

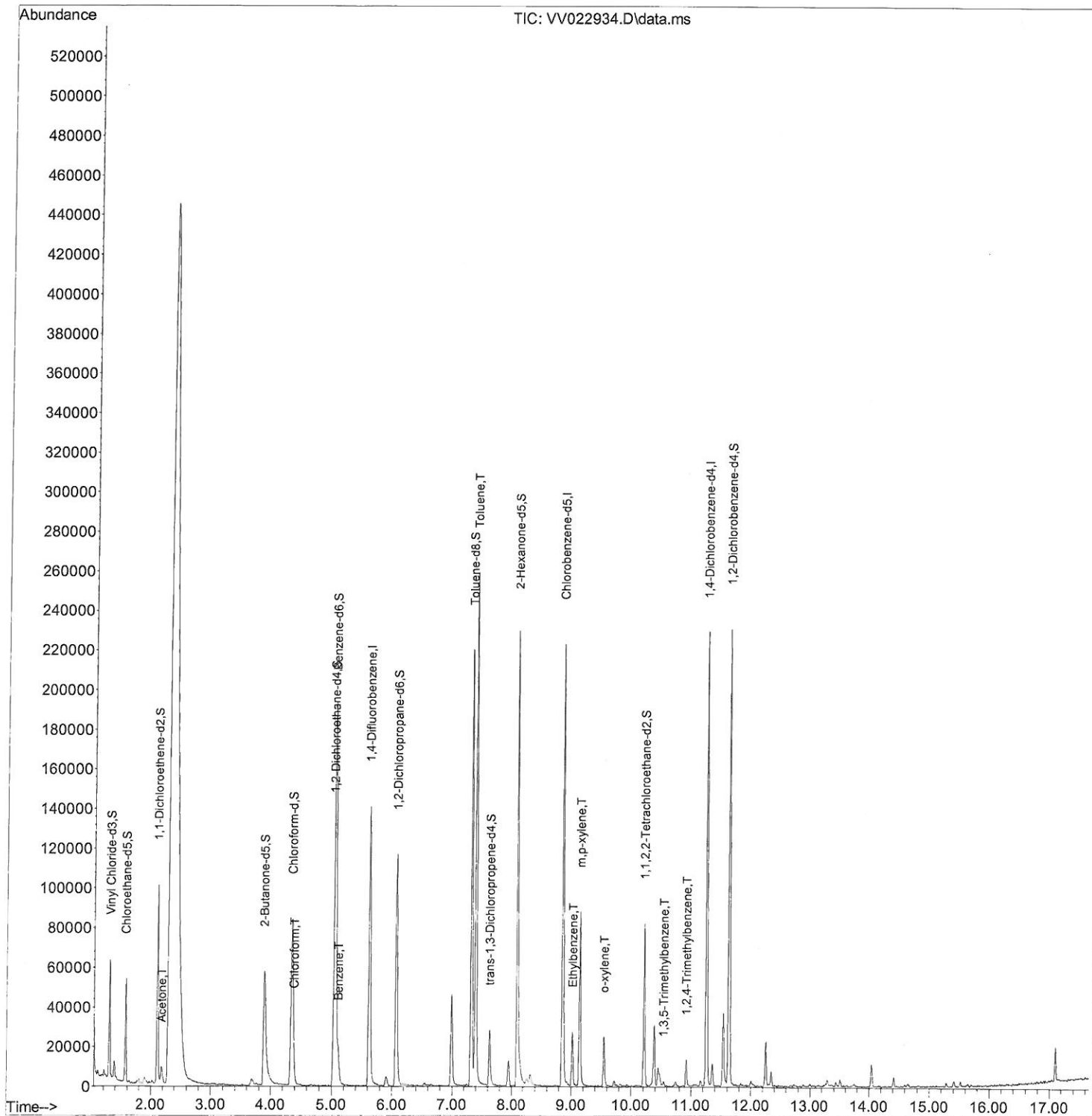
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102021\
Data File : VV022934.D
Acq On : 20 Oct 2021 16:06
Operator : SY/MD
Sample : M4200-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BFGF2

Manual Integrations
APPROVED

MMDadoda
10/22/2021 11:29:11 AM

Quant Time: Oct 21 02:46:36 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100721WMA.M
Quant Title : TRACE VOA SFAM1.0
QLast Update : Thu Oct 21 02:43:59 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

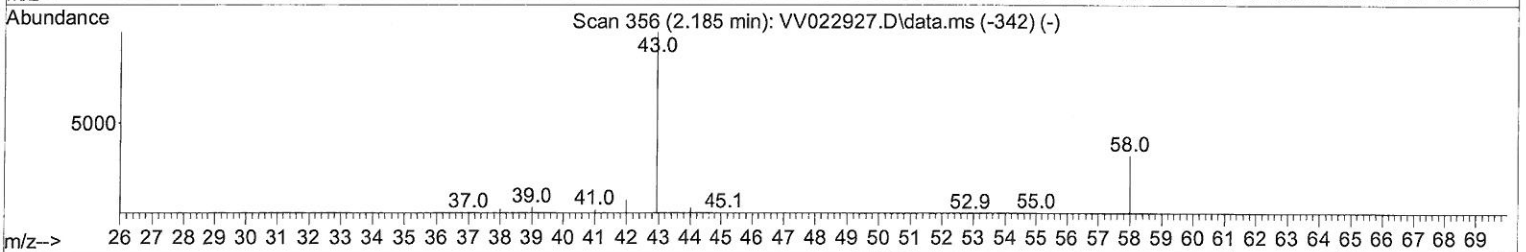
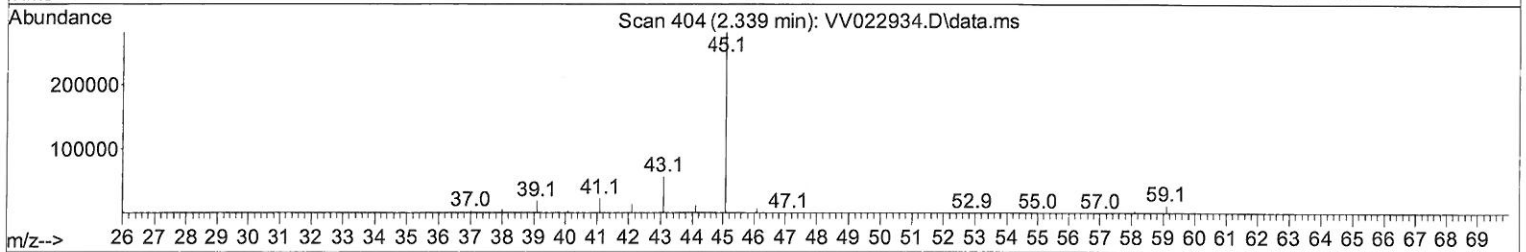
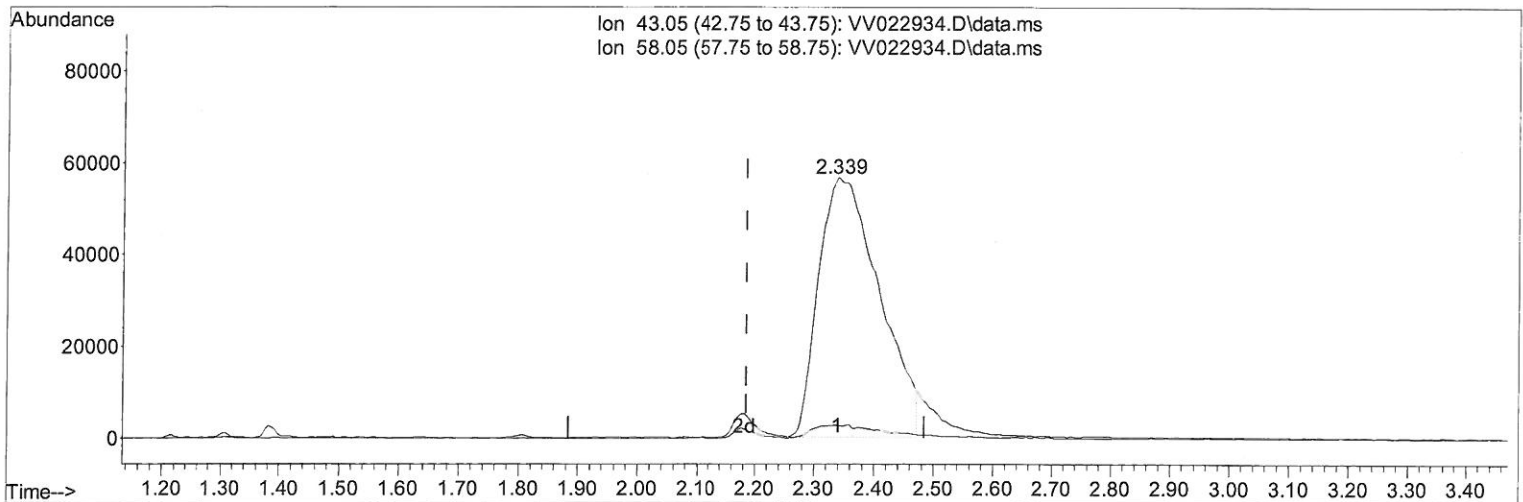
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TIC: VV022934.D\data.ms

(13) Acetone (T)

2.339min (+ 0.155) 315.16 ug/L

response 398980

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	2.34
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

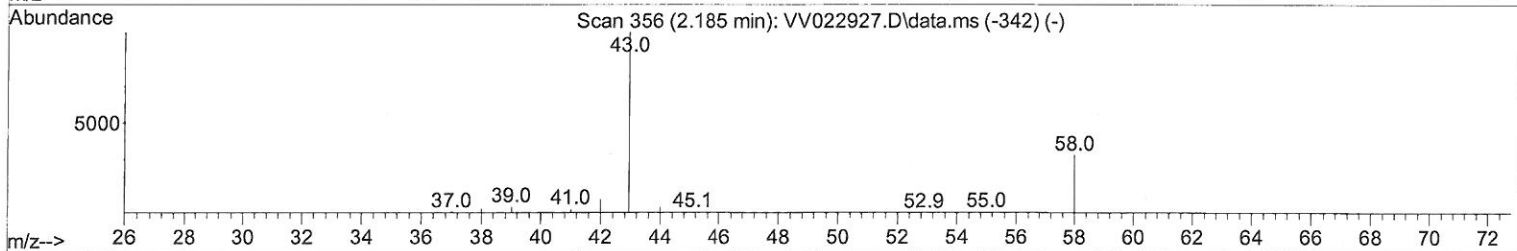
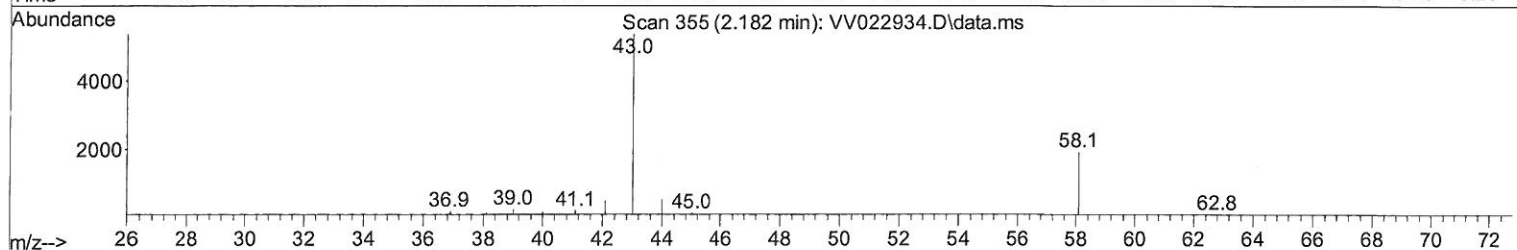
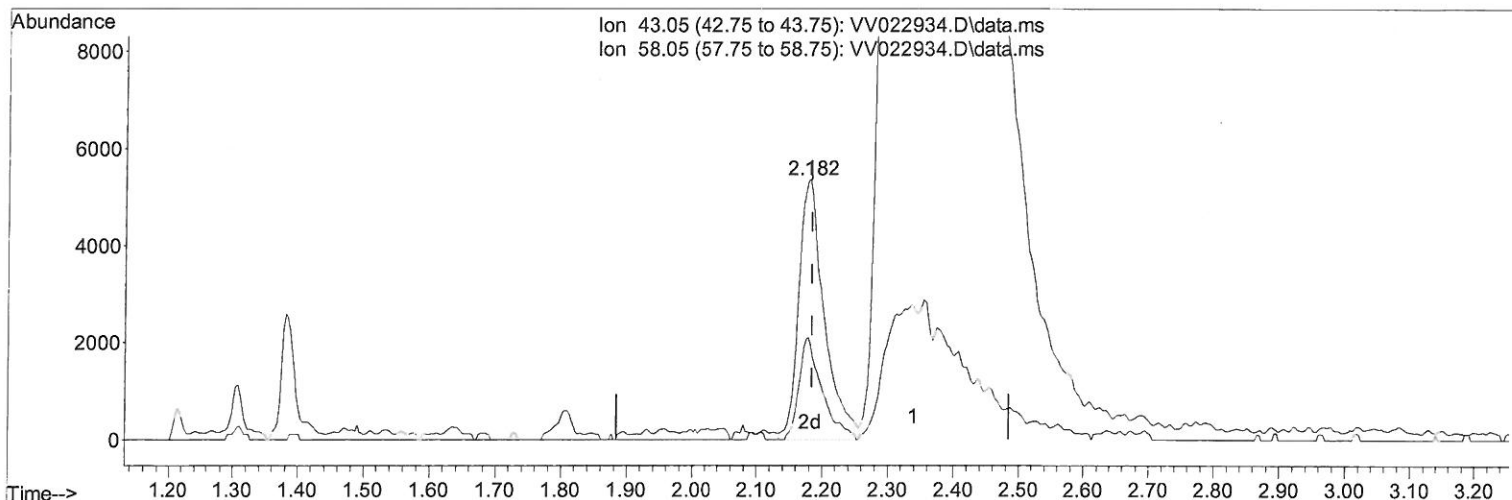
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Manual Integrations
 APPROVED

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 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Oct 21 02:43:59 2021
 Response via : Initial Calibration



TIC: VV022934.D\data.ms

(13) Acetone (T)

2.182min (-0.003) 11.43 ug/L m

response 14465

Ion	Exp%	Act%
43.05	100.00	100.00
58.05	12.10	64.42#
0.00	0.00	0.00
0.00	0.00	0.00

MD
10/27/21

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102021\
 Data File : VV022934.D
 Acq On : 20 Oct 2021 16:06
 Operator : SY/MD
 Sample : M4200-07
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BFGF2

Manual Integrations
 APPROVED

MMDadoda
 10/22/2021 11:29:11 AM

Quant Time: Oct 21 02:46:36 2021
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 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Oct 21 02:43:59 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	125846	5.000	ug/L	0.00
28) Chlorobenzene-d5	8.854	117	128849	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.252	152	61298	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.304	65	33488	3.868	ug/L	0.00
Spiked Amount	5.000	Range 40 - 130	Recovery	=	77.400%	
7) Chloroethane-d5	1.568	69	30314	3.896	ug/L	0.00
Spiked Amount	5.000	Range 65 - 130	Recovery	=	78.000%	
11) 1,1-Dichloroethene-d2	2.108	63	50329	2.963	ug/L	0.00
Spiked Amount	5.000	Range 60 - 125	Recovery	=	59.200%#	
20) 2-Butanone-d5	3.886	46	100984	44.481	ug/L	0.00
Spiked Amount	50.000	Range 40 - 130	Recovery	=	88.960%	
24) Chloroform-d	4.352	84	88722	4.972	ug/L	0.00
Spiked Amount	5.000	Range 70 - 125	Recovery	=	99.400%	
26) 1,2-Dichloroethane-d4	5.034	65	42871	4.950	ug/L	0.00
Spiked Amount	5.000	Range 70 - 130	Recovery	=	99.000%	
32) Benzene-d6	5.053	84	177288	4.622	ug/L	0.00
Spiked Amount	5.000	Range 70 - 125	Recovery	=	92.400%	
36) 1,2-Dichloropropane-d6	6.072	67	58071	5.112	ug/L	0.00
Spiked Amount	5.000	Range 60 - 140	Recovery	=	102.200%	
41) Toluene-d8	7.317	98	146952	4.432	ug/L	0.00
Spiked Amount	5.000	Range 70 - 130	Recovery	=	88.600%	
43) trans-1,3-Dichloroprop...	7.625	79	17417	4.622	ug/L	0.00
Spiked Amount	5.000	Range 55 - 130	Recovery	=	92.400%	
46) 2-Hexanone-d5	8.088	63	77787	51.223	ug/L	0.00
Spiked Amount	50.000	Range 45 - 130	Recovery	=	102.440%	
56) 1,1,2,2-Tetrachloroeth...	10.217	84	38802	5.042	ug/L	0.00
Spiked Amount	5.000	Range 65 - 120	Recovery	=	100.800%	
66) 1,2-Dichlorobenzene-d4	11.625	152	62353	5.486	ug/L	0.00
Spiked Amount	5.000	Range 80 - 120	Recovery	=	109.800%	
Target Compounds						
13) Acetone	2.182	43	14465m	11.426	ug/L	Qvalue > 7070
25) Chloroform	4.381	83	6500	0.377	ug/L	86 10/27/21
33) Benzene	5.105	78	18020	0.519	ug/L	100
42) Toluene	7.391	91	190938	5.420	ug/L	98
52) Ethylbenzene	9.018	91	18731	0.524	ug/L	100
53) m,p-xylene	9.140	106	25320	1.804	ug/L	97
54) o-xylene	9.548	106	6935	0.525	ug/L	91
62) 1,3,5-Trimethylbenzene	10.542	105	1556	0.062	ug/L	93
63) 1,2,4-Trimethylbenzene	10.918	105	7208	0.286	ug/L	97

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