

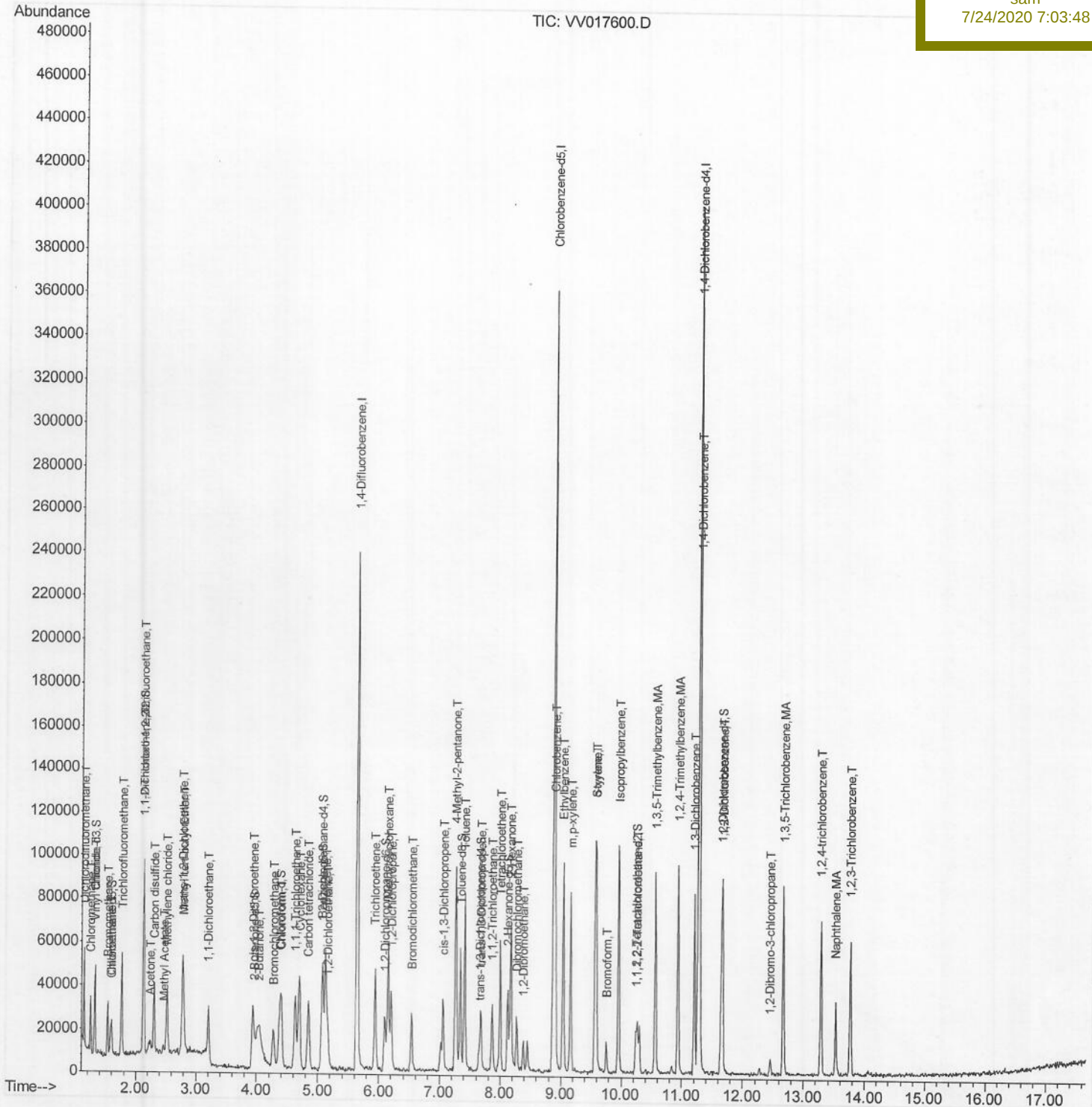
Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV072320\  
Data File : VV017600.D  
Acq On : 23 Jul 2020 12:49  
Operator : SY/MD  
Sample : VSTD00131  
Misc : 25.0mL/MSVOA V/WATER  
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jul 24 01:32:13 2020  
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVTR072320WMA.M  
Quant Title : TRACE VOA SOM01.0  
QLast Update : Fri Jul 24 01:03:26 2020  
Response via : Initial Calibration

**Instrument :**  
MSVOA\_V  
**ClientSampleId :**  
VSTD00131

## Manual Integrations APPROVED

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# Quantitation Report (Qedit)

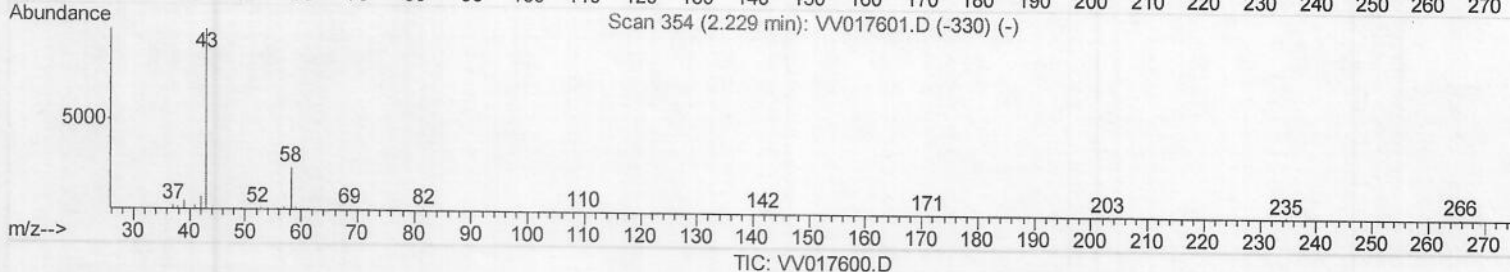
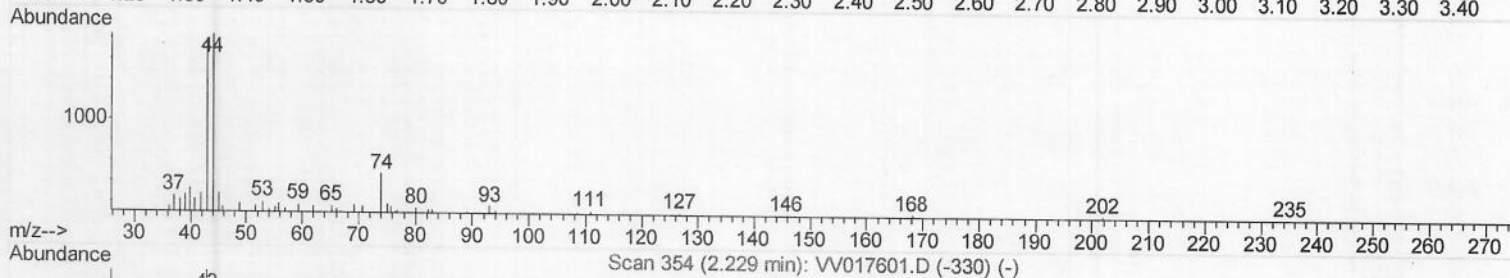
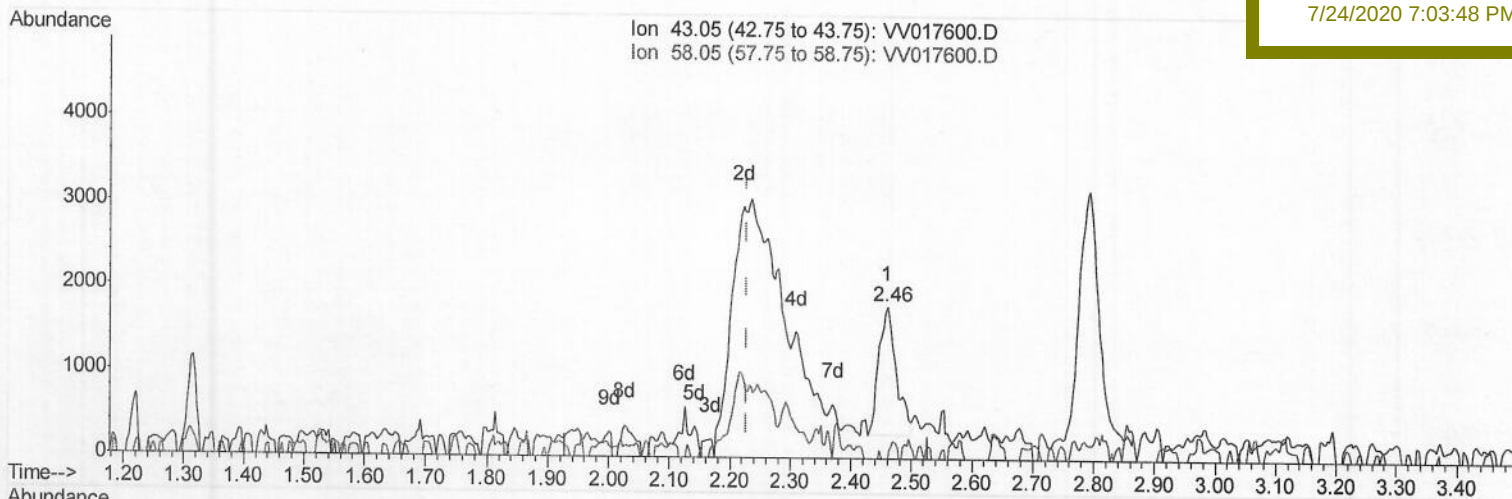
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Manual Integrations  
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(13) Acetone (T)

2.460min (+0.231) 1.70ug/L

response 3032

| Ion   | Exp% | Act% |
|-------|------|------|
| 43.05 | 100  | 100  |
| 58.05 | 9.30 | 5.01 |
| 0.00  | 0.00 | 0.00 |
| 0.00  | 0.00 | 0.00 |

# Quantitation Report (Qedit)

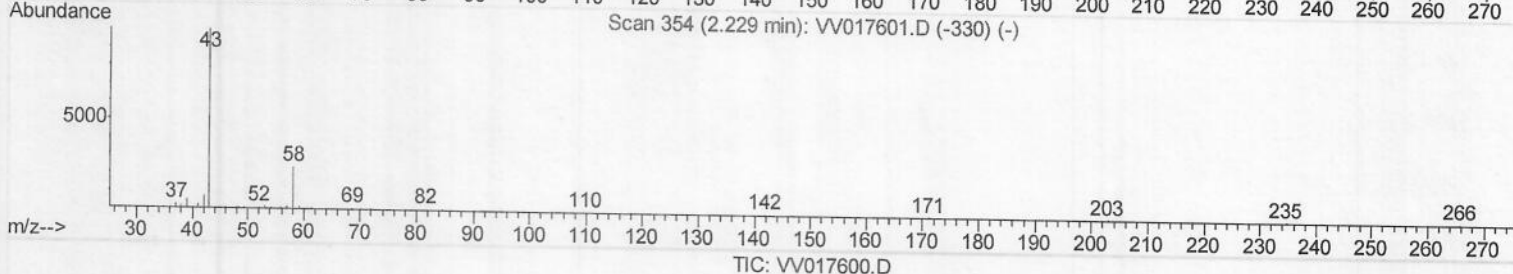
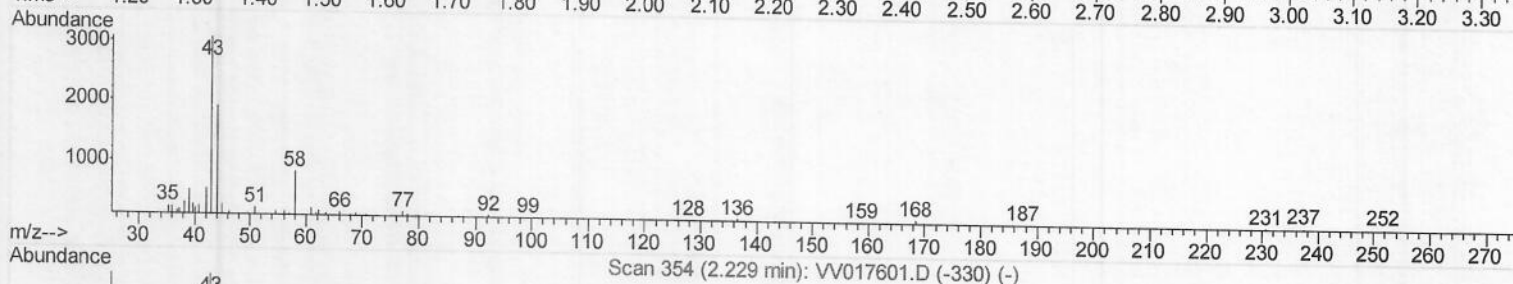
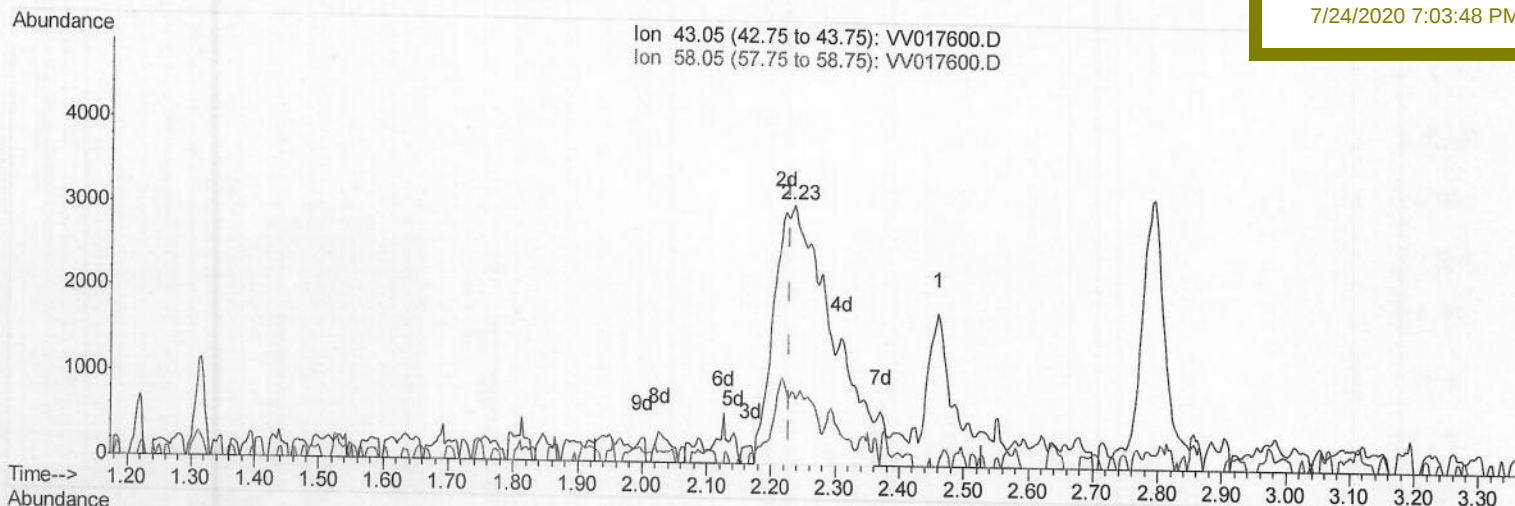
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 ALS Vial : 6 Sample Multiplier: 1

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 QLast Update : Fri Jul 24 01:03:26 2020  
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Instrument :  
 MSVOA\_V  
 ClientSampled :  
 VSTD00131

Manual Integrations  
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(13) Acetone (T)

2.235min (+0.006) 10.63ug/L m

response 18918

| Ion   | Exp% | Act% |
|-------|------|------|
| 43.05 | 100  | 100  |
| 58.05 | 9.30 | 0.80 |
| 0.00  | 0.00 | 0.00 |
| 0.00  | 0.00 | 0.00 |

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# Quantitation Report (Qedit)

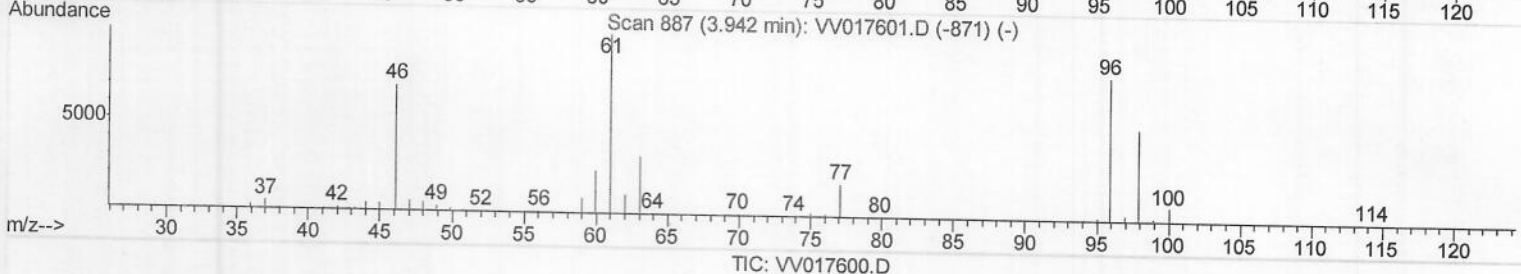
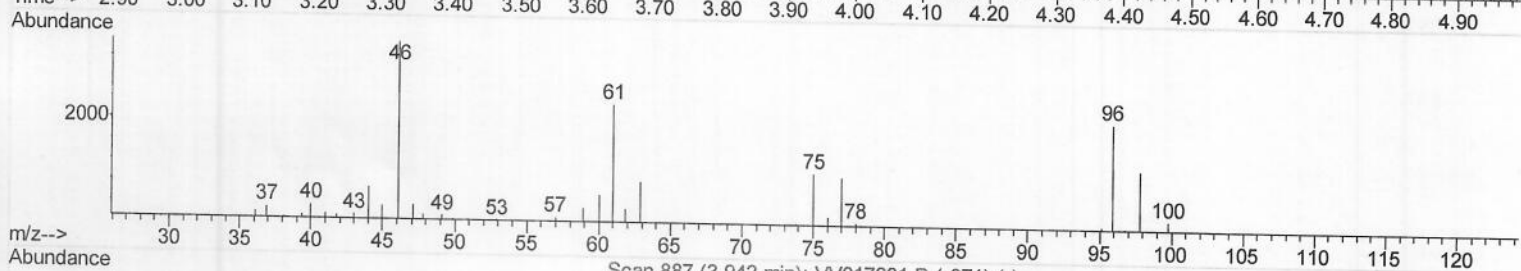
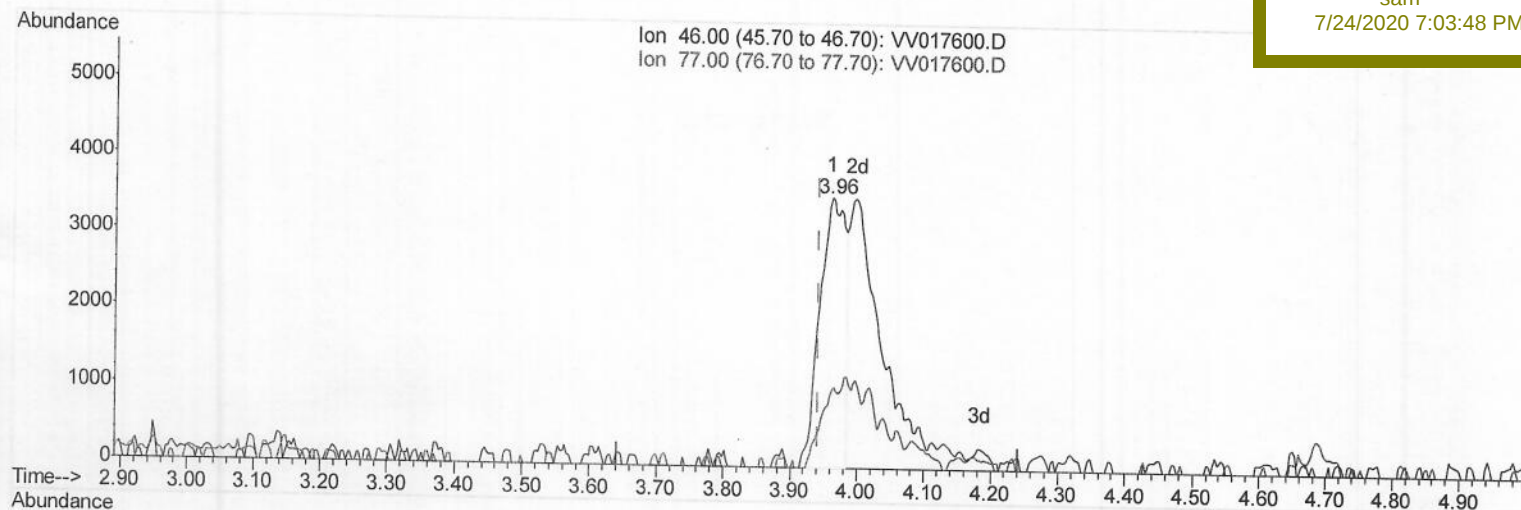
Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV072320\  
 Data File : VV017600.D  
 Acq On : 23 Jul 2020 12:49  
 Operator : SY/MD  
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 ClientSampleId :  
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Manual Integrations  
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(20) 2-Butanone-d5 (S)

3.962min (+0.019) 3.98ug/L

response 9383

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 46.00 | 100   | 100    |
| 77.00 | 11.20 | 20.03# |
| 0.00  | 0.00  | 0.00   |
| 0.00  | 0.00  | 0.00   |

# Quantitation Report (Qedit)

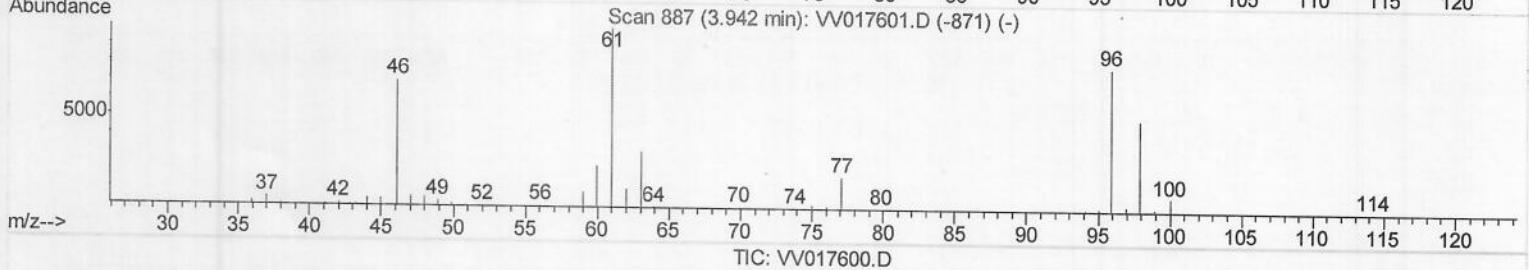
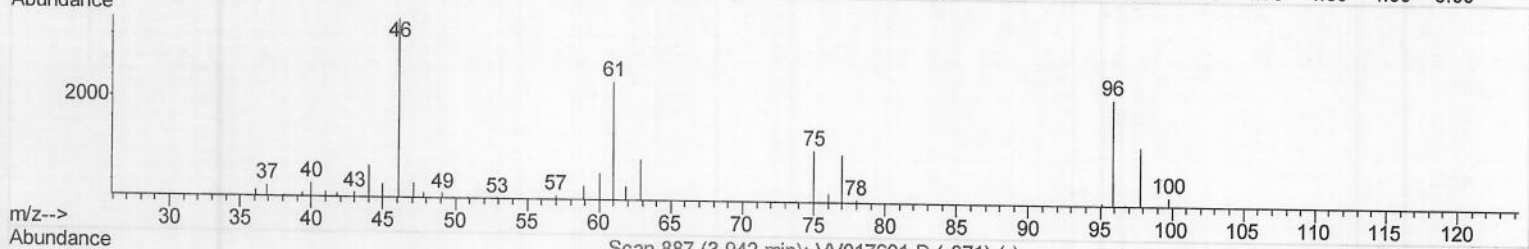
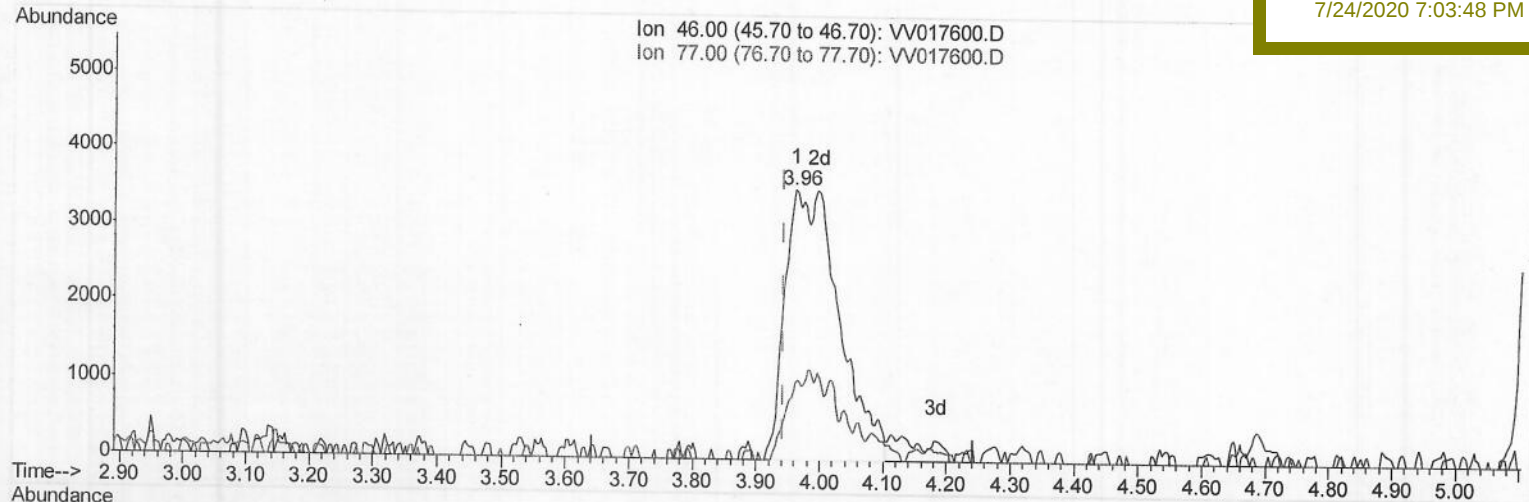
Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV072320\  
 Data File : VV017600.D  
 Acq On : 23 Jul 2020 12:49  
 Operator : SY/MD  
 Sample : VSTD00131  
 Misc : 25.0mL/MSVOA V/WATER  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jul 24 01:29:38 2020  
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 QLast Update : Fri Jul 24 01:03:26 2020  
 Response via : Initial Calibration

Instrument :  
 MSVOA\_V  
 ClientSampleId :  
 VSTD00131

Manual Integrations  
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(20) 2-Butanone-d5 (S)

3.962min (+0.019) 8.83ug/L m

response 20834

| Ion   | Exp%  | Act% |
|-------|-------|------|
| 46.00 | 100   | 100  |
| 77.00 | 11.20 | 9.02 |
| 0.00  | 0.00  | 0.00 |
| 0.00  | 0.00  | 0.00 |

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## Quantitation Report (Qedit)

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV072320\

Data File : VV017600.D

Acq On : 23 Jul 2020 12:49

Operator : SY/MD

Sample : VSTD00131

Misc : 25.0mL/MSVOA V/WATER

ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jul 24 01:29:38 2020

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVTR072320WMA.M

Quant Title : TRACE VOA SOM01.0

QLast Update : Fri Jul 24 01:03:26 2020

Response via : Initial Calibration

Instrument :

MSVOA\_V

ClientSampleId :

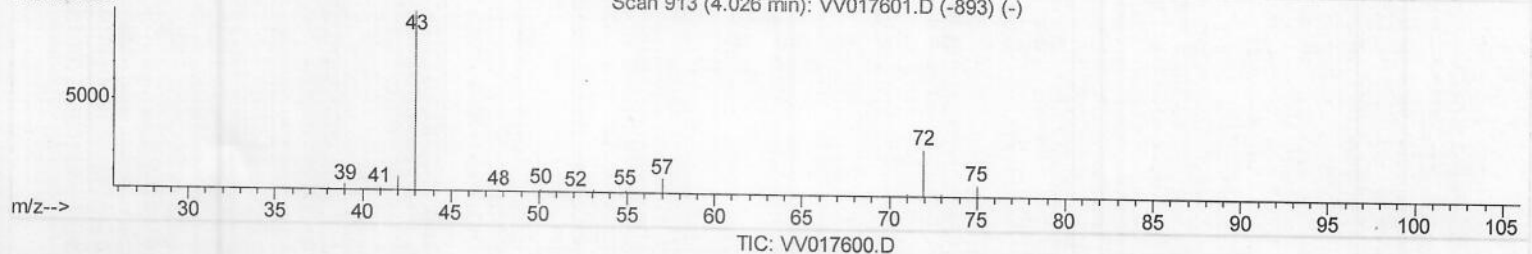
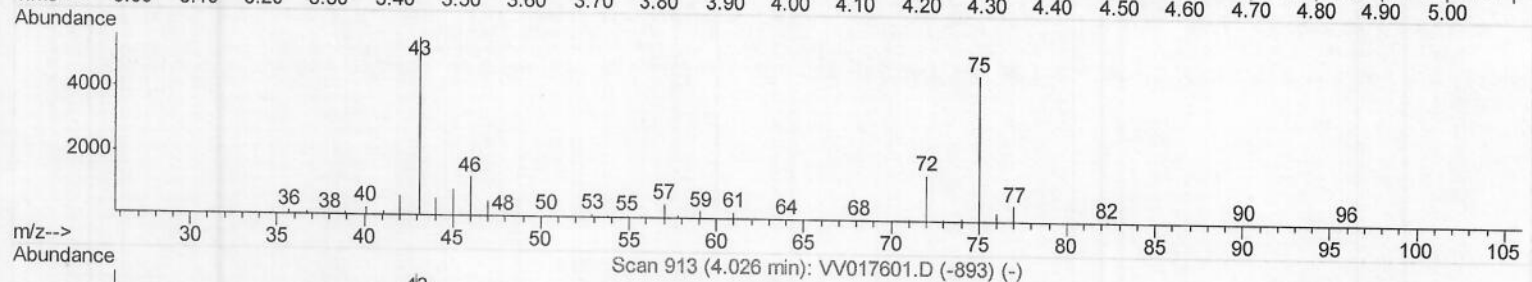
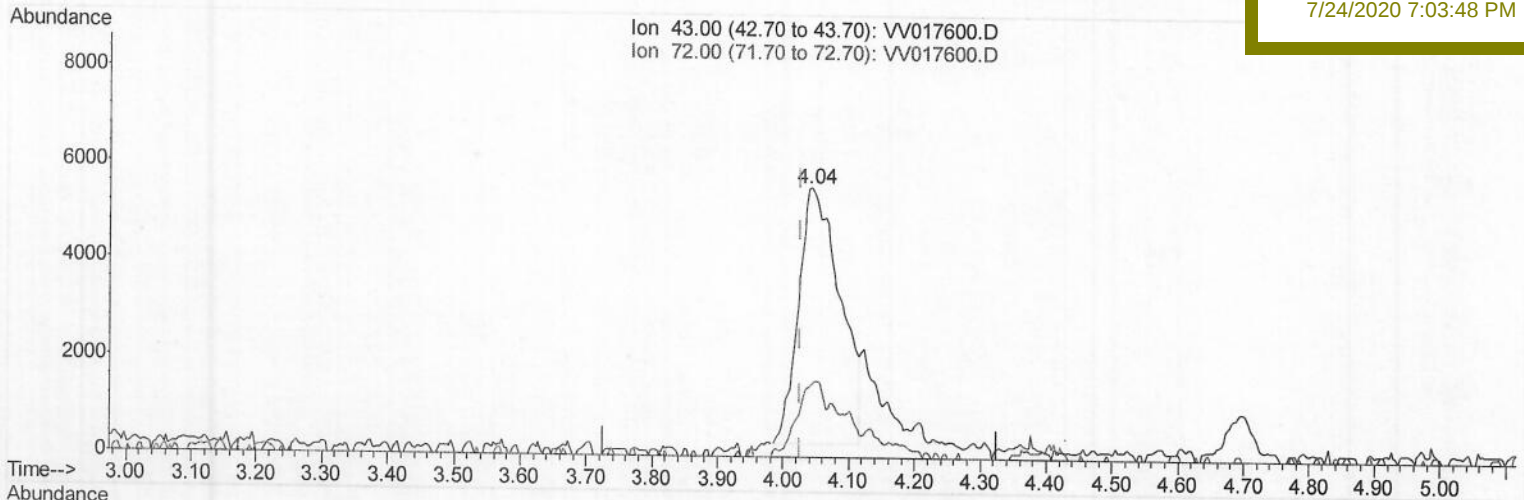
VSTD00131

Manual Integrations

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(21) 2-Butanone (T)

4.042min (+0.016) 8.40ug/L

response 23361

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 43.00 | 100   | 100   |
| 72.00 | 25.90 | 16.49 |
| 0.00  | 0.00  | 0.00  |
| 0.00  | 0.00  | 0.00  |

# Quantitation Report (Qedit)

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV072320\

Data File : VV017600.D

Acq On : 23 Jul 2020 12:49

Operator : SY/MD

Sample : VSTD00131

Misc : 25.0mL/MSVOA V/WATER

ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jul 24 01:29:38 2020

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVTR072320WMA.M

Quant Title : TRACE VOA SOM01.0

QLast Update : Fri Jul 24 01:03:26 2020

Response via : Initial Calibration

Instrument :

MSVOA\_V

Client Sample Id :

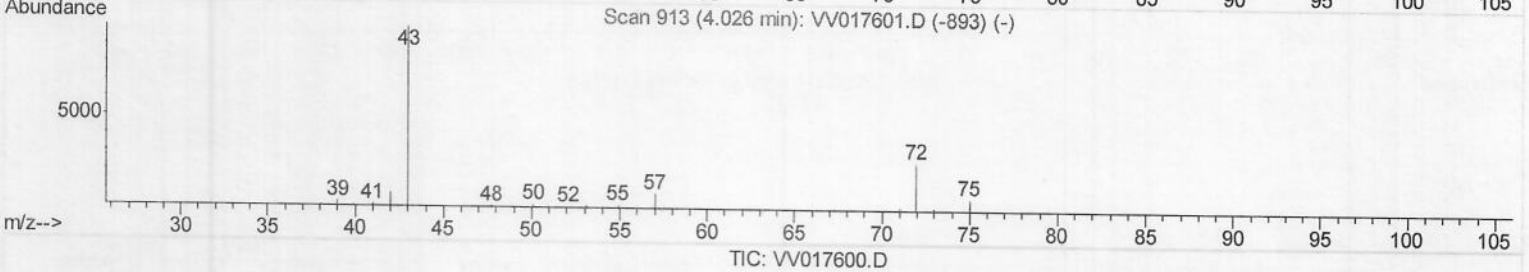
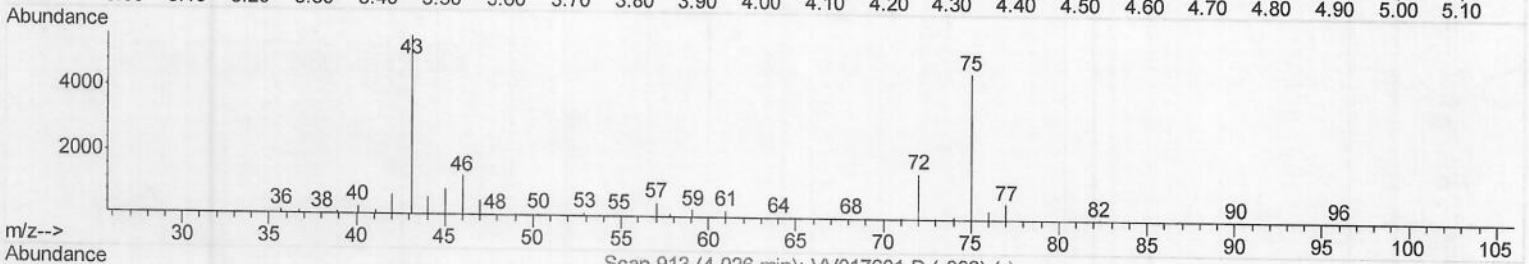
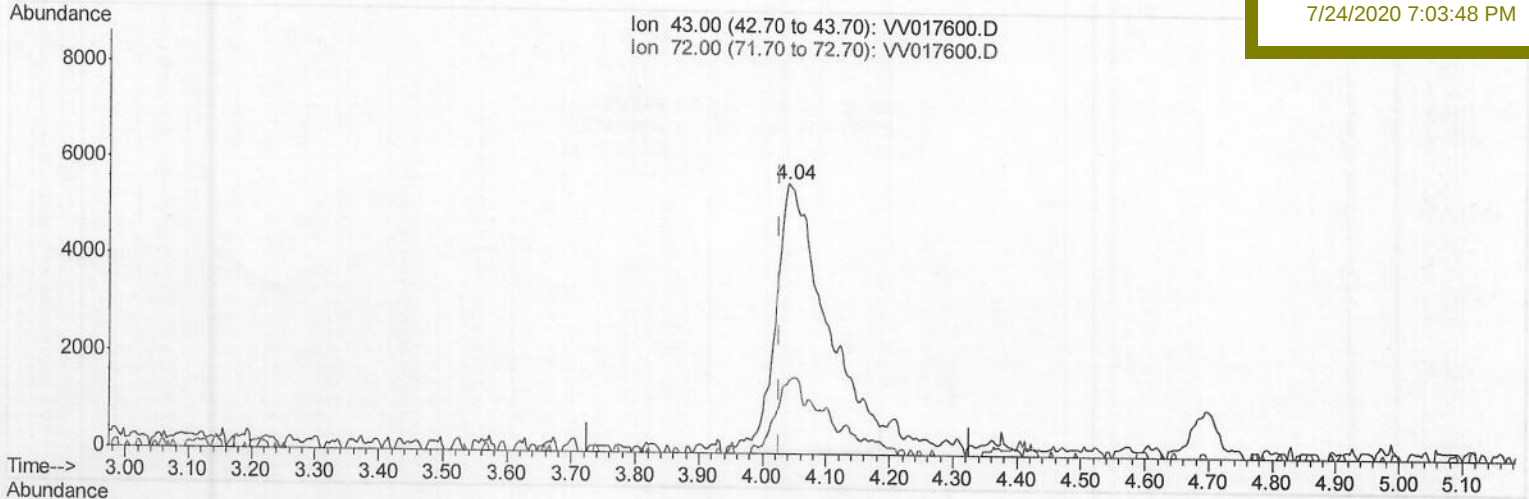
VSTD00131

Manual Integrations

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(21) 2-Butanone (T)

4.042min (+0.016) 10.74ug/L m

response 29855

Ion Exp% Act%

43.00 100 100

72.00 25.90 12.90#

0.00 0.00 0.00

0.00 0.00 0.00

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 Operator : SY/MD  
 Sample : VSTD00131  
 Misc : 25.0mL/MSVOA V/WATER  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 MSVOA\_V  
 Client Sampled :  
 VSTD00131

Quant Time: Jul 24 01:32:13 2020  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA\_V\METHOD\SOMVTR072320WMA.M  
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 Response via : Initial Calibration

Manual Integrations  
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| Internal Standards         | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Difluorobenzene     | 5.64  | 114  | 200353   | 5.00 | ug/L  | 0.00     |
| 28) Chlorobenzene-d5       | 8.87  | 117  | 188139   | 5.00 | ug/L  | 0.00     |
| 61) 1,4-Dichlorobenzene-d4 | 11.27 | 152  | 102114   | 5.00 | ug/L  | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |      |      |
|--------------------------------|-------|-----|--------|------|------|------|
| 4) Vinyl Chloride-d3           | 1.31  | 65  | 9229   | 0.83 | ug/L | 0.00 |
| 7) Chloroethane-d5             | 1.58  | 69  | 6591   | 0.74 | ug/L | 0.00 |
| 11) 1,1-Dichloroethene-d2      | 2.12  | 63  | 20454  | 0.83 | ug/L | 0.00 |
| 20) 2-Butanone-d5              | 3.96  | 46  | 20834m | 8.83 | ug/L | 0.02 |
| 24) Chloroform-d               | 4.38  | 84  | 21125  | 0.83 | ug/L | 0.00 |
| 26) 1,2-Dichloroethane-d4      | 5.06  | 65  | 11732  | 0.88 | ug/L | 0.00 |
| 32) Benzene-d6                 | 5.07  | 84  | 39152  | 0.93 | ug/L | 0.00 |
| 36) 1,2-Dichloropropane-d6     | 6.10  | 67  | 11063  | 0.94 | ug/L | 0.00 |
| 41) Toluene-d8                 | 7.34  | 98  | 36122  | 0.88 | ug/L | 0.00 |
| 45) trans-1,3-Dichloropropene- | 7.65  | 79  | 4795   | 0.85 | ug/L | 0.00 |
| 48) 2-Hexanone-d5              | 8.12  | 63  | 12557  | 7.40 | ug/L | 0.00 |
| 59) 1,1,2,2-Tetrachloroethane- | 10.24 | 84  | 7480   | 0.89 | ug/L | 0.00 |
| 65) 1,2-Dichlorobenzene-d4     | 11.65 | 152 | 14166  | 0.84 | ug/L | 0.00 |

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## Target Compounds

|                                |      |     |        |        |        | Qvalue |
|--------------------------------|------|-----|--------|--------|--------|--------|
| 2) Dichlorodifluoromethane     | 1.14 | 85  | 22748  | 0.971  | ug/L   | 96     |
| 3) Chloromethane               | 1.24 | 50  | 14607  | 0.927  | ug/L   | 95     |
| 5) Vinyl chloride              | 1.32 | 62  | 14256  | 0.901  | ug/L   | 99     |
| 6) Bromomethane                | 1.53 | 94  | 9530   | 0.985  | ug/L   | 97     |
| 8) Chloroethane                | 1.59 | 64  | 8726   | 0.955  | ug/L   | 96     |
| 9) Trichlorofluoromethane      | 1.76 | 101 | 26186  | 0.907  | ug/L   | 100    |
| 10) 1,1,2-Trichloro-1,2,2-trif | 2.13 | 101 | 12022  | 0.895  | ug/L   | 98     |
| 12) 1,1-Dichloroethene         | 2.13 | 96  | 10977  | 0.924  | ug/L   | 97     |
| 13) Acetone                    | 2.23 | 43  | 18918m | 10.635 | ug/L   | 97     |
| 14) Carbon disulfide           | 2.31 | 76  | 32006  | 0.881  | ug/L # | 94     |
| 15) Methyl Acetate             | 2.46 | 43  | 3032   | 0.856  | ug/L   | 95     |
| 16) Methylene chloride         | 2.52 | 84  | 12446  | 0.871  | ug/L   | 93     |
| 17) Methyl tert-butyl Ether    | 2.79 | 73  | 26379  | 0.887  | ug/L   | 95     |
| 18) trans-1,2-Dichloroethene   | 2.77 | 96  | 11780  | 0.930  | ug/L   | 88     |
| 19) 1,1-Dichloroethane         | 3.21 | 63  | 23841  | 0.957  | ug/L   | 99     |
| 21) 2-Butanone                 | 4.04 | 43  | 29855m | 10.739 | ug/L   | 97     |
| 22) cis-1,2-Dichloroethene     | 3.93 | 96  | 15034  | 1.056  | ug/L   | 84     |
| 23) Bromochloromethane         | 4.27 | 128 | 7107   | 1.115  | ug/L   | 91     |
| 25) Chloroform                 | 4.40 | 83  | 28292  | 1.046  | ug/L   | 87     |
| 27) 1,2-Dichloroethane         | 5.16 | 62  | 17411  | 0.983  | ug/L # | 94     |
| 29) 1,1,1-Trichloroethane      | 4.63 | 97  | 26561  | 1.011  | ug/L   | 96     |
| 30) Cyclohexane                | 4.69 | 56  | 20967  | 1.027  | ug/L   | 99     |
| 31) Carbon tetrachloride       | 4.85 | 117 | 23650  | 0.978  | ug/L   | 100    |
| 33) Benzene                    | 5.12 | 78  | 52167  | 1.041  | ug/L   | 100    |
| 34) Trichloroethene            | 5.94 | 95  | 15334  | 1.019  | ug/L   | 95     |
| 35) Methylcyclohexane          | 6.15 | 83  | 21834  | 0.968  | ug/L   | 96     |
| 37) 1,2-Dichloropropane        | 6.20 | 63  | 12335  | 1.049  | ug/L   | 98     |
| 38) Bromodichloromethane       | 6.53 | 83  | 18645  | 1.048  | ug/L   | 98     |
| 39) cis-1,3-Dichloropropene    | 7.05 | 75  | 16892  | 0.918  | ug/L   | 91     |
| 40) 4-Methyl-2-pentanone       | 7.25 | 43  | 63437  | 10.133 | ug/L   | 97     |

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 Response via : Initial Calibration

Instrument :  
 MSVOA\_V  
 Client Sampled :  
 VSTD00131

Manual Integrations  
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| Internal Standards             | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 42) Toluene                    | 7.41  | 91   | 58169    | 1.054 | ug/L  | 97       |
| 43) 1,3,5-Trimethylbenzene     | 10.56 | 105  | 48738    | 0.899 | ug/L  | 98       |
| 44) 1,2,4-Trimethylbenzene     | 10.94 | 105  | 49655    | 0.916 | ug/L  | 98       |
| 46) trans-1,3-Dichloropropene  | 7.68  | 75   | 15801    | 0.968 | ug/L  | 99       |
| 47) 1,1,2-Trichloroethane      | 7.86  | 97   | 9455     | 1.072 | ug/L  | 91       |
| 49) Tetrachloroethene          | 7.99  | 164  | 13226    | 1.030 | ug/L  | 97       |
| 50) 2-Hexanone                 | 8.17  | 43   | 45899    | 9.967 | ug/L  | 98       |
| 51) Dibromochloromethane       | 8.27  | 129  | 11406    | 0.959 | ug/L  | 93       |
| 52) 1,2-Dibromoethane          | 8.38  | 107  | 8930     | 1.039 | ug/L  | 93       |
| 53) Chlorobenzene              | 8.90  | 112  | 39112    | 1.060 | ug/L  | 98       |
| 54) Ethylbenzene               | 9.03  | 91   | 62524    | 0.995 | ug/L  | 96       |
| 55) m,p-xylene                 | 9.16  | 106  | 22918    | 0.962 | ug/L  | 92       |
| 56) o-xylene                   | 9.56  | 106  | 23067    | 0.995 | ug/L  | 91       |
| 57) Styrene                    | 9.58  | 104  | 37302    | 0.963 | ug/L  | 97       |
| 58) Isopropylbenzene           | 9.95  | 105  | 62884    | 0.971 | ug/L  | 99       |
| 60) 1,1,2,2-Tetrachloroethane  | 10.26 | 83   | 10121    | 1.118 | ug/L  | 96       |
| 62) Bromoform                  | 9.75  | 173  | 6048     | 0.907 | ug/L  | 97       |
| 63) 1,3-Dichlorobenzene        | 11.20 | 146  | 32689    | 1.024 | ug/L  | 95       |
| 64) 1,4-Dichlorobenzene        | 11.29 | 146  | 33368    | 1.051 | ug/L  | 97       |
| 66) 1,2-Dichlorobenzene        | 11.67 | 146  | 29729    | 1.021 | ug/L  | 97       |
| 67) 1,2-Dibromo-3-chloropropan | 12.45 | 75   | 1750     | 0.961 | ug/L  | 91       |
| 68) 1,3,5-Trichlorobenzene     | 12.67 | 180  | 25718    | 0.993 | ug/L  | 97       |
| 69) 1,2,4-trichlorobenzene     | 13.29 | 180  | 20914    | 0.973 | ug/L  | 98       |
| 70) Naphthalene                | 13.53 | 128  | 26370    | 0.866 | ug/L  | 98       |
| 71) 1,2,3-Trichlorobenzene     | 13.77 | 180  | 18556    | 0.999 | ug/L  | 99       |

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