



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 05:54 PM JST

PDB ID : 7DE4 / pdb_00007de4
Title : Hybrid cluster protein (HCP) from Escherichia coli
Authors : Fujishiro, T.; Ooi, M.; Takaoka, K.; Takahashi, Y.
Deposited on : 2020-11-02
Resolution : 3.61 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 : 1.8.5 (274361), CSD as541be (2020)
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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

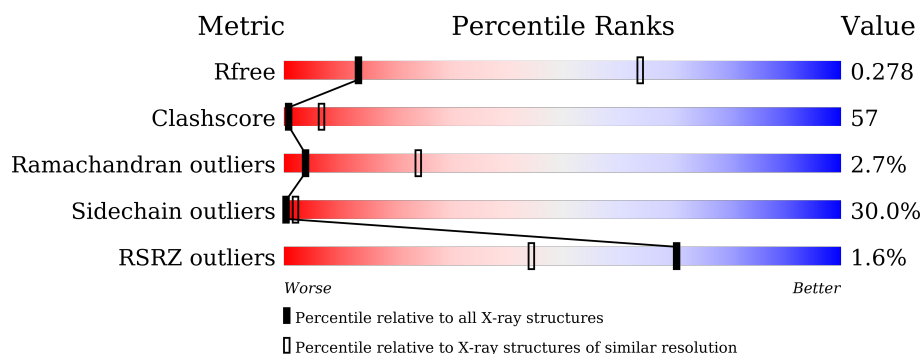
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.61 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	1062 (3.72-3.52)
Clashscore	190562	1092 (3.72-3.52)
Ramachandran outliers	187476	1057 (3.72-3.52)
Sidechain outliers	187428	1055 (3.72-3.52)
RSRZ outliers	180081	1060 (3.72-3.52)

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	558	2% 32% 45% 18% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	FSO	A	602	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4237 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.

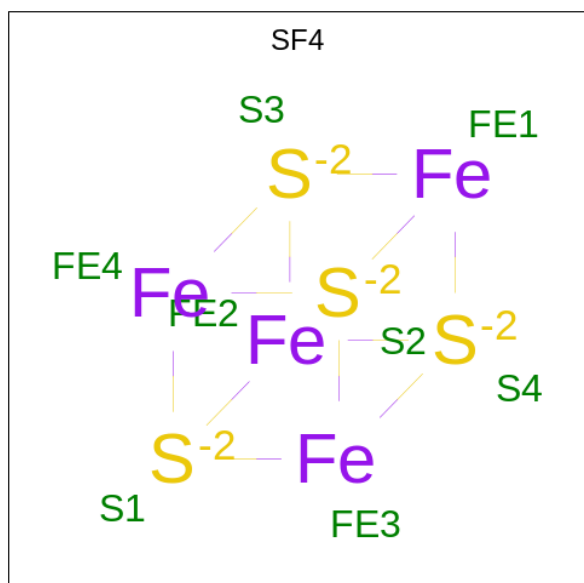
- Molecule 1 is a protein called Hydroxylamine reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	550	4219	2682	717	793	27	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

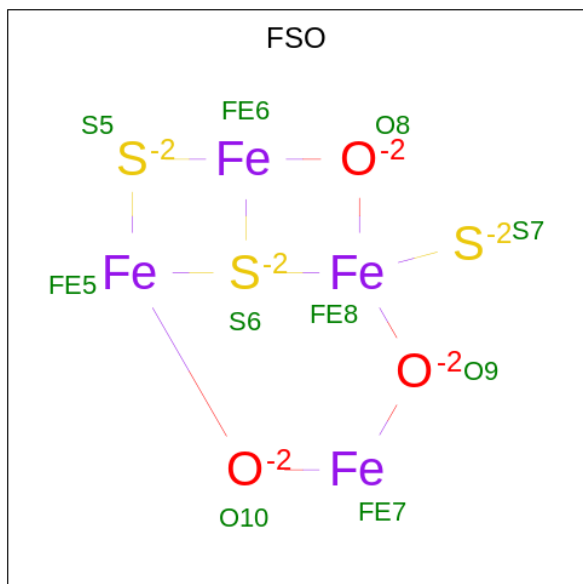
Chain	Residue	Modelled	Actual	Comment	Reference
A	551	LEU	-	expression tag	UNP P75825
A	552	GLU	-	expression tag	UNP P75825
A	553	HIS	-	expression tag	UNP P75825
A	554	HIS	-	expression tag	UNP P75825
A	555	HIS	-	expression tag	UNP P75825
A	556	HIS	-	expression tag	UNP P75825
A	557	HIS	-	expression tag	UNP P75825
A	558	HIS	-	expression tag	UNP P75825

- Molecule 2 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 3 is IRON/SULFUR/OXYGEN HYBRID CLUSTER (CCD ID: FSO) (formula: $\text{Fe}_4\text{O}_3\text{S}_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	Fe	O	S	0	0
			10	4	3	3		

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	142.54Å 142.54Å 117.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.29 – 3.61 46.29 – 3.61	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.29-3.61) 99.5 (46.29-3.61)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.203 , 0.278 0.203 , 0.278	Depositor DCC
R_{free} test set	717 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	116.3	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 154.8	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	4237	wwPDB-VP
Average B, all atoms (Å ²)	129.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% 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: SF4, FSO

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	0.69	1/4306 (0.0%)	1.68	77/5848 (1.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	271	HIS	CE1-NE2	5.10	1.37	1.32

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	299	ASN	CB-CA-C	-13.44	88.33	111.45
1	A	231	PRO	N-CA-C	-9.34	96.09	111.68
1	A	512	THR	CB-CA-C	-8.83	94.55	109.65
1	A	433	CYS	N-CA-CB	8.76	125.12	110.41
1	A	540	VAL	CB-CA-C	-8.69	100.99	111.94
1	A	204	SER	N-CA-C	-8.61	103.62	114.56
1	A	122	SER	CA-C-N	8.26	134.70	120.58
1	A	122	SER	C-N-CA	8.26	134.70	120.58
1	A	216	SER	CA-C-N	8.16	132.01	120.42
1	A	216	SER	C-N-CA	8.16	132.01	120.42
1	A	168	HIS	CB-CA-C	8.15	125.14	110.72
1	A	123	ASP	N-CA-C	8.01	121.54	111.69
1	A	523	ASP	CB-CA-C	-7.92	98.68	110.95
1	A	347	ASP	CA-CB-CG	-7.81	104.79	112.60
1	A	434	GLY	N-CA-C	-7.75	105.70	115.31
1	A	347	ASP	CB-CA-C	7.71	123.67	110.94
1	A	264	THR	CA-CB-OG1	-7.68	98.07	109.60
1	A	259	GLU	CB-CA-C	-7.64	98.89	110.88
1	A	433	CYS	CB-CA-C	-7.53	94.96	109.95
1	A	54	GLU	CB-CG-CD	7.37	125.13	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	THR	CB-CA-C	-7.30	97.81	110.45
1	A	194	PRO	N-CA-C	7.16	127.21	112.47
1	A	358	GLN	CB-CA-C	6.94	122.65	110.85
1	A	223	THR	CB-CA-C	-6.84	97.54	110.67
1	A	538	THR	CA-C-N	6.79	134.50	121.54
1	A	538	THR	C-N-CA	6.79	134.50	121.54
1	A	539	THR	CA-CB-OG1	-6.55	99.77	109.60
1	A	283	LEU	N-CA-C	-6.47	107.25	114.62
1	A	549	SER	CA-C-N	6.42	133.25	121.70
1	A	549	SER	C-N-CA	6.42	133.25	121.70
1	A	273	GLU	CB-CA-C	6.38	121.72	110.37
1	A	122	SER	N-CA-C	6.37	118.86	109.24
1	A	287	LYS	N-CA-C	-6.36	105.52	113.28
1	A	117	ASP	CB-CA-C	6.28	122.73	110.67
1	A	298	GLN	CB-CA-C	6.23	124.14	110.18
1	A	450	PRO	CB-CA-C	-6.16	103.02	111.21
1	A	217	ILE	N-CA-CB	-6.13	101.38	110.58
1	A	54	GLU	CB-CA-C	6.13	120.47	111.65
1	A	385	ASP	CB-CA-C	6.12	122.66	110.17
1	A	194	PRO	N-CA-CB	-6.01	96.94	103.25
1	A	53	ARG	CB-CA-C	5.96	120.24	110.88
1	A	357	GLN	CB-CG-CD	-5.92	102.53	112.60
1	A	494	LYS	CA-C-N	-5.90	110.81	121.14
1	A	494	LYS	C-N-CA	-5.90	110.81	121.14
1	A	399	ILE	N-CA-CB	-5.87	102.46	110.56
1	A	382	GLY	CA-C-N	5.70	128.97	120.31
1	A	382	GLY	C-N-CA	5.70	128.97	120.31
1	A	53	ARG	CA-C-N	-5.67	115.69	122.44
1	A	53	ARG	C-N-CA	-5.67	115.69	122.44
1	A	75	THR	CA-CB-OG1	-5.67	101.10	109.60
1	A	379	THR	CB-CA-C	5.63	121.16	109.95
1	A	417	ASP	CA-CB-CG	-5.62	106.98	112.60
1	A	239	THR	CB-CA-C	-5.61	100.26	109.80
1	A	185	LYS	CB-CG-CD	5.61	124.20	111.30
1	A	9	THR	CB-CA-C	-5.59	100.70	113.33
1	A	377	ARG	N-CA-C	-5.57	105.21	111.28
1	A	124	ASP	CA-C-N	5.55	132.14	121.54
1	A	124	ASP	C-N-CA	5.55	132.14	121.54
1	A	454	ASP	CA-CB-CG	5.47	118.07	112.60
1	A	434	GLY	CA-C-N	-5.47	112.41	120.28
1	A	434	GLY	C-N-CA	-5.47	112.41	120.28
1	A	27	LYS	CB-CA-C	5.46	119.59	109.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	ASP	CA-C-N	5.39	131.84	121.54
1	A	123	ASP	C-N-CA	5.39	131.84	121.54
1	A	512	THR	N-CA-CB	5.36	117.85	109.97
1	A	22	GLN	CB-CG-CD	-5.29	103.62	112.60
1	A	398	HIS	CB-CA-C	5.26	120.40	110.62
1	A	240	ALA	CB-CA-C	-5.19	110.61	116.63
1	A	80	ASP	CB-CA-C	-5.18	101.30	109.80
1	A	87	TYR	N-CA-C	-5.17	105.73	111.36
1	A	70	PHE	CB-CA-C	5.16	119.46	110.68
1	A	123	ASP	CB-CA-C	-5.15	101.28	110.70
1	A	454	ASP	CB-CA-C	5.10	118.06	109.89
1	A	28	THR	CA-CB-OG1	-5.09	101.96	109.60
1	A	236	VAL	CB-CA-C	-5.05	105.74	112.46
1	A	523	ASP	CA-CB-CG	5.02	117.62	112.60
1	A	512	THR	OG1-CB-CG2	5.01	119.33	109.30

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	4219	0	4179	476	0
2	A	8	0	0	0	0
3	A	10	0	0	6	0
All	All	4237	0	4179	476	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 57.

All (476) 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:148:ILE:CG2	1:A:195:ALA:HB1	1.67	1.24
1:A:232:THR:HG21	1:A:279:GLY:HA2	1.17	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LEU:HB3	1:A:128:LEU:HD11	1.17	1.16
1:A:154:LEU:HB2	1:A:519:PHE:HE2	1.00	1.13
1:A:281:PRO:HA	1:A:284:ARG:HD3	1.25	1.13
1:A:401:LEU:HD12	1:A:429:LEU:O	1.49	1.10
1:A:48:TRP:CH2	1:A:112:ASP:HB3	1.85	1.10
1:A:189:TRP:HH2	1:A:200:LEU:HD23	1.10	1.09
1:A:148:ILE:HG22	1:A:195:ALA:HB1	1.33	1.06
1:A:48:TRP:HH2	1:A:112:ASP:HB3	1.12	1.05
1:A:544:MET:HA	1:A:544:MET:HE3	1.38	1.05
1:A:51:LYS:HE3	1:A:51:LYS:HA	1.40	1.04
1:A:154:LEU:HB2	1:A:519:PHE:CE2	1.92	1.02
1:A:401:LEU:HD23	1:A:486:LEU:CD2	1.89	1.02
1:A:232:THR:HB	1:A:284:ARG:HH22	1.25	1.00
1:A:258:LEU:HD13	1:A:286:PHE:CD2	1.96	0.99
1:A:120:LEU:HD21	1:A:135:PHE:CZ	1.98	0.99
1:A:298:GLN:HG3	1:A:407:GLY:O	1.63	0.98
1:A:232:THR:CG2	1:A:279:GLY:HA2	1.94	0.98
1:A:125:LEU:HB3	1:A:128:LEU:CD1	1.93	0.98
1:A:176:ASP:HB3	1:A:179:ILE:HG13	1.46	0.97
1:A:148:ILE:HG21	1:A:195:ALA:HB1	1.47	0.97
1:A:120:LEU:HD21	1:A:135:PHE:HZ	1.28	0.97
1:A:203:CYS:O	1:A:203:CYS:SG	2.23	0.97
1:A:539:THR:CG2	1:A:542:GLU:HB2	1.95	0.97
1:A:5:GLN:HG3	1:A:493:GLN:NE2	1.81	0.96
1:A:99:LYS:C	1:A:103:LEU:HD12	1.91	0.95
1:A:223:THR:HG21	1:A:228:HIS:CE1	2.01	0.95
1:A:44:GLY:C	1:A:95:ARG:HH12	1.75	0.94
1:A:154:LEU:CB	1:A:519:PHE:HE2	1.79	0.94
1:A:189:TRP:CH2	1:A:200:LEU:HD23	2.01	0.94
1:A:151:LEU:CD1	1:A:524:LEU:HD22	1.98	0.93
1:A:239:THR:OG1	1:A:290:VAL:HG11	1.70	0.92
1:A:377:ARG:O	1:A:381:LEU:HD13	1.71	0.90
1:A:151:LEU:HD13	1:A:524:LEU:CD2	2.01	0.90
1:A:261:THR:O	1:A:264:THR:HG22	1.72	0.90
1:A:435:LYS:HD3	1:A:454:ASP:OD2	1.70	0.90
1:A:77:VAL:HG22	1:A:316:ASN:HD22	1.37	0.89
1:A:16:ASN:HB3	1:A:518:GLY:O	1.72	0.89
1:A:281:PRO:HA	1:A:284:ARG:CD	2.02	0.89
1:A:398:HIS:O	1:A:426:CYS:HB2	1.71	0.89
1:A:223:THR:CG2	1:A:228:HIS:CE1	2.57	0.88
1:A:258:LEU:HD13	1:A:286:PHE:HD2	1.34	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:LEU:HD21	1:A:62:VAL:HG13	1.56	0.88
1:A:151:LEU:CD1	1:A:524:LEU:CD2	2.52	0.88
1:A:125:LEU:CB	1:A:128:LEU:HD11	2.02	0.88
1:A:208:GLY:O	1:A:501:THR:HG22	1.73	0.88
1:A:213:LYS:O	1:A:216:SER:HB2	1.74	0.88
1:A:151:LEU:HD13	1:A:524:LEU:HD21	1.56	0.87
1:A:189:TRP:HH2	1:A:200:LEU:CD2	1.88	0.87
1:A:10:ILE:HG22	1:A:17:GLY:C	2.01	0.86
1:A:401:LEU:HD23	1:A:486:LEU:HD22	1.57	0.86
1:A:160:LYS:HD2	1:A:493:GLN:HG2	1.57	0.86
1:A:435:LYS:O	1:A:439:ASN:HB2	1.76	0.85
1:A:539:THR:HG21	1:A:542:GLU:HB2	1.60	0.84
1:A:264:THR:HG23	1:A:266:VAL:H	1.42	0.84
1:A:125:LEU:O	1:A:128:LEU:HD12	1.77	0.84
1:A:99:LYS:O	1:A:103:LEU:HD12	1.77	0.84
1:A:244:ILE:HG23	1:A:310:PRO:HG2	1.59	0.84
1:A:513:GLY:HA3	1:A:537:ILE:HG12	1.58	0.84
1:A:213:LYS:O	1:A:216:SER:CB	2.26	0.84
1:A:45:LEU:N	1:A:95:ARG:HH12	1.76	0.83
1:A:125:LEU:CB	1:A:128:LEU:CD1	2.57	0.83
1:A:257:LEU:O	1:A:261:THR:HG23	1.77	0.82
1:A:492:GLU:HG3	3:A:602:FSO:S7	2.18	0.82
1:A:48:TRP:CH2	1:A:112:ASP:CB	2.62	0.82
1:A:216:SER:HA	1:A:462:TYR:CE2	2.14	0.82
1:A:120:LEU:CD2	1:A:135:PHE:HZ	1.93	0.81
1:A:539:THR:HG23	1:A:542:GLU:HB2	1.60	0.81
1:A:423:PRO:HG2	1:A:426:CYS:SG	2.21	0.81
1:A:77:VAL:CG2	1:A:316:ASN:HD22	1.93	0.81
1:A:236:VAL:HG11	1:A:440:LYS:HE3	1.63	0.81
1:A:315:SER:HA	1:A:336:VAL:CG2	2.11	0.81
1:A:175:TYR:HD2	1:A:180:TYR:HE2	1.27	0.80
1:A:414:TYR:CE1	1:A:418:PHE:HD2	2.00	0.80
1:A:151:LEU:HD11	1:A:524:LEU:HD22	1.63	0.79
1:A:514:PRO:HA	1:A:537:ILE:HD11	1.65	0.79
1:A:189:TRP:CH2	1:A:200:LEU:CD2	2.65	0.79
1:A:47:ALA:O	1:A:50:VAL:HG12	1.83	0.79
1:A:77:VAL:CG2	1:A:316:ASN:ND2	2.46	0.78
1:A:281:PRO:CA	1:A:284:ARG:HD3	2.11	0.78
1:A:329:ARG:HD3	1:A:359:MET:CE	2.12	0.78
1:A:9:THR:HG21	1:A:17:GLY:HA2	1.65	0.78
1:A:264:THR:CG2	1:A:266:VAL:HG12	2.14	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:539:THR:HG21	1:A:542:GLU:CD	2.09	0.78
1:A:208:GLY:HA2	1:A:501:THR:HG22	1.66	0.77
1:A:211:ASN:HB3	1:A:501:THR:CG2	2.15	0.77
1:A:398:HIS:HB2	1:A:483:PRO:HB2	1.67	0.77
1:A:218:LEU:HD12	1:A:218:LEU:O	1.85	0.75
1:A:5:GLN:HG3	1:A:493:GLN:CD	2.12	0.75
1:A:115:MET:HE3	1:A:117:ASP:H	1.52	0.74
1:A:81:SER:N	1:A:82:PRO:HD2	2.02	0.74
1:A:289:LEU:HD23	1:A:289:LEU:O	1.88	0.74
1:A:332:THR:HB	1:A:336:VAL:HG12	1.68	0.74
1:A:9:THR:CG2	1:A:17:GLY:HA2	2.18	0.73
1:A:108:ASN:HB3	1:A:112:ASP:HA	1.70	0.73
1:A:182:GLN:OE1	1:A:210:MET:HB2	1.88	0.73
1:A:258:LEU:CD1	1:A:286:PHE:CD2	2.69	0.73
1:A:539:THR:HG21	1:A:542:GLU:CB	2.18	0.73
1:A:48:TRP:CZ2	1:A:112:ASP:HB2	2.24	0.72
1:A:154:LEU:HG	1:A:154:LEU:O	1.89	0.72
1:A:297:TRP:HB2	1:A:317:CYS:HB3	1.72	0.72
1:A:254:LEU:O	1:A:258:LEU:HG	1.90	0.71
1:A:89:ARG:HH11	1:A:89:ARG:CG	2.02	0.71
1:A:544:MET:HA	1:A:544:MET:CE	2.18	0.71
1:A:51:LYS:HB2	1:A:51:LYS:NZ	2.05	0.71
1:A:232:THR:CB	1:A:284:ARG:HH22	2.01	0.71
1:A:234:VAL:HG21	1:A:373:VAL:HG11	1.72	0.71
1:A:539:THR:HG21	1:A:542:GLU:CG	2.22	0.70
1:A:521:THR:OG1	1:A:523:ASP:HB2	1.90	0.70
1:A:175:TYR:CD2	1:A:180:TYR:HE2	2.09	0.70
1:A:259:GLU:O	1:A:262:GLU:HB2	1.91	0.70
1:A:16:ASN:CB	1:A:518:GLY:O	2.40	0.70
1:A:405:CYS:HB3	3:A:602:FSO:S6	2.31	0.70
1:A:401:LEU:CD1	1:A:429:LEU:O	2.36	0.70
1:A:58:ILE:HD12	1:A:58:ILE:H	1.57	0.69
1:A:45:LEU:HD21	1:A:62:VAL:CG1	2.22	0.69
1:A:48:TRP:O	1:A:51:LYS:HB3	1.92	0.69
1:A:155:CYS:SG	1:A:203:CYS:O	2.51	0.69
1:A:264:THR:HG21	1:A:266:VAL:HG12	1.74	0.69
1:A:48:TRP:CZ2	1:A:112:ASP:CB	2.76	0.68
1:A:76:ASN:O	1:A:318:ILE:HG22	1.94	0.68
1:A:401:LEU:CD2	1:A:486:LEU:CD2	2.68	0.68
1:A:273:GLU:OE1	1:A:458:CYS:N	2.26	0.68
1:A:315:SER:HA	1:A:336:VAL:HG23	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LEU:C	1:A:159:LEU:HD23	2.18	0.68
1:A:234:VAL:HG23	1:A:373:VAL:HG12	1.76	0.67
1:A:165:TYR:CE1	1:A:459:ASN:HB3	2.30	0.67
1:A:243:CYS:O	1:A:309:GLY:HA3	1.93	0.67
1:A:398:HIS:O	1:A:426:CYS:CB	2.43	0.67
1:A:51:LYS:HE3	1:A:51:LYS:CA	2.18	0.67
1:A:234:VAL:HG23	1:A:373:VAL:CG1	2.24	0.67
1:A:89:ARG:NH1	1:A:89:ARG:HG3	2.09	0.67
1:A:126:GLY:O	1:A:129:GLN:HB3	1.95	0.67
1:A:175:TYR:HD2	1:A:180:TYR:CE2	2.13	0.67
1:A:492:GLU:OE1	1:A:494:LYS:N	2.28	0.67
1:A:11:ARG:HH21	1:A:11:ARG:HG2	1.59	0.66
1:A:297:TRP:CB	1:A:317:CYS:HB3	2.25	0.66
1:A:387:LEU:HD21	1:A:427:LEU:HD11	1.77	0.66
1:A:208:GLY:CA	1:A:501:THR:HG22	2.26	0.66
1:A:211:ASN:HB3	1:A:501:THR:HG21	1.77	0.66
1:A:77:VAL:HG11	1:A:335:ILE:HD12	1.78	0.66
1:A:10:ILE:CG2	1:A:17:GLY:O	2.44	0.66
1:A:234:VAL:CG2	1:A:373:VAL:CG1	2.75	0.65
1:A:168:HIS:O	1:A:172:LEU:HD12	1.96	0.65
1:A:108:ASN:CB	1:A:112:ASP:HA	2.26	0.64
1:A:239:THR:CB	1:A:290:VAL:HG11	2.27	0.64
1:A:115:MET:HG3	1:A:118:LEU:H	1.63	0.64
1:A:310:PRO:HD3	1:A:356:ALA:HA	1.79	0.64
1:A:16:ASN:CG	1:A:518:GLY:O	2.41	0.64
1:A:151:LEU:HD22	1:A:197:MET:HG3	1.79	0.63
1:A:258:LEU:CD1	1:A:286:PHE:HD2	2.07	0.63
1:A:48:TRP:HZ2	1:A:112:ASP:HB2	1.62	0.63
1:A:272:GLY:HA2	1:A:436:TYR:HE2	1.63	0.63
1:A:223:THR:HG22	1:A:228:HIS:CE1	2.32	0.63
1:A:162:ALA:HB2	1:A:497:VAL:HG21	1.81	0.63
1:A:9:THR:HG22	1:A:18:CYS:N	2.15	0.62
1:A:230:THR:HB	1:A:231:PRO:CD	2.29	0.62
1:A:94:LEU:O	1:A:98:LEU:HB2	1.99	0.62
1:A:525:LEU:HD12	1:A:525:LEU:C	2.24	0.62
1:A:10:ILE:CG2	1:A:17:GLY:C	2.73	0.62
1:A:182:GLN:O	1:A:186:ILE:HG13	2.00	0.62
1:A:264:THR:HG23	1:A:266:VAL:HG12	1.82	0.61
1:A:275:LEU:N	1:A:276:PRO:CD	2.62	0.61
1:A:169:ALA:HB2	1:A:218:LEU:HB2	1.82	0.61
1:A:208:GLY:C	1:A:501:THR:HG22	2.24	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:LEU:HD11	1:A:524:LEU:CD2	2.25	0.61
1:A:480:ASN:OD1	1:A:480:ASN:N	2.32	0.61
1:A:204:SER:O	1:A:500:LEU:HD13	2.01	0.61
1:A:381:LEU:N	1:A:381:LEU:HD12	2.14	0.61
1:A:151:LEU:CD1	1:A:524:LEU:HD21	2.25	0.61
1:A:120:LEU:CD2	1:A:135:PHE:CZ	2.73	0.61
1:A:432:ALA:O	1:A:435:LYS:HB3	2.01	0.61
1:A:89:ARG:CG	1:A:89:ARG:NH1	2.60	0.60
1:A:223:THR:HG22	1:A:228:HIS:ND1	2.16	0.60
1:A:470:THR:O	1:A:474:LYS:HG3	2.00	0.60
1:A:492:GLU:CD	3:A:602:FSO:S7	2.84	0.60
1:A:152:ARG:NH2	1:A:193:TRP:HB3	2.16	0.60
1:A:414:TYR:HE1	1:A:418:PHE:CD2	2.20	0.59
1:A:401:LEU:HB3	1:A:486:LEU:HD23	1.85	0.59
1:A:275:LEU:N	1:A:276:PRO:HD3	2.17	0.59
1:A:304:PHE:HB2	1:A:326:TYR:CE2	2.36	0.59
1:A:9:THR:HG21	1:A:17:GLY:CA	2.32	0.59
1:A:248:GLY:HA3	1:A:314:THR:OG1	2.02	0.59
1:A:125:LEU:C	1:A:128:LEU:HD12	2.27	0.59
1:A:10:ILE:HG22	1:A:18:CYS:N	2.18	0.59
1:A:218:LEU:HD12	1:A:218:LEU:C	2.28	0.59
1:A:5:GLN:CG	1:A:493:GLN:NE2	2.60	0.59
1:A:148:ILE:HG21	1:A:195:ALA:CB	2.28	0.58
1:A:186:ILE:HG21	1:A:207:ILE:HD11	1.85	0.58
1:A:414:TYR:CE1	1:A:418:PHE:CD2	2.87	0.58
1:A:400:PHE:CZ	1:A:544:MET:HE2	2.39	0.58
1:A:321:PRO:HG3	1:A:341:VAL:HG21	1.85	0.58
1:A:39:ILE:O	1:A:43:GLN:HG3	2.03	0.58
1:A:329:ARG:HD3	1:A:359:MET:HE2	1.85	0.58
1:A:329:ARG:HD3	1:A:359:MET:HE1	1.85	0.57
1:A:331:TRP:CZ3	1:A:342:ARG:HD2	2.38	0.57
1:A:38:LEU:CD1	1:A:42:LEU:HD23	2.35	0.57
1:A:60:HIS:CE1	1:A:175:TYR:CE2	2.92	0.57
1:A:217:ILE:O	1:A:220:ALA:HB3	2.04	0.57
1:A:315:SER:HA	1:A:336:VAL:HG21	1.86	0.57
1:A:5:GLN:O	1:A:493:GLN:NE2	2.37	0.57
1:A:234:VAL:CG2	1:A:373:VAL:HG11	2.33	0.57
1:A:38:LEU:HD11	1:A:42:LEU:HD23	1.85	0.56
1:A:77:VAL:HG23	1:A:316:ASN:ND2	2.19	0.56
1:A:4:VAL:HG22	1:A:4:VAL:O	2.05	0.56
1:A:49:ALA:C	1:A:51:LYS:H	2.13	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:ILE:HG22	1:A:195:ALA:CB	2.23	0.56
1:A:208:GLY:O	1:A:501:THR:CG2	2.52	0.56
1:A:338:TRP:HD1	1:A:341:VAL:HG23	1.70	0.56
1:A:405:CYS:CB	3:A:602:FSO:S6	2.92	0.56
1:A:47:ALA:HB1	1:A:113:ASN:HD21	1.71	0.56
1:A:226:TYR:OH	1:A:252:LYS:HA	2.05	0.56
1:A:480:ASN:ND2	1:A:508:LYS:H	2.03	0.56
1:A:381:LEU:H	1:A:381:LEU:CD1	2.18	0.56
1:A:176:ASP:OD1	1:A:177:ASN:N	2.39	0.56
1:A:76:ASN:HA	1:A:79:PHE:CZ	2.41	0.55
1:A:321:PRO:HB3	1:A:326:TYR:HE1	1.71	0.55
1:A:134:GLU:O	1:A:134:GLU:HG3	2.05	0.55
1:A:148:ILE:CD1	1:A:197:MET:HB2	2.37	0.55
1:A:401:LEU:CB	1:A:486:LEU:HD23	2.36	0.55
1:A:313:MET:HE3	1:A:318:ILE:HG13	1.89	0.55
1:A:77:VAL:CG1	1:A:335:ILE:HD12	2.36	0.55
1:A:239:THR:OG1	1:A:290:VAL:CG1	2.51	0.55
1:A:60:HIS:O	1:A:64:SER:HB3	2.06	0.55
1:A:51:LYS:CA	1:A:51:LYS:CE	2.85	0.54
1:A:414:TYR:HE1	1:A:418:PHE:HD2	1.47	0.54
1:A:92:ILE:HG23	1:A:95:ARG:HE	1.72	0.54
1:A:42:LEU:HD21	1:A:69:ALA:HB3	1.90	0.54
1:A:193:TRP:N	1:A:194:PRO:HD2	2.22	0.54
1:A:8:GLN:OE1	1:A:408:ALA:N	2.40	0.54
1:A:44:GLY:C	1:A:95:ARG:NH1	2.57	0.54
1:A:147:ASN:O	1:A:524:LEU:HD11	2.08	0.54
1:A:81:SER:N	1:A:82:PRO:CD	2.71	0.54
1:A:363:PRO:HD2	1:A:364:TYR:H	1.73	0.54
1:A:517:PRO:HB2	1:A:520:LEU:HD23	1.89	0.54
1:A:9:THR:CG2	1:A:18:CYS:H	2.21	0.54
1:A:9:THR:HG22	1:A:10:ILE:H	1.73	0.54
1:A:45:LEU:N	1:A:95:ARG:NH1	2.53	0.54
1:A:95:ARG:HG3	1:A:95:ARG:HH11	1.72	0.54
1:A:154:LEU:CB	1:A:519:PHE:CE2	2.69	0.54
1:A:7:GLU:OE2	1:A:409:ARG:NH1	2.41	0.54
1:A:418:PHE:CZ	1:A:487:VAL:HG21	2.43	0.53
1:A:418:PHE:CE1	1:A:544:MET:HE1	2.43	0.53
1:A:51:LYS:HE2	1:A:53:ARG:HG3	1.89	0.53
1:A:412:ARG:O	1:A:413:HIS:C	2.52	0.53
1:A:213:LYS:O	1:A:216:SER:HB3	2.05	0.53
1:A:402:LEU:CD1	1:A:419:ALA:HB2	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:LEU:H	1:A:307:PHE:HE2	1.56	0.53
1:A:296:GLY:N	1:A:436:TYR:CD2	2.76	0.53
1:A:320:ASP:HA	1:A:338:TRP:CZ2	2.44	0.53
1:A:295:SER:OG	1:A:295:SER:O	2.24	0.53
1:A:103:LEU:CD2	1:A:108:ASN:OD1	2.57	0.53
1:A:245:LEU:O	1:A:311:ILE:HA	2.09	0.53
1:A:76:ASN:HA	1:A:79:PHE:HZ	1.74	0.53
1:A:297:TRP:O	1:A:300:GLN:HG2	2.09	0.53
1:A:51:LYS:HB2	1:A:51:LYS:HZ2	1.74	0.52
1:A:230:THR:HB	1:A:231:PRO:HD3	1.91	0.52
1:A:370:LEU:HD12	1:A:370:LEU:H	1.73	0.52
1:A:381:LEU:N	1:A:381:LEU:CD1	2.71	0.52
1:A:175:TYR:HB2	1:A:180:TYR:CE2	2.44	0.52
1:A:405:CYS:SG	3:A:602:FSO:S6	3.07	0.52
1:A:508:LYS:HG2	1:A:532:PHE:HA	1.92	0.52
1:A:88:ALA:O	1:A:92:ILE:HG13	2.09	0.52
1:A:391:VAL:CG2	1:A:396:LEU:HB3	2.39	0.52
1:A:513:GLY:HA3	1:A:537:ILE:CG1	2.34	0.52
1:A:60:HIS:ND1	1:A:175:TYR:CE2	2.78	0.52
1:A:321:PRO:CB	1:A:326:TYR:HE1	2.22	0.52
1:A:398:HIS:H	1:A:426:CYS:HB3	1.74	0.52
1:A:375:PHE:HB2	1:A:454:ASP:O	2.10	0.52
1:A:503:LEU:HD22	1:A:532:PHE:CG	2.45	0.52
1:A:225:LYS:HB3	1:A:226:TYR:CD2	2.45	0.51
1:A:9:THR:HG22	1:A:18:CYS:H	1.75	0.51
1:A:147:ASN:ND2	1:A:523:ASP:CB	2.72	0.51
1:A:369:HIS:C	1:A:369:HIS:CD2	2.88	0.51
1:A:457:GLN:O	1:A:458:CYS:C	2.53	0.51
1:A:42:LEU:HD21	1:A:69:ALA:CB	2.41	0.51
1:A:99:LYS:C	1:A:103:LEU:CD1	2.77	0.51
1:A:261:THR:C	1:A:264:THR:HG22	2.35	0.51
1:A:453:VAL:HG12	1:A:453:VAL:O	2.10	0.51
1:A:537:ILE:HD12	1:A:538:THR:H	1.76	0.51
1:A:391:VAL:HG23	1:A:396:LEU:HB3	1.92	0.51
1:A:110:ARG:HD2	1:A:111:VAL:H	1.75	0.51
1:A:493:GLN:CD	1:A:493:GLN:N	2.69	0.51
1:A:245:LEU:N	1:A:307:PHE:HE2	2.07	0.51
1:A:77:VAL:HG22	1:A:316:ASN:ND2	2.13	0.50
1:A:539:THR:HG23	1:A:539:THR:O	2.09	0.50
1:A:45:LEU:CD2	1:A:62:VAL:HG13	2.36	0.50
1:A:327:ASP:O	1:A:341:VAL:HG13	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:GLU:CG	3:A:602:FSO:S7	2.90	0.50
1:A:396:LEU:HD23	1:A:426:CYS:HA	1.92	0.50
1:A:462:TYR:CE1	1:A:466:ILE:HD13	2.47	0.50
1:A:544:MET:HG3	1:A:548:LEU:CD2	2.41	0.50
1:A:321:PRO:CG	1:A:341:VAL:HG21	2.42	0.50
1:A:514:PRO:HA	1:A:537:ILE:CD1	2.41	0.50
1:A:245:LEU:N	1:A:307:PHE:CE2	2.80	0.50
1:A:147:ASN:ND2	1:A:523:ASP:HB2	2.27	0.49
1:A:52:ALA:HA	1:A:56:GLY:HA3	1.95	0.49
1:A:70:PHE:CE2	1:A:160:LYS:HA	2.48	0.49
1:A:326:TYR:HD1	1:A:326:TYR:C	2.20	0.49
1:A:147:ASN:HD21	1:A:523:ASP:CB	2.26	0.49
1:A:247:SER:HB2	1:A:271:HIS:ND1	2.28	0.49
1:A:181:ALA:O	1:A:185:LYS:HG3	2.12	0.49
1:A:192:THR:O	1:A:192:THR:OG1	2.30	0.49
1:A:49:ALA:C	1:A:51:LYS:N	2.71	0.49
1:A:166:MET:O	1:A:169:ALA:HB3	2.13	0.49
1:A:297:TRP:CG	1:A:317:CYS:HB3	2.47	0.49
1:A:414:TYR:CD1	1:A:414:TYR:C	2.90	0.49
1:A:44:GLY:O	1:A:45:LEU:C	2.55	0.48
1:A:505:LEU:O	1:A:505:LEU:HD22	2.13	0.48
1:A:34:LEU:HD12	1:A:81:SER:HB3	1.95	0.48
1:A:45:LEU:O	1:A:45:LEU:HD23	2.13	0.48
1:A:166:MET:CG	1:A:214:VAL:HG21	2.44	0.48
1:A:33:ASP:HB3	1:A:130:ARG:HG3	1.96	0.48
1:A:538:THR:O	1:A:539:THR:HG22	2.14	0.48
1:A:103:LEU:HD21	1:A:112:ASP:OD2	2.13	0.48
1:A:258:LEU:HD23	1:A:268:VAL:HG21	1.96	0.48
1:A:321:PRO:HD2	1:A:338:TRP:CE2	2.49	0.48
1:A:387:LEU:HD21	1:A:427:LEU:CD1	2.43	0.48
1:A:102:CYS:O	1:A:105:VAL:HG23	2.13	0.48
1:A:409:ARG:HD2	1:A:514:PRO:HG2	1.96	0.48
1:A:208:GLY:HA2	1:A:501:THR:CG2	2.40	0.48
1:A:1:MET:O	1:A:19:SER:HA	2.14	0.48
1:A:103:LEU:C	1:A:105:VAL:H	2.21	0.48
1:A:527:VAL:O	1:A:530:GLU:HB3	2.13	0.48
1:A:454:ASP:OD1	1:A:456:GLY:N	2.40	0.47
1:A:108:ASN:OD1	1:A:112:ASP:OD2	2.32	0.47
1:A:186:ILE:HD13	1:A:207:ILE:HG13	1.95	0.47
1:A:326:TYR:C	1:A:326:TYR:CD1	2.92	0.47
1:A:225:LYS:HB3	1:A:226:TYR:CE2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:PHE:CZ	1:A:544:MET:CE	2.97	0.47
1:A:70:PHE:CD1	1:A:187:MET:HE1	2.49	0.47
1:A:166:MET:HG3	1:A:214:VAL:HG21	1.96	0.47
1:A:321:PRO:HA	1:A:326:TYR:CE1	2.48	0.47
1:A:407:GLY:H	1:A:412:ARG:HD3	1.79	0.47
1:A:422:VAL:HG12	1:A:451:ARG:HH21	1.80	0.47
1:A:540:VAL:HG12	1:A:540:VAL:O	2.14	0.47
1:A:9:THR:CG2	1:A:17:GLY:CA	2.92	0.47
1:A:242:LYS:C	1:A:266:VAL:HG23	2.40	0.47
1:A:363:PRO:CD	1:A:364:TYR:H	2.28	0.47
1:A:418:PHE:CZ	1:A:544:MET:HE1	2.50	0.47
1:A:120:LEU:HD21	1:A:135:PHE:CE2	2.49	0.47
1:A:407:GLY:HA3	1:A:412:ARG:CZ	2.43	0.47
1:A:446:ILE:O	1:A:447:GLU:HB2	2.15	0.47
1:A:48:TRP:O	1:A:51:LYS:CB	2.62	0.47
1:A:232:THR:CG2	1:A:279:GLY:CA	2.82	0.47
1:A:285:LYS:HE3	1:A:286:PHE:CZ	2.50	0.47
1:A:380:LEU:O	1:A:383:ALA:HB3	2.15	0.47
1:A:381:LEU:HD12	1:A:381:LEU:H	1.79	0.47
1:A:129:GLN:O	1:A:131:GLN:N	2.48	0.46
1:A:261:THR:HB	1:A:264:THR:HG21	1.96	0.46
1:A:99:LYS:O	1:A:103:LEU:CD1	2.55	0.46
1:A:4:VAL:HB	1:A:17:GLY:HA3	1.96	0.46
1:A:114:PRO:O	1:A:114:PRO:HG2	2.15	0.46
1:A:433:CYS:C	1:A:435:LYS:N	2.70	0.46
1:A:480:ASN:HD22	1:A:508:LYS:H	1.63	0.46
1:A:10:ILE:CG2	1:A:18:CYS:HA	2.45	0.46
1:A:152:ARG:HH21	1:A:193:TRP:HE3	1.64	0.46
1:A:293:TYR:C	1:A:293:TYR:CD1	2.94	0.46
1:A:167:GLU:C	1:A:169:ALA:H	2.22	0.46
1:A:268:VAL:HG12	1:A:289:LEU:HB2	1.98	0.46
1:A:4:VAL:CG2	1:A:518:GLY:HA3	2.46	0.46
1:A:300:GLN:HG2	1:A:300:GLN:H	1.50	0.46
1:A:103:LEU:HD23	1:A:108:ASN:OD1	2.16	0.45
1:A:275:LEU:HD22	1:A:373:VAL:HG21	1.98	0.45
1:A:388:ILE:O	1:A:392:SER:HB2	2.16	0.45
1:A:99:LYS:CG	1:A:103:LEU:HD11	2.45	0.45
1:A:245:LEU:HB3	1:A:307:PHE:CZ	2.52	0.45
1:A:380:LEU:HD23	1:A:380:LEU:HA	1.78	0.45
1:A:432:ALA:O	1:A:435:LYS:HE2	2.16	0.45
1:A:239:THR:HG1	1:A:290:VAL:HG11	1.78	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:PRO:HD2	1:A:338:TRP:NE1	2.32	0.45
1:A:103:LEU:C	1:A:105:VAL:N	2.72	0.45
1:A:218:LEU:O	1:A:222:GLU:HG3	2.16	0.45
1:A:45:LEU:C	1:A:45:LEU:HD23	2.41	0.45
1:A:283:LEU:HD12	1:A:283:LEU:HA	1.79	0.45
1:A:92:ILE:O	1:A:96:GLU:HG2	2.17	0.45
1:A:521:THR:HG23	1:A:524:LEU:HD12	1.98	0.45
1:A:125:LEU:HB2	1:A:128:LEU:CD1	2.44	0.45
1:A:433:CYS:C	1:A:435:LYS:H	2.24	0.45
1:A:212:PHE:O	1:A:213:LYS:C	2.56	0.45
1:A:412:ARG:HG3	1:A:514:PRO:HG3	1.99	0.45
1:A:88:ALA:O	1:A:89:ARG:C	2.61	0.44
1:A:301:GLN:OE1	1:A:319:ILE:HD12	2.16	0.44
1:A:477:CYS:SG	1:A:478:GLY:N	2.90	0.44
1:A:38:LEU:O	1:A:42:LEU:HB2	2.17	0.44
1:A:154:LEU:CA	1:A:519:PHE:CE2	3.01	0.44
1:A:77:VAL:HG12	1:A:78:ASN:HB2	1.98	0.44
1:A:151:LEU:HD21	1:A:201:LEU:HD23	2.00	0.44
1:A:27:LYS:HE2	1:A:35:GLN:OE1	2.17	0.44
1:A:95:ARG:HH11	1:A:95:ARG:CG	2.30	0.44
1:A:226:TYR:HA	1:A:282:GLU:HB3	2.00	0.44
1:A:462:TYR:CD1	1:A:462:TYR:C	2.96	0.44
1:A:66:ALA:N	1:A:67:PRO:HD2	2.33	0.44
1:A:113:ASN:OD1	1:A:114:PRO:HD2	2.18	0.44
1:A:254:LEU:HD11	1:A:268:VAL:HG11	2.00	0.44
1:A:38:LEU:HD11	1:A:42:LEU:CD2	2.48	0.44
1:A:227:GLY:HA3	1:A:281:PRO:HG2	1.99	0.44
1:A:282:GLU:O	1:A:282:GLU:HG3	2.17	0.44
1:A:427:LEU:HD21	1:A:449:LEU:HD13	1.99	0.44
1:A:431:LEU:HD13	1:A:431:LEU:O	2.17	0.44
1:A:519:PHE:CD1	1:A:519:PHE:N	2.84	0.44
1:A:74:LEU:HD12	1:A:160:LYS:HG2	1.99	0.43
1:A:175:TYR:CD1	1:A:175:TYR:N	2.86	0.43
1:A:4:VAL:O	1:A:4:VAL:CG2	2.66	0.43
1:A:252:LYS:HE2	1:A:252:LYS:HB2	1.52	0.43
1:A:266:VAL:HG11	1:A:353:ILE:HG23	2.00	0.43
1:A:298:GLN:CG	1:A:407:GLY:O	2.51	0.43
1:A:300:GLN:HG3	1:A:319:ILE:CG1	2.48	0.43
1:A:321:PRO:CB	1:A:326:TYR:CE1	3.01	0.43
1:A:7:GLU:CD	1:A:7:GLU:O	2.62	0.43
1:A:11:ARG:HG2	1:A:11:ARG:NH2	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:GLU:O	1:A:94:LEU:HD12	2.17	0.43
1:A:167:GLU:C	1:A:169:ALA:N	2.77	0.43
1:A:243:CYS:HB2	1:A:307:PHE:CZ	2.53	0.43
1:A:83:ARG:NH2	1:A:343:HIS:HD2	2.17	0.43
1:A:163:ALA:O	1:A:164:ALA:C	2.62	0.43
1:A:297:TRP:CE3	1:A:297:TRP:C	2.97	0.43
1:A:56:GLY:O	1:A:57:ILE:O	2.37	0.43
1:A:122:SER:OG	1:A:131:GLN:NE2	2.52	0.43
1:A:431:LEU:O	1:A:432:ALA:HB2	2.19	0.43
1:A:57:ILE:HG23	1:A:58:ILE:N	2.33	0.43
1:A:471:LEU:HD23	1:A:471:LEU:HA	1.89	0.43
1:A:2:PHE:CD1	1:A:2:PHE:C	2.97	0.42
1:A:245:LEU:CD2	1:A:293:TYR:CD1	3.02	0.42
1:A:457:GLN:O	1:A:460:ASP:HB2	2.19	0.42
1:A:39:ILE:HG13	1:A:73:THR:HG21	2.01	0.42
1:A:152:ARG:CA	1:A:200:LEU:HD13	2.49	0.42
1:A:462:TYR:CE1	1:A:466:ILE:CD1	3.01	0.42
1:A:37:LEU:HD21	1:A:120:LEU:HD11	2.01	0.42
1:A:244:ILE:HD12	1:A:257:LEU:HD11	2.02	0.42
1:A:256:ASN:C	1:A:258:LEU:N	2.76	0.42
1:A:275:LEU:HB3	1:A:276:PRO:HD3	2.02	0.42
1:A:27:LYS:HD3	1:A:32:SER:OG	2.19	0.42
1:A:234:VAL:O	1:A:370:LEU:HA	2.20	0.42
1:A:81:SER:H	1:A:82:PRO:HD2	1.79	0.42
1:A:435:LYS:HB3	1:A:435:LYS:HE2	1.67	0.42
1:A:175:TYR:N	1:A:175:TYR:HD1	2.18	0.42
1:A:211:ASN:HD22	1:A:501:THR:CG2	2.33	0.42
1:A:236:VAL:HG23	1:A:369:HIS:O	2.19	0.42
1:A:276:PRO:O	1:A:277:ALA:C	2.60	0.42
1:A:402:LEU:HD23	1:A:402:LEU:HA	1.92	0.42
1:A:182:GLN:CD	1:A:210:MET:HB2	2.45	0.41
1:A:344:LEU:HD21	1:A:348:ASP:O	2.20	0.41
1:A:28:THR:HB	1:A:30:GLU:H	1.85	0.41
1:A:117:ASP:O	1:A:120:LEU:HB2	2.20	0.41
1:A:230:THR:CB	1:A:231:PRO:HD3	2.50	0.41
1:A:311:ILE:HB	1:A:330:ILE:HG12	2.03	0.41
1:A:355:GLN:O	1:A:355:GLN:HG2	2.20	0.41
1:A:136:THR:O	1:A:193:TRP:CH2	2.73	0.41
1:A:201:LEU:HD22	1:A:201:LEU:HA	1.92	0.41
1:A:449:LEU:HD23	1:A:449:LEU:HA	1.74	0.41
1:A:136:THR:O	1:A:193:TRP:HH2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:THR:HG23	1:A:266:VAL:N	2.22	0.41
1:A:287:LYS:HE2	1:A:287:LYS:HB3	1.56	0.41
1:A:414:TYR:CZ	1:A:537:ILE:HG22	2.56	0.41
1:A:517:PRO:HB3	1:A:519:PHE:CE1	2.55	0.41
1:A:390:LEU:HB3	1:A:396:LEU:HB2	2.02	0.41
1:A:37:LEU:HB2	1:A:132:ALA:HB2	2.03	0.41
1:A:387:LEU:HD22	1:A:396:LEU:CD1	2.51	0.41
1:A:43:GLN:CD	1:A:191:GLY:HA2	2.46	0.41
1:A:493:GLN:CD	1:A:493:GLN:H	2.28	0.41
1:A:321:PRO:CA	1:A:326:TYR:CE1	3.04	0.41
1:A:70:PHE:CE1	1:A:187:MET:HE1	2.57	0.40
1:A:537:ILE:H	1:A:537:ILE:HG13	1.58	0.40
1:A:115:MET:CE	1:A:117:ASP:H	2.26	0.40
1:A:232:THR:HB	1:A:284:ARG:NH2	2.09	0.40
1:A:258:LEU:HB3	1:A:286:PHE:CD2	2.57	0.40
1:A:138:ASN:HB3	1:A:149:LEU:HD11	2.04	0.40
1:A:159:LEU:HD23	1:A:159:LEU:O	2.21	0.40
1:A:226:TYR:O	1:A:281:PRO:HD2	2.22	0.40
1:A:245:LEU:HB3	1:A:307:PHE:CE2	2.56	0.40
1:A:261:THR:OG1	1:A:288:HIS:CD2	2.74	0.40
1:A:543:ASP:O	1:A:547:LEU:HG	2.21	0.40
1:A:151:LEU:HD22	1:A:197:MET:CG	2.47	0.40
1:A:192:THR:C	1:A:194:PRO:HD2	2.47	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	548/558 (98%)	479 (87%)	54 (10%)	15 (3%)	4 24

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	ILE
1	A	124	ASP
1	A	130	ARG
1	A	194	PRO
1	A	539	THR
1	A	315	SER
1	A	509	ASN
1	A	537	ILE
1	A	11	ARG
1	A	125	LEU
1	A	4	VAL
1	A	431	LEU
1	A	538	THR
1	A	363	PRO
1	A	476	GLY

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	446/454 (98%)	312 (70%)	134 (30%)	0 2

All (134) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	5	GLN
1	A	9	THR
1	A	11	ARG
1	A	12	THR
1	A	27	LYS
1	A	28	THR
1	A	32	SER
1	A	37	LEU
1	A	42	LEU
1	A	50	VAL
1	A	51	LYS

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Mol	Chain	Res	Type
1	A	53	ARG
1	A	54	GLU
1	A	57	ILE
1	A	58	ILE
1	A	61	ASP
1	A	74	LEU
1	A	89	ARG
1	A	95	ARG
1	A	98	LEU
1	A	99	LYS
1	A	101	GLN
1	A	105	VAL
1	A	106	ASP
1	A	108	ASN
1	A	110	ARG
1	A	111	VAL
1	A	112	ASP
1	A	120	LEU
1	A	121	VAL
1	A	123	ASP
1	A	127	GLU
1	A	128	LEU
1	A	130	ARG
1	A	134	GLU
1	A	139	LYS
1	A	144	ILE
1	A	149	LEU
1	A	152	ARG
1	A	154	LEU
1	A	155	CYS
1	A	156	LEU
1	A	159	LEU
1	A	168	HIS
1	A	171	VAL
1	A	172	LEU
1	A	174	GLN
1	A	175	TYR
1	A	179	ILE
1	A	192	THR
1	A	194	PRO
1	A	201	LEU
1	A	203	CYS

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Mol	Chain	Res	Type
1	A	204	SER
1	A	205	MET
1	A	214	VAL
1	A	216	SER
1	A	217	ILE
1	A	218	LEU
1	A	219	ASP
1	A	223	THR
1	A	225	LYS
1	A	236	VAL
1	A	237	LYS
1	A	239	THR
1	A	242	LYS
1	A	243	CYS
1	A	244	ILE
1	A	245	LEU
1	A	247	SER
1	A	252	LYS
1	A	262	GLU
1	A	268	VAL
1	A	275	LEU
1	A	284	ARG
1	A	287	LYS
1	A	289	LEU
1	A	298	GLN
1	A	300	GLN
1	A	301	GLN
1	A	306	ARG
1	A	316	ASN
1	A	318	ILE
1	A	319	ILE
1	A	326	TYR
1	A	327	ASP
1	A	328	ASP
1	A	330	ILE
1	A	335	ILE
1	A	352	VAL
1	A	357	GLN
1	A	358	GLN
1	A	359	MET
1	A	367	ILE
1	A	369	HIS

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Mol	Chain	Res	Type
1	A	370	LEU
1	A	371	ILE
1	A	372	THR
1	A	378	GLN
1	A	380	LEU
1	A	391	VAL
1	A	392	SER
1	A	396	LEU
1	A	398	HIS
1	A	421	SER
1	A	422	VAL
1	A	426	CYS
1	A	431	LEU
1	A	433	CYS
1	A	435	LYS
1	A	441	LEU
1	A	449	LEU
1	A	452	LEU
1	A	453	VAL
1	A	466	ILE
1	A	467	LEU
1	A	474	LYS
1	A	479	VAL
1	A	480	ASN
1	A	484	LEU
1	A	485	SER
1	A	505	LEU
1	A	508	LYS
1	A	510	ILE
1	A	521	THR
1	A	525	LEU
1	A	534	LEU
1	A	535	ARG
1	A	536	SER
1	A	537	ILE
1	A	538	THR
1	A	548	LEU
1	A	549	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
1	A	131	GLN
1	A	147	ASN
1	A	174	GLN
1	A	211	ASN
1	A	228	HIS
1	A	278	HIS
1	A	292	ASN
1	A	299	ASN
1	A	316	ASN
1	A	343	HIS
1	A	357	GLN
1	A	369	HIS
1	A	378	GLN
1	A	413	HIS
1	A	439	ASN
1	A	493	GLN
1	A	509	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](#)

2 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	SF4	A	601	1	0,12,12	-	-	-		
3	FSO	A	602	1	1,12,12	7.01	1 (100%)	-		

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	SF4	A	601	1	-	-	0/6/5/5
3	FSO	A	602	1	-	-	0/2/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	FSO	S6-FE8	-7.01	2.11	2.29

There are no bond angle outliers.

There are no chirality outliers.

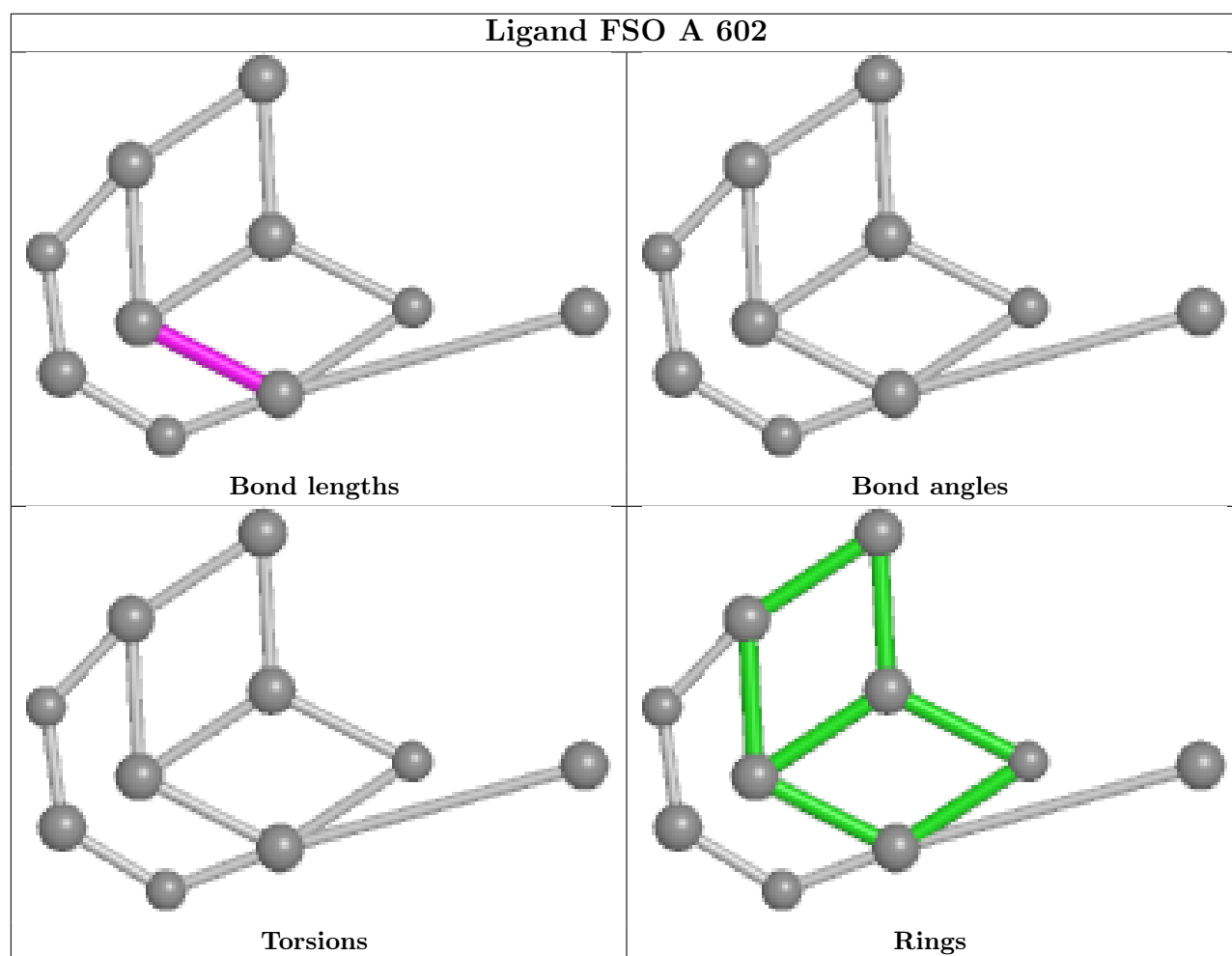
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	FSO	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 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 [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	550/558 (98%)	-0.14	9 (1%)	70 44	77, 122, 185, 306	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	LEU	3.3
1	A	485	SER	2.9
1	A	26	GLY	2.6
1	A	14	ALA	2.4
1	A	129	GLN	2.3
1	A	510	ILE	2.3
1	A	112	ASP	2.2
1	A	125	LEU	2.1
1	A	302	VAL	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](#)

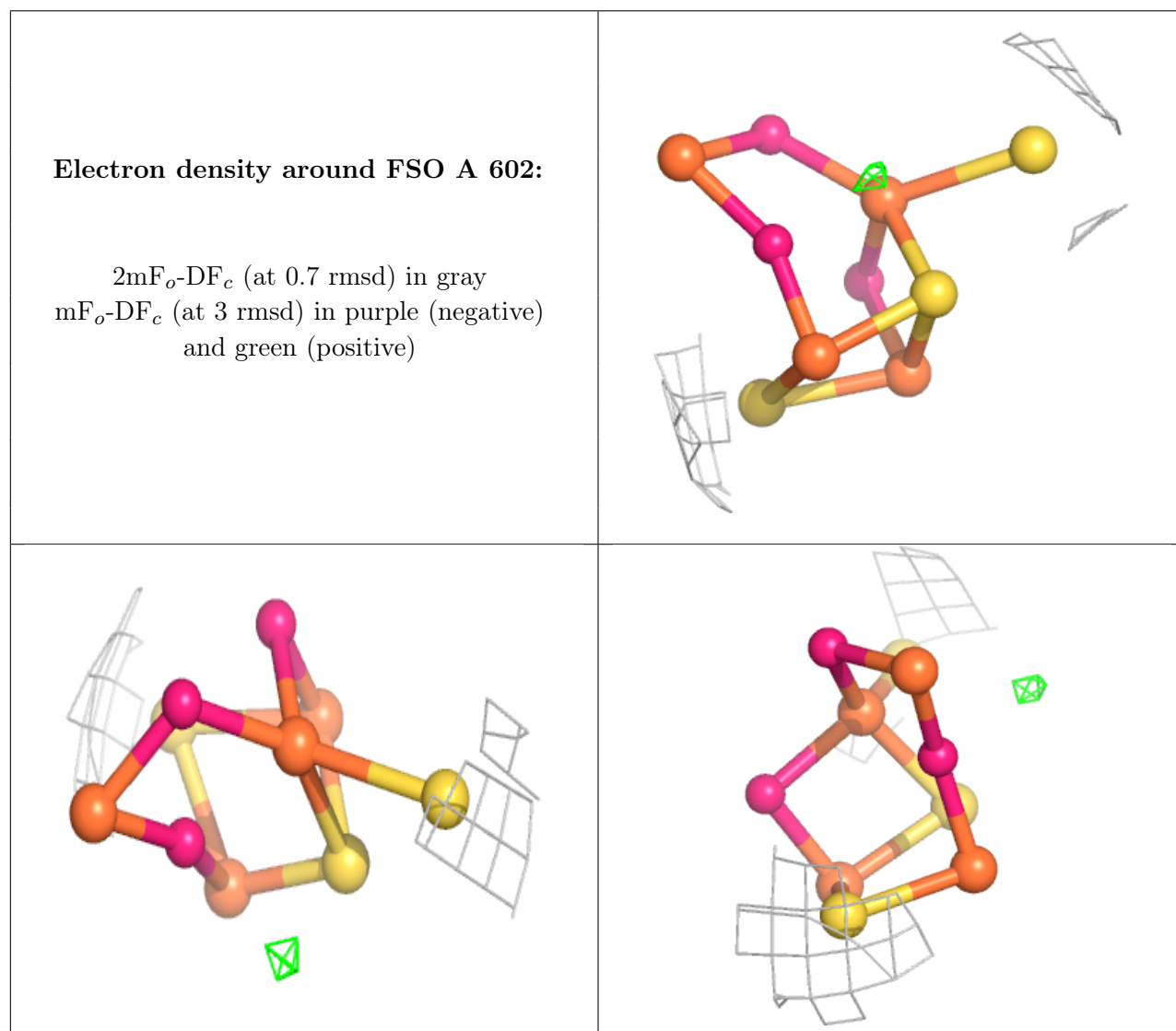
There are no oligosaccharides in this entry.

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
2	SF4	A	601	8/8	0.99	0.03	100,114,136,148	0
3	FSO	A	602	10/10	0.99	0.04	123,136,164,182	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.