

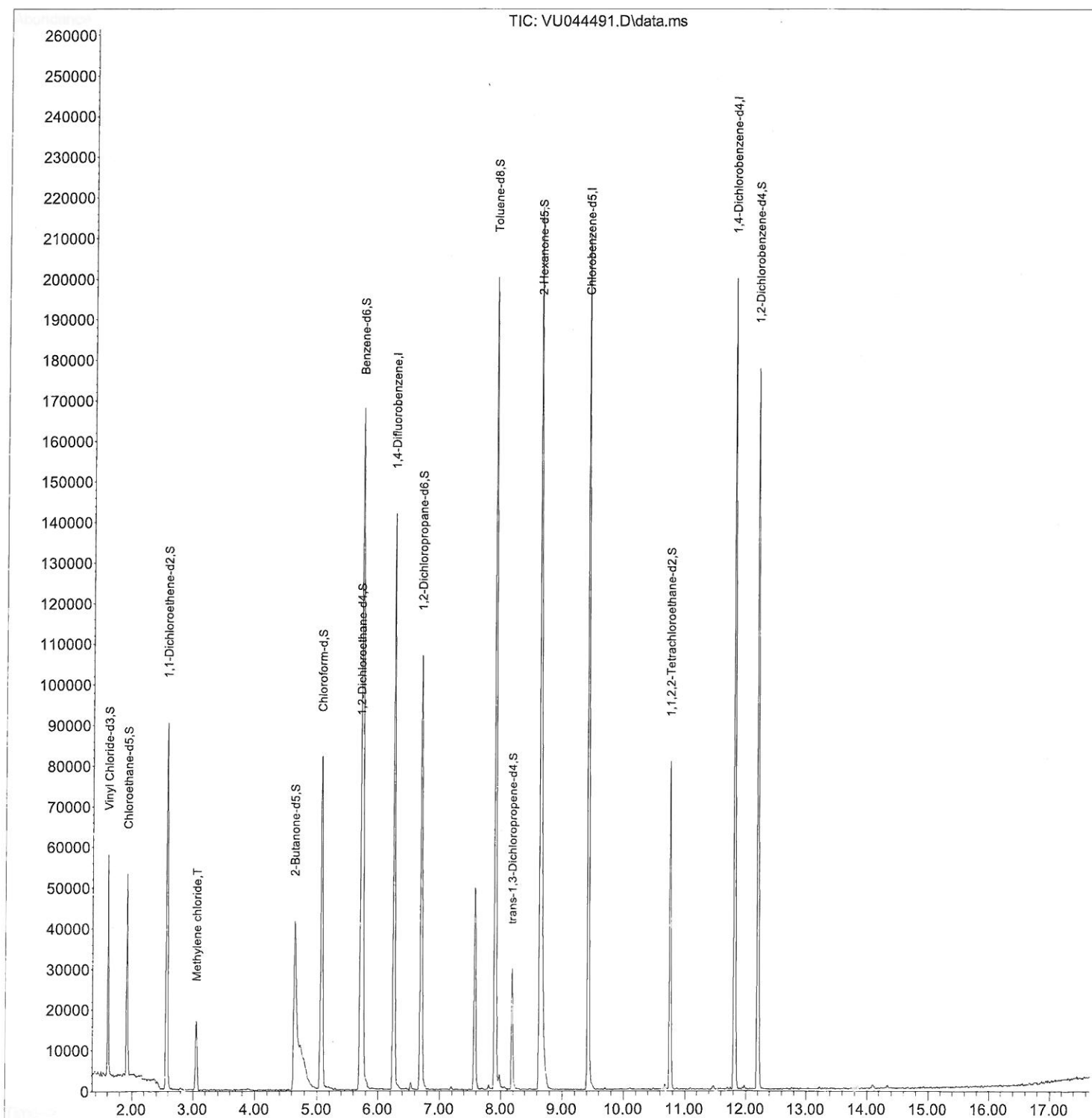
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU082321\
Data File : VU044491.D
Acq On : 23 Aug 2021 10:22
Operator : SY/MD
Sample : VU0823WBL01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
VBLK010

Manual Integrations
APPROVED

SAM
8/24/2021 11:20:50 AM

Quant Time: Aug 24 03:57:27 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR081821WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Tue Aug 24 03:56:17 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU082321\
 Data File : VU044491.D
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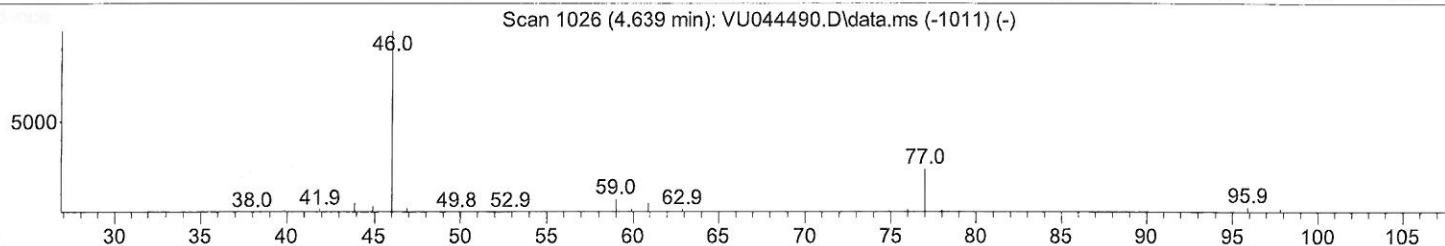
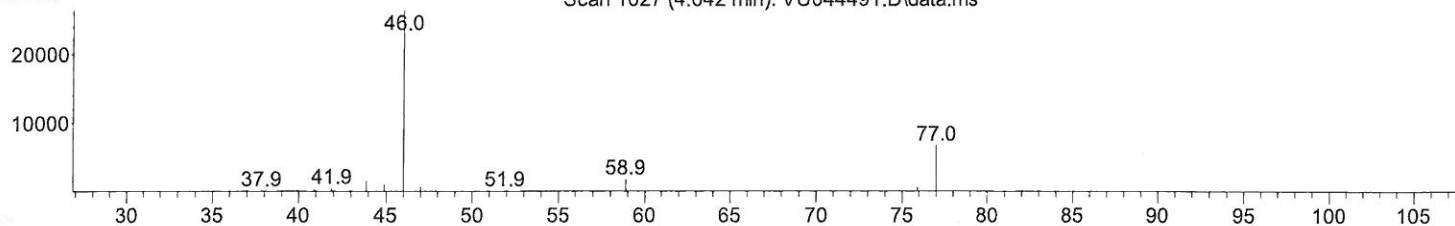
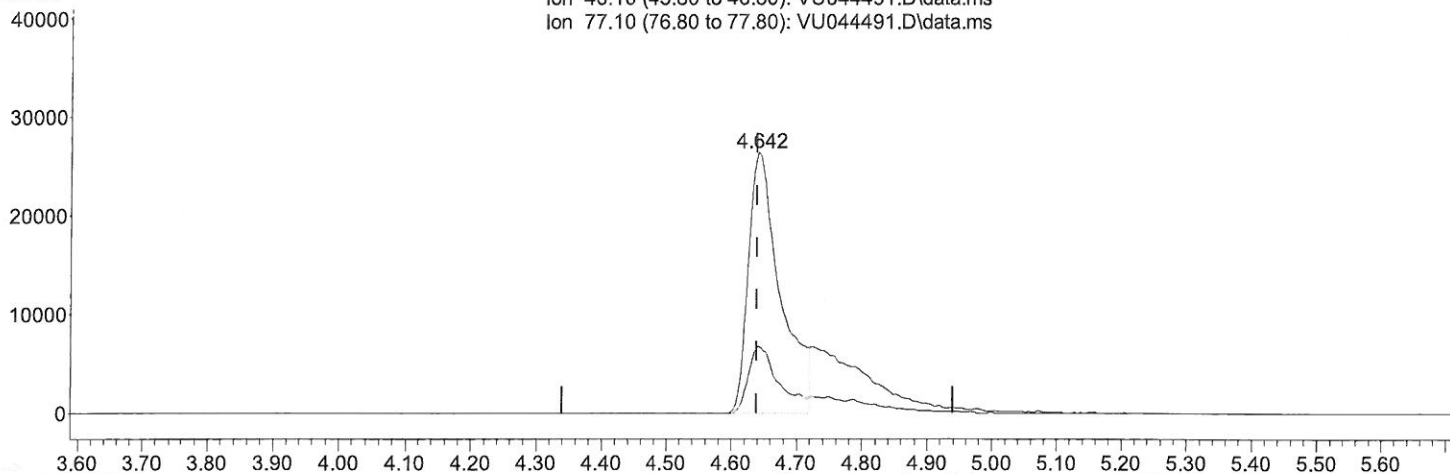
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Ion 46.10 (45.80 to 46.80): VU044491.D\data.ms
 Ion 77.10 (76.80 to 77.80): VU044491.D\data.ms



(20) 2-Butanone-d5 (S)

4.642min (+ 0.003) 36.27 ug/L

response 90457

Ion	Exp%	Act%
46.10	100.00	100.00
77.10	25.00	22.30
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

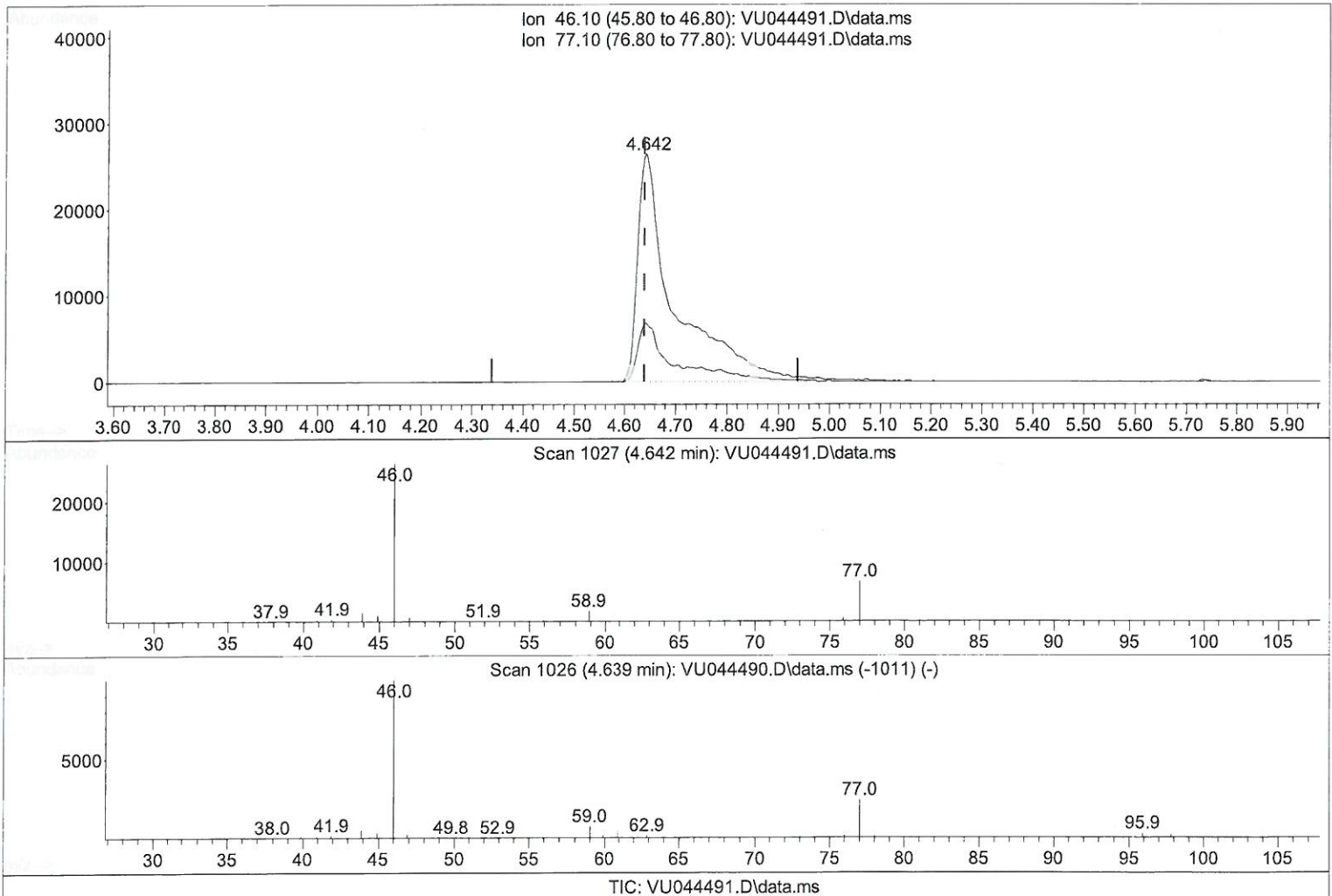
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Operator : SY/MD
Sample : VU0823WBL01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_U
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Manual Integrations
APPROVED

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Quant Title : TRACE VOA SFAM1.0
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Response via : Initial Calibration



(20) 2-Butanone-d5 (S)

4.642min (+ 0.003) 52.87 ug/L m

response 131852

Ion	Exp%	Act%
46.10	100.00	100.00
77.10	25.00	15.30#
0.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU082321\
 Data File : VU044491.D
 Acq On : 23 Aug 2021 10:22
 Operator : SY/MD
 Sample : VU0823WBL01
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VBLK010

Manual Integrations
 APPROVED

SAM
 8/24/2021 11:20:50 AM

Quant Time: Aug 24 03:57:27 2021
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 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Difluorobenzene	6.256	114	111492	5.000	ug/L	0.00
28) Chlorobenzene-d5	9.423	117	113583	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.819	152	50076	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.601	65	37648	4.175	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery =	83.600%		
7) Chloroethane-d5	1.919	69	36770	4.692	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery =	93.800%		
11) 1,1-Dichloroethene-d2	2.572	65	17806	4.642	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery =	92.800%		
20) 2-Butanone-d5	4.642	46	131852m	52.872	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery =	105.740%		
24) Chloroform-d	5.073	84	76872	4.821	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery =	96.400%		
26) 1,2-Dichloroethane-d4	5.710	65	43993	5.060	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery =	101.200%		
32) Benzene-d6	5.735	84	153898	4.812	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery =	96.200%		
36) 1,2-Dichloropropane-d6	6.697	67	51718	5.054	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery =	101.000%		
41) Toluene-d8	7.906	98	130730	4.609	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery =	92.200%		
43) trans-1,3-Dichloroprop...	8.185	79	17304	4.709	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery =	94.200%		
46) 2-Hexanone-d5	8.645	63	105478	53.080	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery =	106.160%		
56) 1,1,2,2-Tetrachloroeth...	10.761	84	42150	4.891	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery =	97.800%		
66) 1,2-Dichlorobenzene-d4	12.198	152	43781	5.175	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery =	103.400%		
Target Compounds						Qvalue
16) Methylene chloride	3.057	84	8092	0.636	ug/L	98

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

ME
 9/7/21