

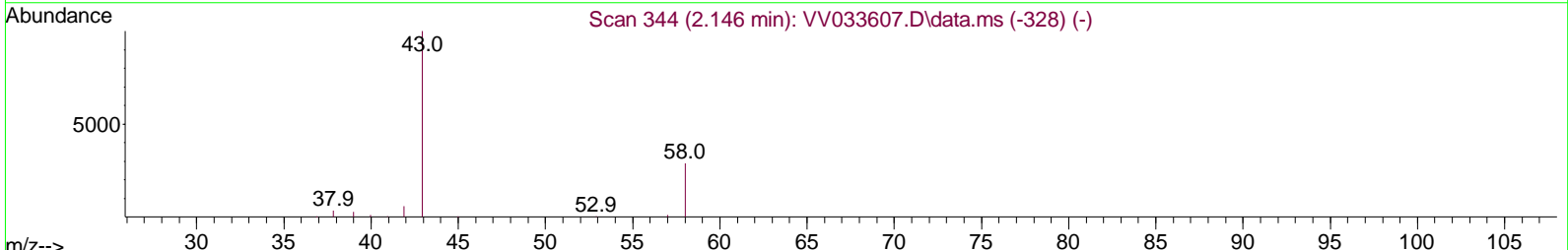
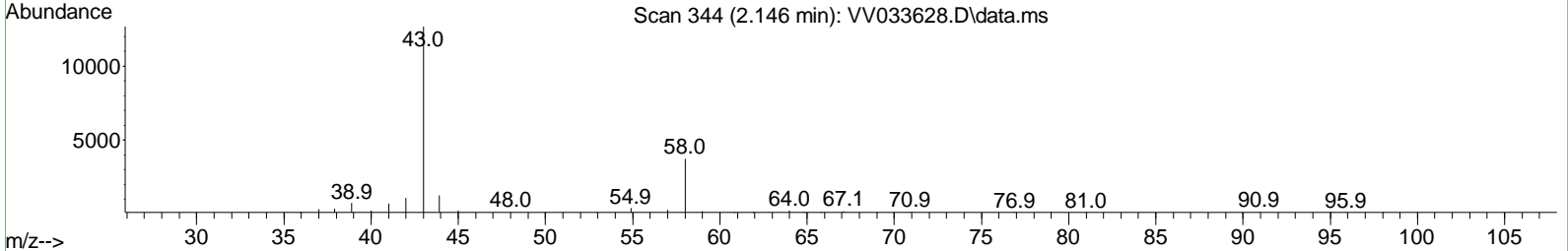
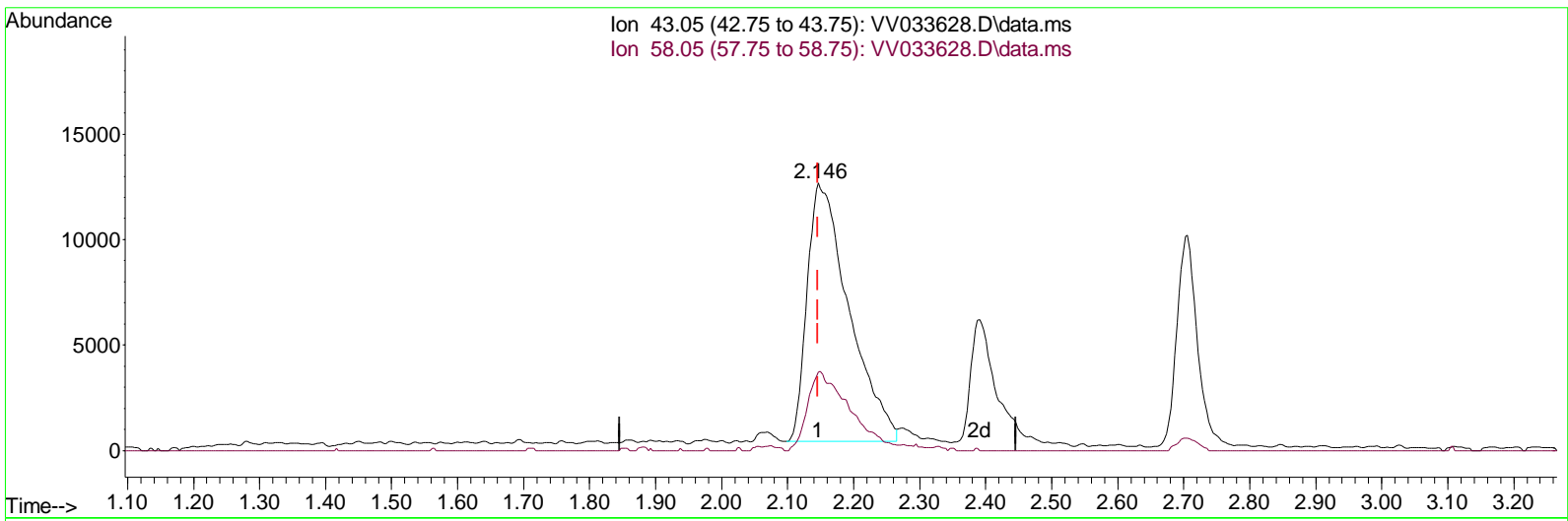
Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV011524\  
 Data File : VV033628.D  
 Acq On : 15 Jan 2024 09:05  
 Operator : SY/MD  
 Sample : VSTDCCC005  
 Mi sc : 25.0mL/MSVOA\_V/WATER  
 ALS Vial : 2 Sample Multiplier: 1

**Instrument :**  
 MSVOA\_V  
**LabSampleID :**  
 VSTDCCC005

**Manual Integrations APPROVED**

Reviewed By : Semsettin Yesilyurt 01/18/2024  
 Supervised By : Mahesh Dadoda 01/18/2024

Quant Time: Jan 16 03:36:01 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR011224WMA.M  
 Quant Title : TRACE VOA SFAM1.0  
 QLast Update : Sat Jan 13 00:29:44 2024  
 Response via : Initial Calibration



TIC: VV033628.D\data.ms

(13) Acetone (T)

2.146min (+ 0.000) 37.56 ug/L

response	53780	
Ion	Exp%	Act%
43.05	100.00	100.00
58.05	25.50	30.82
0.00	0.00	0.00
0.00	0.00	0.00

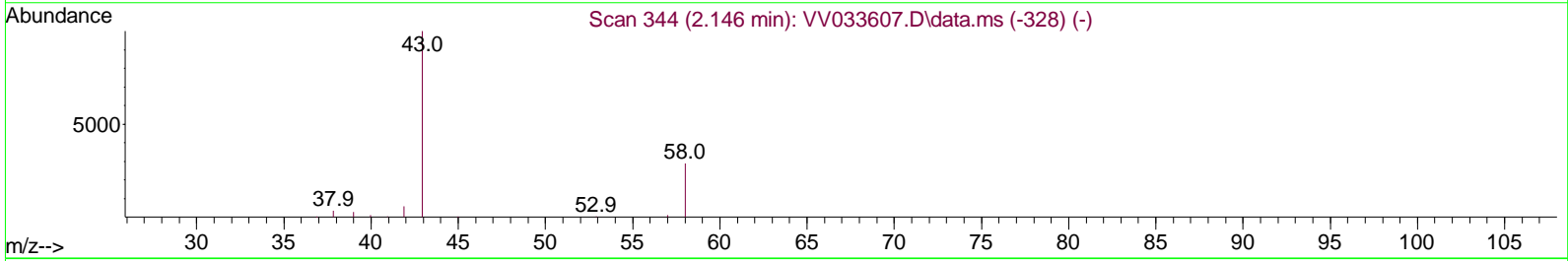
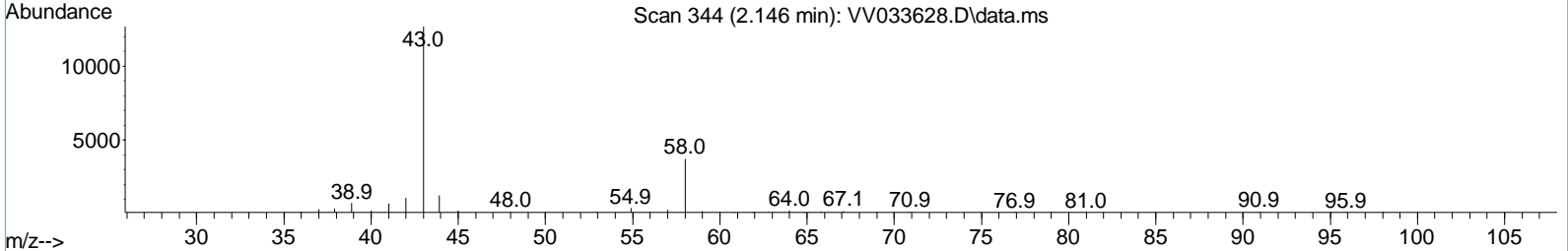
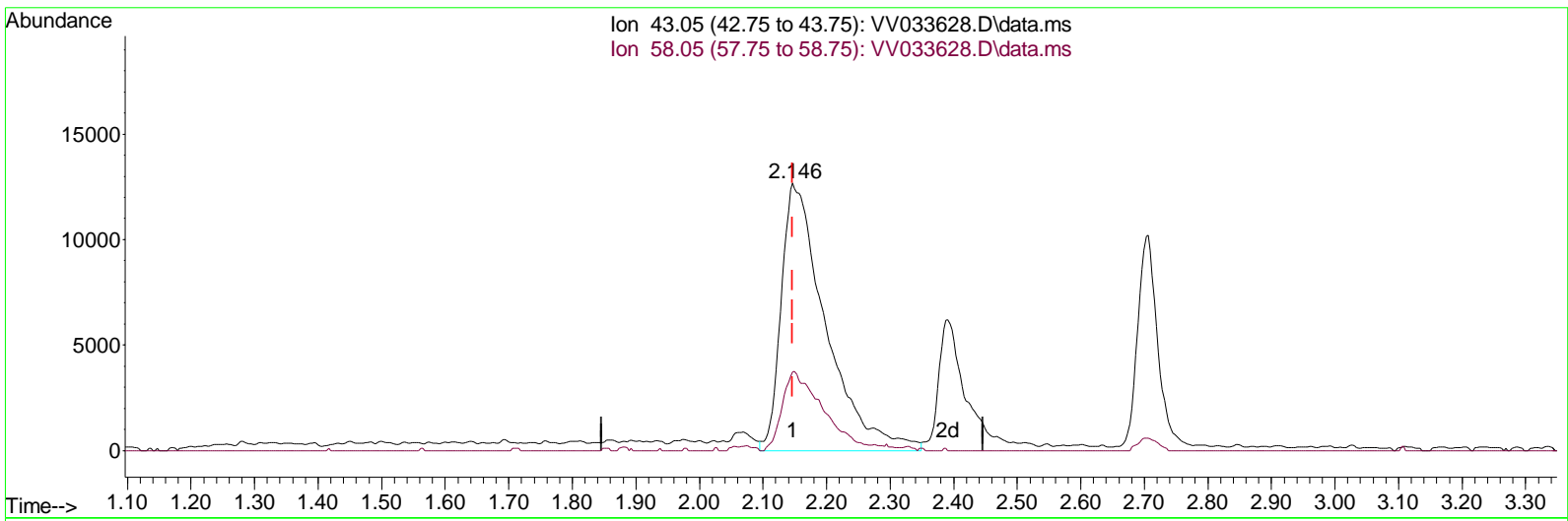
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TIC: VV033628.D\data.ms

(13) Acetone (T)

2.146min (+ 0.000) 42.94 ug/L m

response	61474	
Ion	Exp%	Act%
43.05	100.00	100.00
58.05	25.50	26.96
0.00	0.00	0.00
0.00	0.00	0.00

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
<b>Internal Standards</b>						
1) 1,4-Di fl uorobenzene	5.535	114	131908	5.000	ug/L	0.00
28) Chl orobenzene-d5	8.786	117	117814	5.000	ug/L	0.00
58) 1,4-Di chl orobenzene-d4	11.185	152	59099	5.000	ug/L	0.00
<b>System Moni toring Compounds</b>						
4) Vi nyl Chl ori de-d3	1.282	65	31616	4.383	ug/L	0.00
Spi ked Amount 5.000	Range 40 - 130		Recovery =	87.600%		
7) Chl oroethane-d5	1.536	69	28670	4.482	ug/L	0.00
Spi ked Amount 5.000	Range 65 - 130		Recovery =	89.600%		
11) 1,1-Di chl oroethene-d2	2.060	65	14411	4.433	ug/L	0.00
Spi ked Amount 5.000	Range 60 - 125		Recovery =	88.600%		
20) 2-Butanone-d5	3.818	46	75592	42.862	ug/L	0.00
Spi ked Amount 50.000	Range 40 - 130		Recovery =	85.720%		
24) Chl oroform-d	4.249	84	73508	4.713	ug/L	0.00
Spi ked Amount 5.000	Range 70 - 125		Recovery =	94.200%		
26) 1,2-Di chl oroethane-d4	4.944	65	31179	4.458	ug/L	0.00
Spi ked Amount 5.000	Range 70 - 130		Recovery =	89.200%		
32) Benzene-d6	4.960	84	133295	4.689	ug/L	0.00
Spi ked Amount 5.000	Range 70 - 125		Recovery =	93.800%		
36) 1,2-Di chl oropropane-d6	5.989	67	41796	4.410	ug/L	0.00
Spi ked Amount 5.000	Range 60 - 140		Recovery =	88.200%		
41) Tol uene-d8	7.243	98	120918	4.765	ug/L	0.00
Spi ked Amount 5.000	Range 70 - 130		Recovery =	95.200%		
43) trans-1,3-Di chl oroprop. . .	7.555	79	14144	4.317	ug/L	0.00
Spi ked Amount 5.000	Range 55 - 130		Recovery =	86.400%		
46) 2-Hexanone-d5	8.027	63	55787	43.366	ug/L	0.00
Spi ked Amount 50.000	Range 45 - 130		Recovery =	86.740%		
56) 1,1,2,2-Tetrachl oroeth. . .	10.153	84	24591	4.441	ug/L	0.00
Spi ked Amount 5.000	Range 65 - 120		Recovery =	88.800%		
66) 1,2-Di chl orobenzene-d4	11.561	152	40299	4.441	ug/L	0.00
Spi ked Amount 5.000	Range 80 - 120		Recovery =	88.800%		
<b>Target Compounds</b>						
2) Di chl orodi fl uoromethane	1.108	85	61683	5.132	ug/L	99
3) Chl oromethane	1.217	50	69844	4.847	ug/L	99
5) Vi nyl chl ori de	1.285	62	68627	5.001	ug/L	92
6) Bromomethane	1.490	94	27383	3.990	ug/L	99
8) Chl oroethane	1.552	64	44498	5.340	ug/L	95
9) Tri chl orofl uoromethane	1.716	101	89769	5.384	ug/L	100
10) 1,1,2-Tri chl oro-1,2,2-. . .	2.069	101	49600	5.294	ug/L	98
12) 1,1-Di chl oroethene	2.072	96	48040	5.167	ug/L	97
13) Acetone	2.146	43	61474m	42.938	ug/L	
14) Carbon di sul fi de	2.243	76	170868	4.967	ug/L	100
15) Methyl Acetate	2.391	43	15553	4.545	ug/L	99
16) Methyl ene chl ori de	2.449	84	49001	4.465	ug/L	98
17) Methyl tert-butyl Ether	2.703	73	99281	4.716	ug/L	99
18) trans-1,2-Di chl oroethene	2.696	96	52312	4.980	ug/L	97
19) 1,1-Di chl oroethane	3.111	63	103862	4.977	ug/L	99
21) 2-Butanone	3.899	43	85424	42.985	ug/L	98
22) ci s-1,2-Di chl oroethene	3.812	96	54403	4.947	ug/L	95
23) Bromochl oromethane	4.150	128	20306	5.063	ug/L	98

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**LabSampleId :**  
 VSTDCCC005

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 Quant Title : TRACE VOA SFAM1.0  
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Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
25) Chl oroform	4.275	83	96531	5.037	ug/L	98
27) 1,2-Di chl oroethane	5.043	62	53367	4.888	ug/L	100
29) 1,1,1-Tri chl oroethane	4.510	97	88257	5.283	ug/L	99
30) Cycl ohexane	4.580	56	102489	5.119	ug/L	98
31) Carbon tetrachl ori de	4.735	117	77401	5.435	ug/L	98
33) Benzene	5.008	78	216403	5.174	ug/L	100
34) Tri chl oroethene	5.831	95	56444	5.069	ug/L	98
35) Methyl cycl ohexane	6.047	83	99479	5.278	ug/L	98
37) 1,2-Di chl oropropane	6.092	63	50448	4.856	ug/L	100
38) Bromodi chl oromethane	6.432	83	64102	5.063	ug/L	99
39) ci s-1,3-Di chl oropropene	6.953	75	73087	4.927	ug/L	96
40) 4-Methyl -2-pentanone	7.159	43	219789	46.028	ug/L	98
42) Tol uene	7.313	91	229296	5.183	ug/L	99
44) trans-1,3-Di chl oropropene	7.580	75	57352	4.797	ug/L	98
45) 1,1,2-Tri chl oroethane	7.770	97	28297	4.903	ug/L	97
47) Tetrachl oroethene	7.902	164	43304	5.413	ug/L	97
48) 2-Hexanone	8.079	43	150558	43.039	ug/L	98
49) Di bromochl oromethane	8.175	129	35646	4.961	ug/L	98
50) 1,2-Di bromoethane	8.281	107	26424	4.961	ug/L	97
51) Chl orobenzene	8.812	112	136644	5.042	ug/L	99
52) Ethyl benzene	8.944	91	263207	5.133	ug/L	99
53) m, p-Xyl ene	9.072	106	96756	5.121	ug/L	97
54) o-Xyl ene	9.477	106	93916	5.108	ug/L	99
55) Styrene	9.493	104	151197	5.160	ug/L	99
57) 1,1,2,2-Tetrachl oroethane	10.178	83	31113	4.912	ug/L	97
59) Bromoform	9.667	173	17683	4.957	ug/L	99
60) I sopropyl benzene	9.866	105	246550	5.273	ug/L	99
61) 1,2,3-Tri chl oropropane	10.210	75	22309	4.709	ug/L	95
62) 1,3,5-Tri methyl benzene	10.474	105	215827	5.342	ug/L	100
63) 1,2,4-Tri methyl benzene	10.850	105	218340	5.353	ug/L	100
64) 1,3-Di chl orobenzene	11.117	146	107681	5.207	ug/L	100
65) 1,4-Di chl orobenzene	11.207	146	100897	4.983	ug/L	98
67) 1,2-Di chl orobenzene	11.577	146	91110	5.070	ug/L	96
68) 1,2-Di bromo-3-chl oropr...	12.365	75	4925	4.730	ug/L	93
69) 1,3,5-Tri chl orobenzene	12.580	180	83372	5.193	ug/L	100
70) 1,2,4-tri chl orobenzene	13.197	180	61740	4.616	ug/L	100
71) Naphthal ene	13.442	128	71595	3.207	ug/L	100
72) 1,2,3-Tri chl orobenzene	13.680	180	46293	4.316	ug/L	100

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

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 Misc : 25.0mL/MSVOA\_V/WATER  
 ALS Vial : 2 Sample Multiplier: 1

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 MSVOA\_V  
**Lab Sample ID :**  
 VSTDC005

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