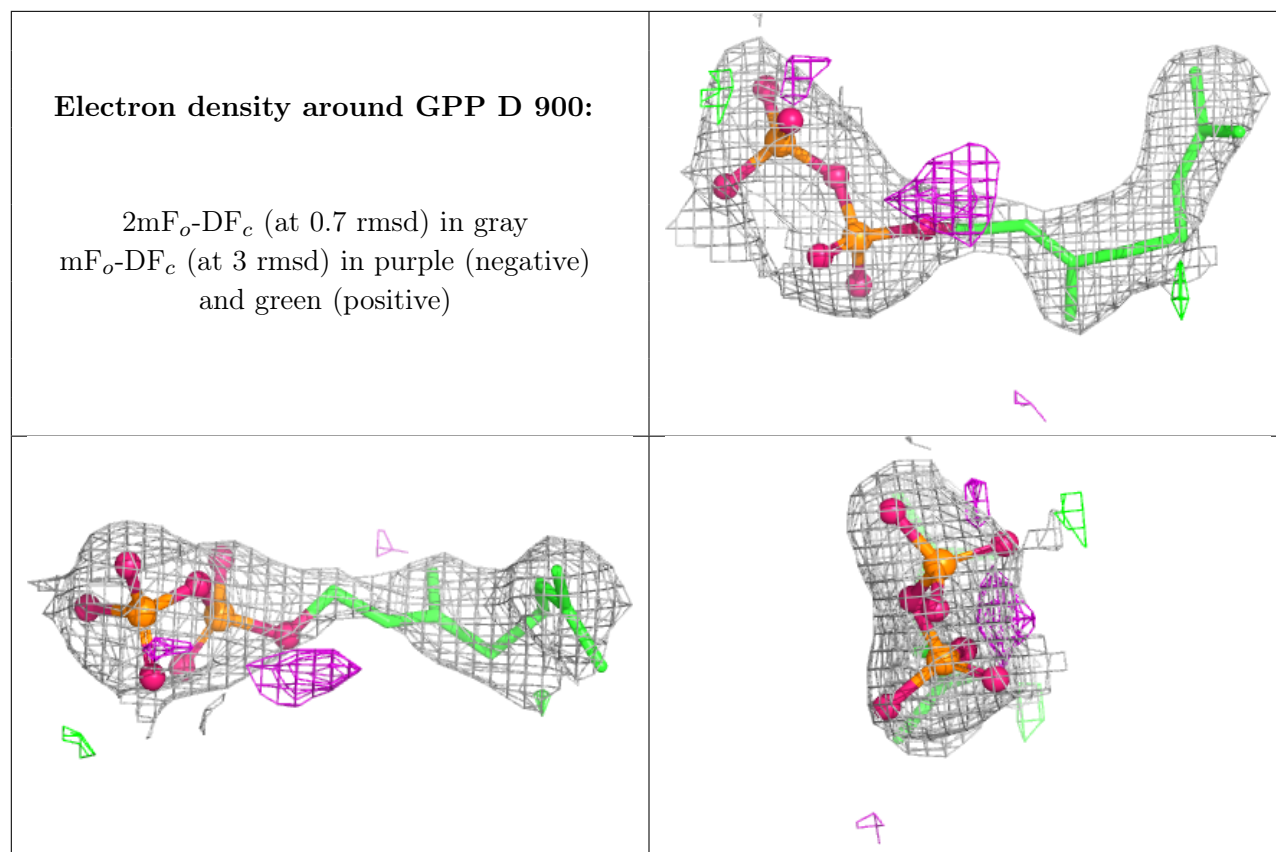
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PPV	A	901	9/9	0.80	0.09	83,84,86,87	0
4	MG	A	905	1/1	0.85	0.07	61,61,61,61	0
4	MG	D	904	1/1	0.87	0.05	57,57,57,57	0
6	GPP	D	900	19/19	0.89	0.11	53,61,68,69	0
4	MG	A	906	1/1	0.90	0.08	96,96,96,96	0
4	MG	D	903	1/1	0.93	0.06	48,48,48,48	0
5	EDO	C	902	4/4	0.95	0.05	35,37,38,38	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.