

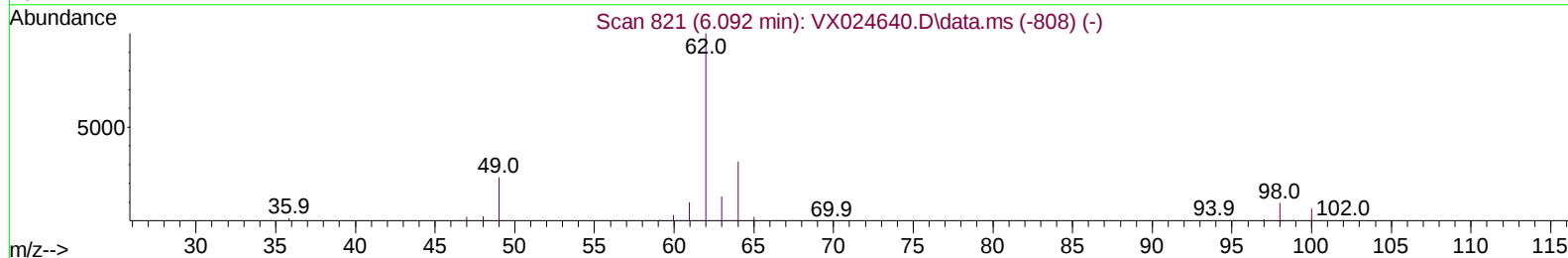
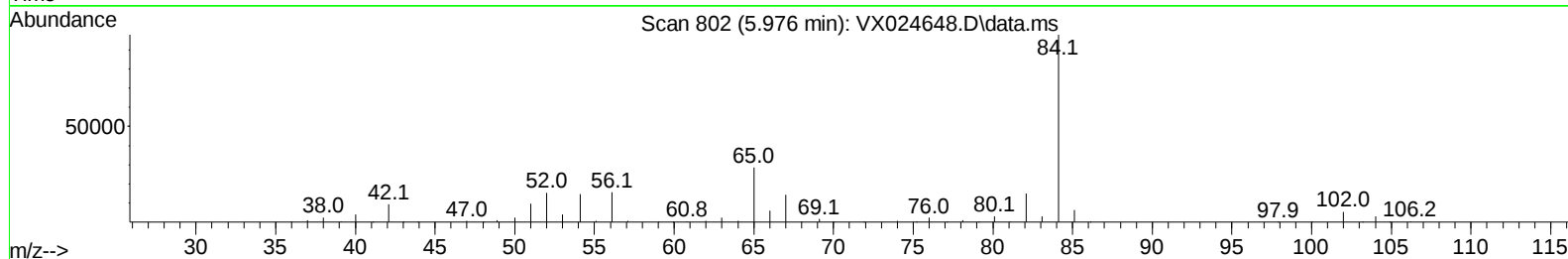
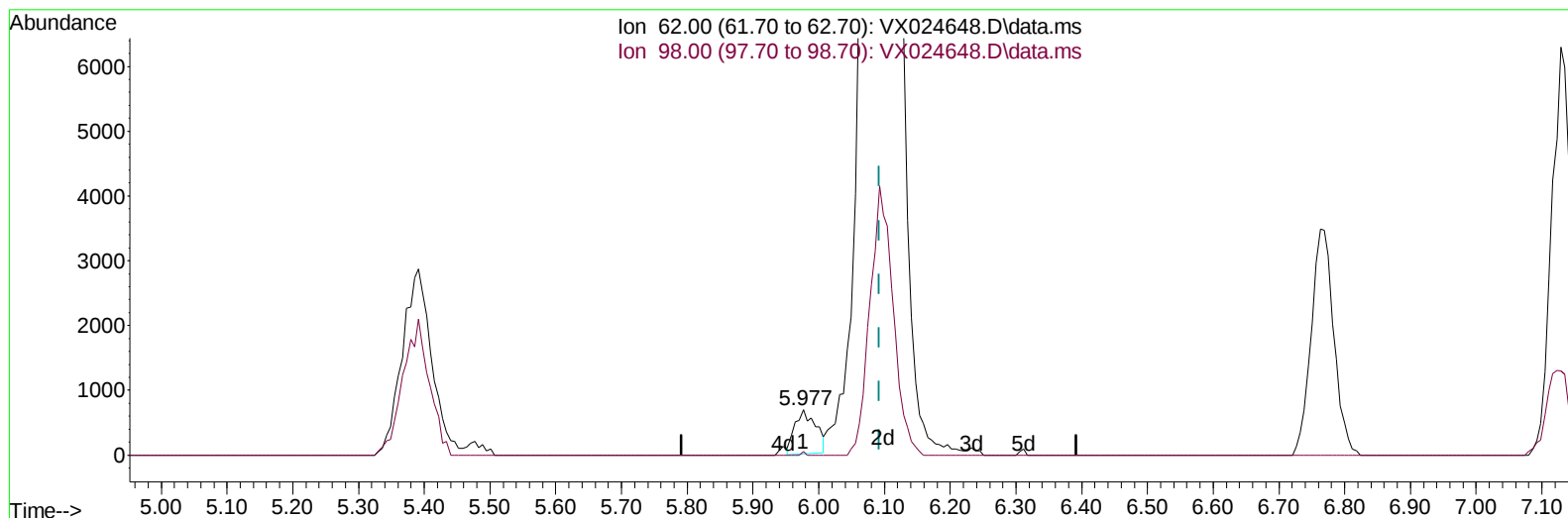
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100721\
Data File : VX024648.D
Acq On : 07 Oct 2021 16:13
Operator : JC/MD
Sample : M4001-04MS 10X
Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EW918MEMS

Manual Integrations
APPROVED

MMDadoda
10/8/2021 9:38:54 AM

Quant Time: Oct 07 23:08:36 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML100721WMA.M
Quant Title : VOC Analysis
QLast Update : Thu Oct 07 23:05:28 2021
Response via : Initial Calibration



TIC: VX024648.D\data.ms

(27) 1,2-Dichloroethane (T)

5.976min (-0.116) 0.65 ug/L

response 1505

Ion	Exp%	Act%
62.00	100.00	100.00
98.00	9.90	8.25
0.00	0.00	0.00
0.00	0.00	0.00

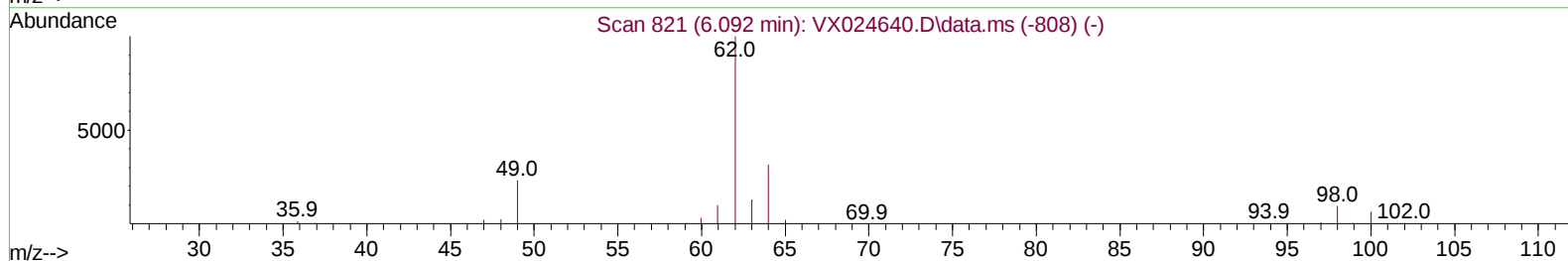
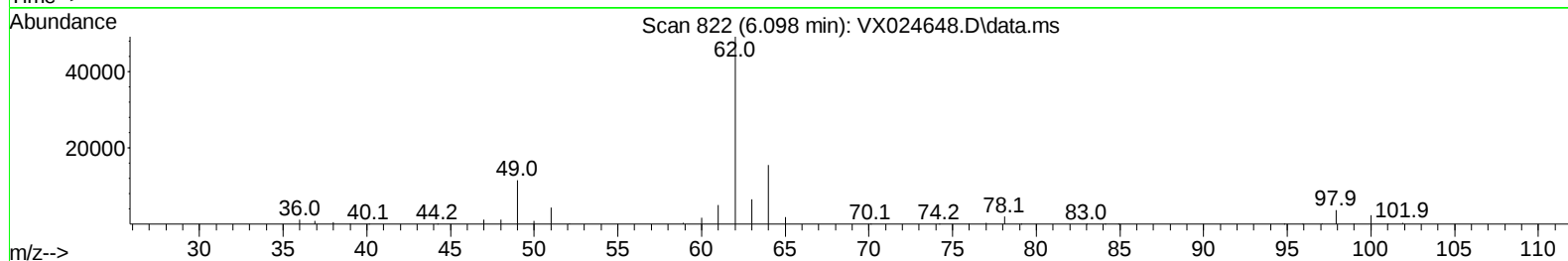
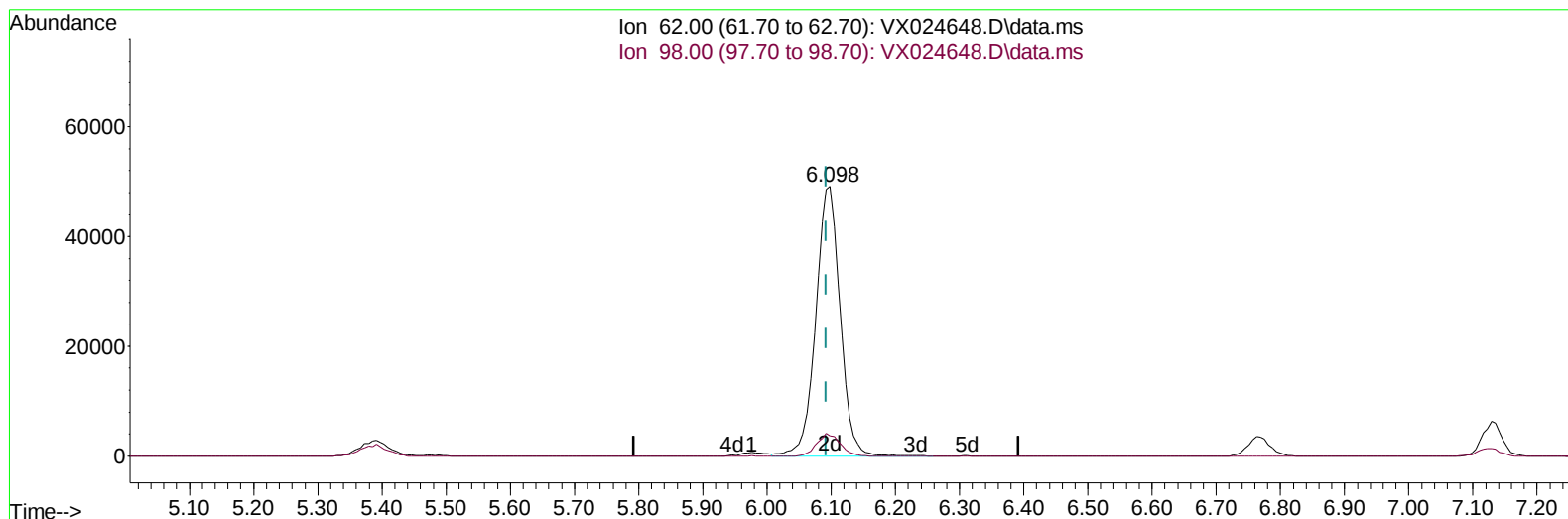
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100721\
Data File : VX024648.D
Acq On : 07 Oct 2021 16:13
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Manual Integrations
APPROVED

MMDadoda
10/8/2021 9:38:54 AM

Quant Time: Oct 07 23:08:36 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML100721WMA.M
Quant Title : VOC Analysis
QLast Update : Thu Oct 07 23:05:28 2021
Response via : Initial Calibration



TIC: VX024648.D\data.ms

(27) 1,2-Dichloroethane (T)

6.098min (+ 0.006) 55.78 ug/L m

response 129166

Ion	Exp%	Act%
62.00	100.00	100.00
98.00	9.90	7.54#
0.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100721\
 Data File : VX024648.D
 Acq On : 07 Oct 2021 16:13
 Operator : JC/MD
 Sample : M4001-04MS 10X
 Misc : 5.04g/5.0mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
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Manual Integrations
 APPROVED

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Quant Time: Oct 07 23:08:36 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM100721WMA.M
 Quant Title : VOC Analysis
 QLast Update : Thu Oct 07 23:05:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Difluorobenzene	6.769	114	216948	50.000	ug/L	0.00
28) Chlorobenzene-d5	10.055	117	198311	50.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	12.024	152	104352	50.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.368	65	57189	42.575	ug/L	0.00
Spiked Amount 50.000	Range 60	- 135	Recovery	=	85.160%	
7) Chloroethane-d5	1.648	69	44535	50.366	ug/L	-0.01
Spiked Amount 50.000	Range 70	- 130	Recovery	=	100.740%	
11) 1,1-Dichloroethene-d2	2.300	63	129030	48.924	ug/L	0.00
Spiked Amount 50.000	Range 60	- 125	Recovery	=	97.840%	
21) 2-Butanone-d5	4.471	46	117619	106.508	ug/L	0.01
Spiked Amount 100.000	Range 40	- 130	Recovery	=	106.510%	
24) Chloroform-d	5.068	84	139383	51.235	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recovery	=	102.480%	
26) 1,2-Dichloroethane-d4	5.964	65	98527	51.209	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recovery	=	102.420%	
32) Benzene-d6	5.983	84	261956	46.892	ug/L	0.00
Spiked Amount 50.000	Range 70	- 125	Recovery	=	93.780%	
36) 1,2-Dichloropropane-d6	7.318	67	83209	47.623	ug/L	0.00
Spiked Amount 50.000	Range 70	- 120	Recovery	=	95.240%	
41) Toluene-d8	8.653	98	232527	42.723	ug/L	0.00
Spiked Amount 50.000	Range 80	- 120	Recovery	=	85.440%	
43) trans-1,3-Dichloroprop...	8.952	79	39679	45.832	ug/L	0.00
Spiked Amount 50.000	Range 60	- 125	Recovery	=	91.660%	
47) 2-Hexanone-d5	9.391	63	79843	97.337	ug/L	0.00
Spiked Amount 100.000	Range 45	- 130	Recovery	=	97.340%	
56) 1,1,2,2-Tetrachloroeth...	11.195	84	122165	50.767	ug/L	0.00
Spiked Amount 50.000	Range 65	- 120	Recovery	=	101.540%	
66) 1,2-Dichlorobenzene-d4	12.323	152	97585	47.795	ug/L	0.00
Spiked Amount 50.000	Range 80	- 120	Recovery	=	95.580%	
Target Compounds						
						Qvalue
2) Dichlorodifluoromethane	1.166	85	93674	48.843	ug/L	99
3) Chloromethane	1.294	50	75025	47.913	ug/L	100
5) Vinyl chloride	1.374	62	77716	49.194	ug/L	99
6) Bromomethane	1.599	94	46170	52.962	ug/L	99
8) Chloroethane	1.672	64	45318	54.949	ug/L	98
9) Trichlorofluoromethane	1.880	101	123564	55.494	ug/L	98
10) 1,1,2-Trichloro-1,2,2-...	2.325	101	74965	54.096	ug/L	96
12) 1,1-Dichloroethene	2.313	96	66969	53.744	ug/L	90
13) Acetone	2.392	43	97299	89.135	ug/L	91
14) Carbon disulfide	2.508	76	174025	48.373	ug/L	99
15) Methyl Acetate	2.709	43	90179	51.851	ug/L	93
16) Methylene chloride	2.788	84	77756	51.863	ug/L	87
17) trans-1,2-Dichloroethene	3.093	96	71877	50.877	ug/L	92
18) Methyl tert-butyl Ether	3.123	73	245402	52.812	ug/L	96
19) 1,1-Dichloroethane	3.611	63	136767	51.812	ug/L	97
20) cis-1,2-Dichloroethene	4.495	96	81354	51.021	ug/L	96
22) 2-Butanone	4.574	43	136685	94.590	ug/L	92
23) Bromochloromethane	4.910	128	43011	51.616	ug/L	96

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Instrument :
 MSVOA_X
 ClientSampleId :
 EW918MEMS

Manual Integrations
 APPROVED

MMDadoda
 10/8/2021 9:38:54 AM

Quant Time: Oct 07 23:08:36 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM100721WMA.M
 Quant Title : VOC Analysis
 QLast Update : Thu Oct 07 23:05:28 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
25) Chloroform	5.105	83	151416	50.731	ug/L	97
27) 1,2-Dichloroethane	6.098	62	129166m	55.778	ug/L	
29) Cyclohexane	5.477	56	124225	50.469	ug/L	97
30) 1,1,1-Trichloroethane	5.391	97	134487	51.715	ug/L	100
31) Carbon tetrachloride	5.684	117	117346	51.433	ug/L	99
33) Benzene	6.044	78	309608	50.524	ug/L	100
34) Trichloroethene	7.129	95	84660	50.193	ug/L	98
35) Methylcyclohexane	7.385	83	132051	49.246	ug/L	98
37) 1,2-Dichloropropane	7.434	63	78960	48.925	ug/L	99
38) Bromodichloromethane	7.830	83	104909	50.886	ug/L	98
39) cis-1,3-Dichloropropene	8.372	75	119603	49.426	ug/L	100
40) 4-Methyl-2-pentanone	8.580	43	251683	102.486	ug/L	91
42) Toluene	8.720	91	325348	47.567	ug/L	99
44) trans-1,3-Dichloropropene	8.982	75	120790	51.107	ug/L	99
45) 1,1,2-Trichloroethane	9.159	97	78294	49.895	ug/L	96
46) Tetrachloroethene	9.275	164	64474	49.038	ug/L	95
48) 2-Hexanone	9.433	43	200149	99.910	ug/L #	90
49) Dibromochloromethane	9.525	129	79886	48.955	ug/L	97
50) 1,2-Dibromoethane	9.616	107	83502	49.273	ug/L	95
51) Chlorobenzene	10.079	112	209965	50.932	ug/L	98
52) Ethylbenzene	10.195	91	361932	51.392	ug/L	99
53) m,p-Xylene	10.305	106	143440	52.133	ug/L	94
54) o-Xylene	10.646	106	140137	50.992	ug/L	96
55) Styrene	10.659	104	242658	52.898	ug/L	94
57) 1,1,2,2-Tetrachloroethane	11.213	83	127418	51.376	ug/L	100
59) Bromoform	10.805	173	57919	49.055	ug/L	99
60) Isopropylbenzene	10.963	105	366145	50.953	ug/L	99
61) 1,2,3-Trichloropropane	11.244	75	107703	50.285	ug/L	99
62) 1,3,5-Trimethylbenzene	11.457	105	323832	51.308	ug/L	99
63) 1,2,4-Trimethylbenzene	11.756	105	320104	51.882	ug/L	98
64) 1,3-Dichlorobenzene	11.969	146	168856	50.388	ug/L	95
65) 1,4-Dichlorobenzene	12.042	146	170755	51.011	ug/L	99
67) 1,2-Dichlorobenzene	12.335	146	172546	52.148	ug/L	96
68) 1,2-Dibromo-3-chloropr...	12.945	75	30364	52.738	ug/L	97
69) 1,3,5-Trichlorobenzene	13.115	180	125252	52.034	ug/L	99
70) 1,2,4-trichlorobenzene	13.591	180	108401	53.225	ug/L	98
71) Naphthalene	13.780	128	388011	57.072	ug/L	99
72) 1,2,3-Trichlorobenzene	13.963	180	115125	54.780	ug/L	99

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

Instrument :
MSVOA_X
ClientSampleId :
EW918MEMS

MMDadoda
10/8/2021 9:38:54 AM

TIC: VX024648.D\data.ms

Abundance

Time-->

1,1-Dichloroethane, T
1,1,1-Trichloroethane, T
1,1,2-Dichloroethane, T
1,1,2,2-Tetrachloroethane, T
1,1,2,2,3-Pentachloroethane, T
1,1,2,2,3,3-Hexachloroethane, T
1,1,2,2,3,3,4-Heptachloroethane, T
1,1,2,2,3,3,4,4-Octachloroethane, T
1,1,2,2,3,3,4,4,5-Nonachloroethane, T
1,1,2,2,3,3,4,4,5,5-Decachloroethane, T
1,1,2,2,3,3,4,4,5,5,6-Undecachloroethane, T
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1,1,2,2,3,3,4,4,5,5,6,6,7-Trichloroethane, T
1,1,2,2,3,3,4,4,5,5,6,6,7,7-Tetrachloroethane, T
1,1,2,2,3,3,4,4,5,5,6,6,7,7,8-Octachloroethane, T
1,1,2,2,3,3,4,4,5,5,6,6,7,7,8,8-Nonachloroethane, T
1,1,2,2,3,3,4,4,5,5,6,6,7,7,8,8,9-Decachloroethane, T
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1,1,2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,20,20,21,21,22,22,23,23-Nonachloroethane, T
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1,1,2,2,3,3,4,4,5,5,6,6,7,7,8,8,9,9,10,10,11,11,12,12,13,13,14,14,15,15,16,16,17,17,18,18,19,19,20,20,21,21,22,22,23,23,24,24,25,25,26-Tetrachloroethane, T
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