

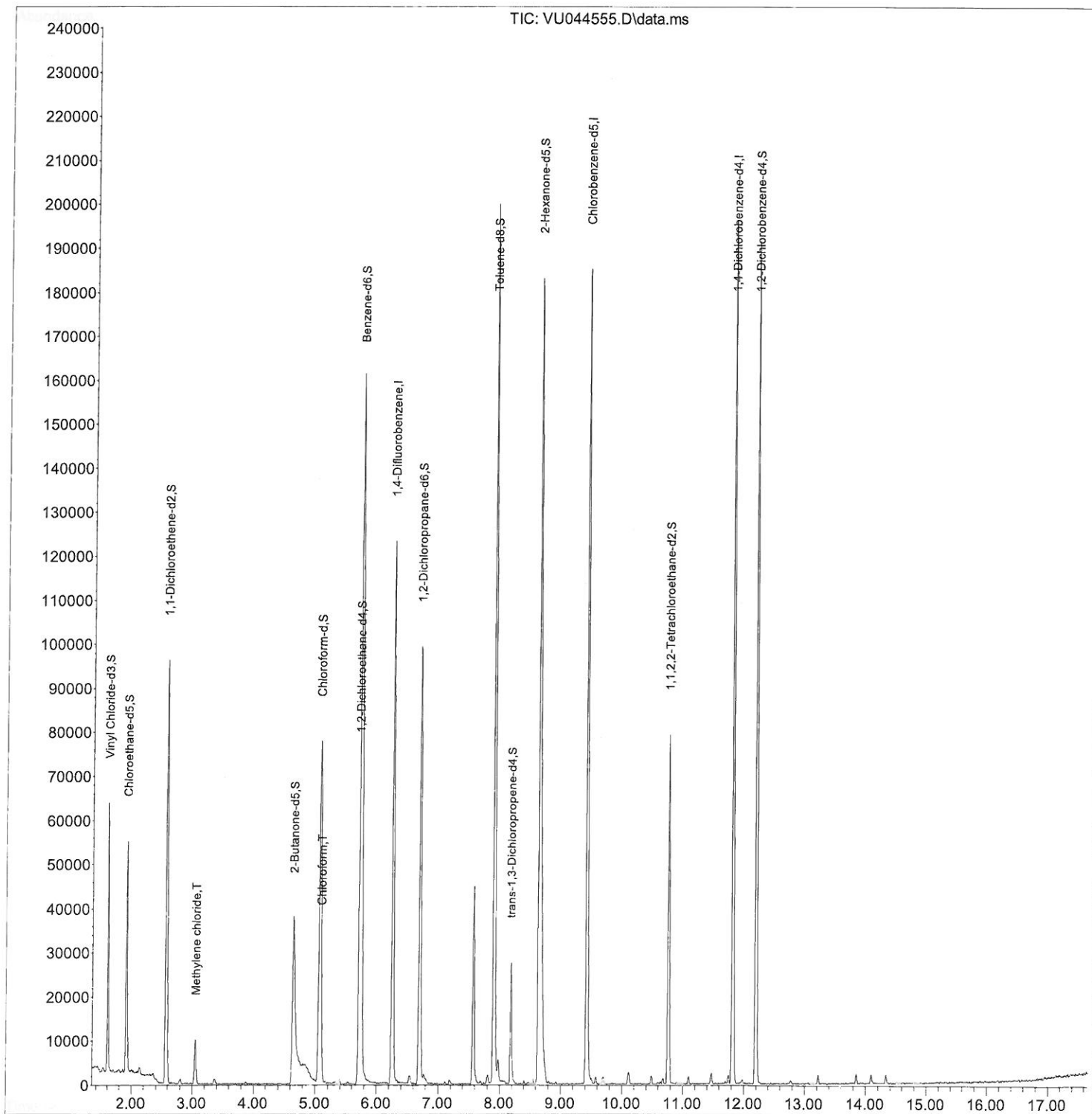
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090221\
Data File : VU044555.D
Acq On : 02 Sep 2021 16:04
Operator : SY/MD
Sample : VU0902WBL01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
VBLK013

Manual Integrations
APPROVED

MMDadoda
9/7/2021 9:07:36 AM

Quant Time: Sep 03 02:44:43 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR081821WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Fri Sep 03 02:43:14 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU090221\
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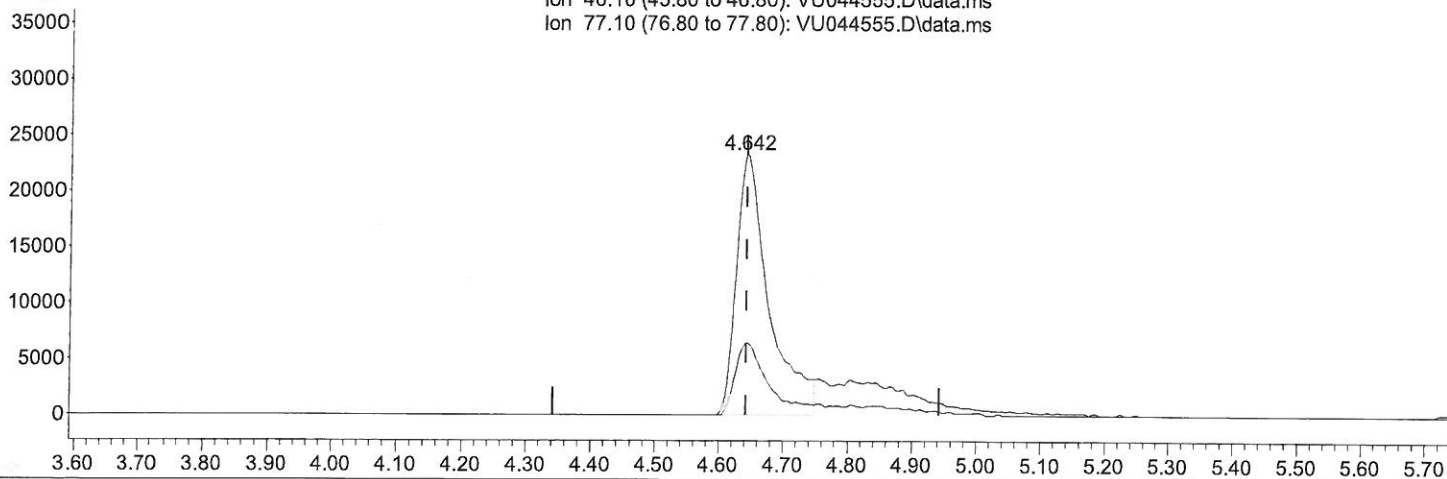
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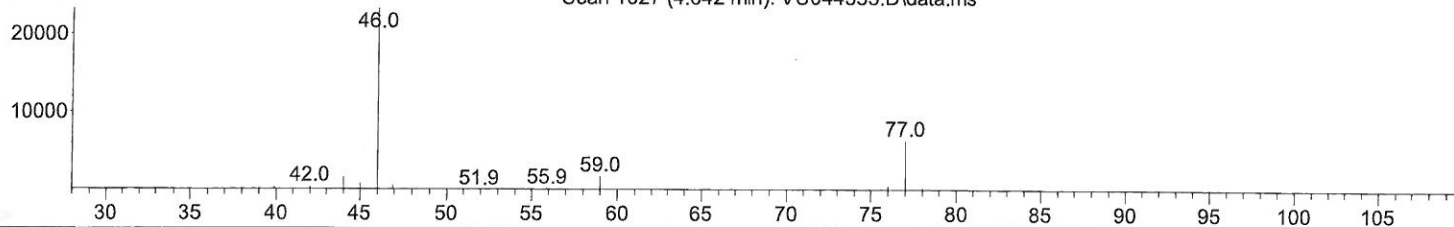
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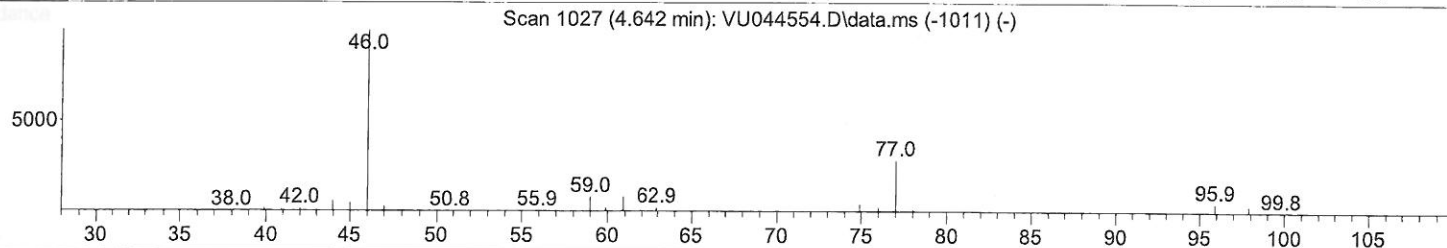
Ion 46.10 (45.80 to 46.80): VU044555.D\data.ms
 Ion 77.10 (76.80 to 77.80): VU044555.D\data.ms



Scan 1027 (4.642 min): VU044555.D\data.ms



Scan 1027 (4.642 min): VU044554.D\data.ms (-1011) (-)



TIC: VU044555.D\data.ms

(20) 2-Butanone-d5 (S)

4.642min (-0.000) 35.30 ug/L

response 81761

Ion	Exp%	Act%
46.10	100.00	100.00
77.10	25.00	25.63
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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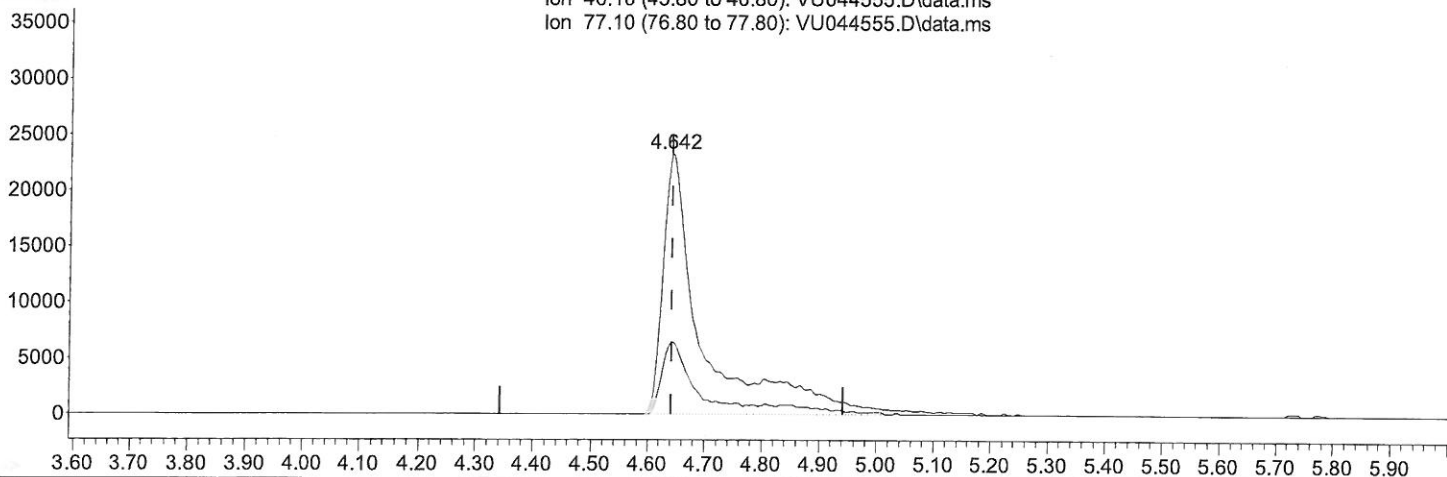
Instrument :
 MSVOA_U
 ClientSampleId :
 VBLK013

Manual Integrations
 APPROVED

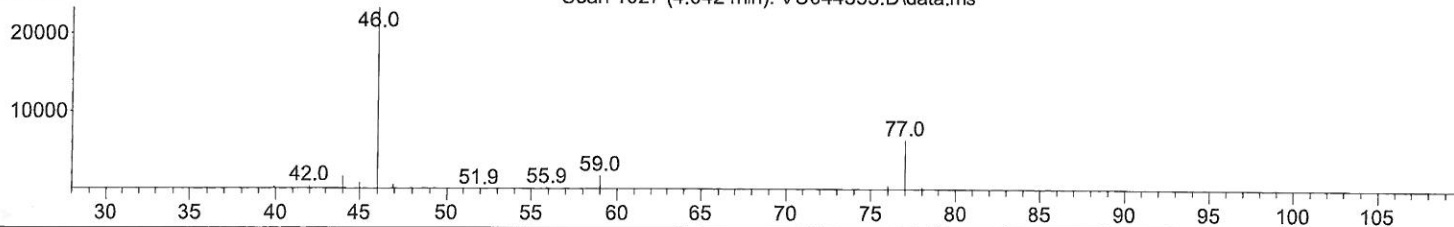
MMDadoda
 9/7/2021 9:07:36 AM

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 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR081821WMA.M
 Quant Title : TRACE VOA SFAM1.0
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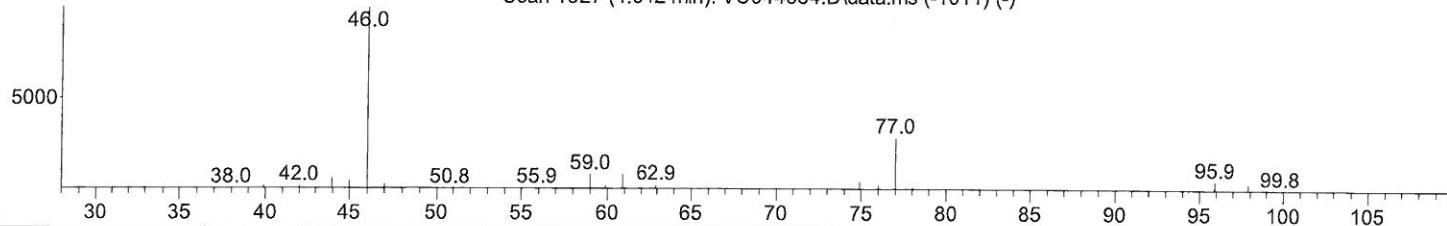
Ion 46.10 (45.80 to 46.80): VU044555.D\data.ms
 Ion 77.10 (76.80 to 77.80): VU044555.D\data.ms



Scan 1027 (4.642 min): VU044555.D\data.ms



Scan 1027 (4.642 min): VU044554.D\data.ms (-1011) (-)



TIC: VU044555.D\data.ms

(20) 2-Butanone-d5 (S)

4.642min (-0.000) 48.55 ug/L m

response 112445

Ion	Exp%	Act%
46.10	100.00	100.00
77.10	25.00	18.64
0.00	0.00	0.00
0.00	0.00	0.00

7 mg
 9/10/21

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Manual Integrations
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Difluorobenzene	6.256	114	103554	5.000	ug/L	0.00
28) Chlorobenzene-d5	9.423	117	107289	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.819	152	52739	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.601	65	42115	5.029	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery = 100.600%			
7) Chloroethane-d5	1.919	69	39978	5.492	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery = 109.800%			
11) 1,1-Dichloroethene-d2	2.572	65	17877	5.018	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery = 100.400%			
20) 2-Butanone-d5	4.642	46	112445m	48.546	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery = 97.100%			
24) Chloroform-d	5.070	84	73361	4.954	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 99.000%			
26) 1,2-Dichloroethane-d4	5.707	65	41274	5.111	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 102.200%			
32) Benzene-d6	5.736	84	154436	5.112	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery = 102.200%			
36) 1,2-Dichloropropane-d6	6.697	67	46791	4.841	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery = 96.800%			
41) Toluene-d8	7.903	98	136535	5.097	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery = 102.000%			
43) trans-1,3-Dichloroprop...	8.186	79	15917	4.586	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery = 91.800%			
46) 2-Hexanone-d5	8.645	63	96464	51.392	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery = 102.780%			
56) 1,1,2,2-Tetrachloroeth...	10.761	84	41922	5.150	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery = 103.000%			
66) 1,2-Dichlorobenzene-d4	12.198	152	51199	5.746	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery = 115.000%			
Target Compounds						
16) Methylene chloride	3.054	84	5116	0.433	ug/L	95
25) Chloroform	5.092	83	4312	0.261	ug/L	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

MD
 9/10/21