Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV112321\

Data File: VV023684.D

Acq On : 23 Nov 2021 20:30

Operator : SY/MD Sample : M4723-14

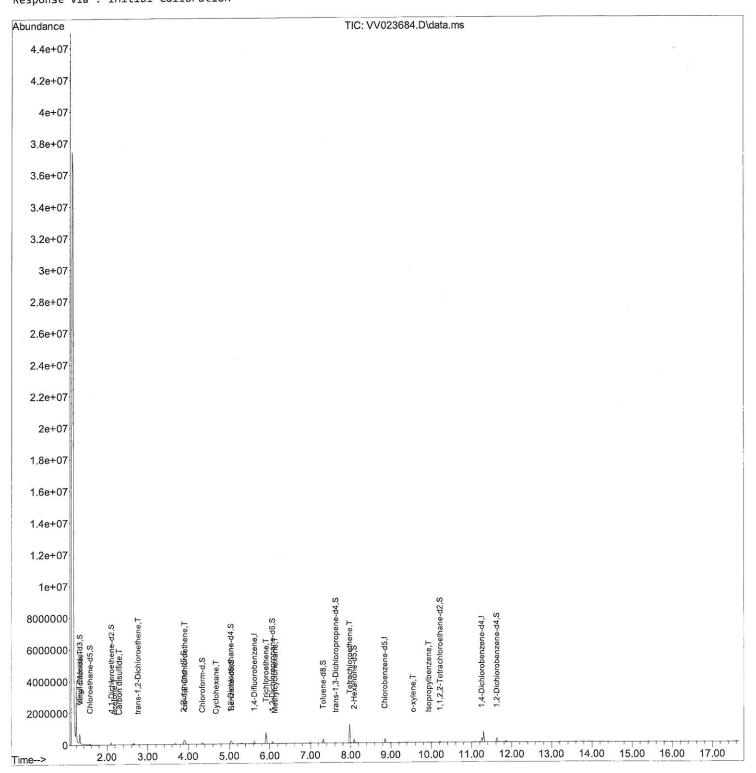
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Nov 24 05:01:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR112321WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 24 04:42:45 2021 Response via : Initial Calibration Instrument : MSVOA\_V ClientSampleId : C0G43

# **Manual IntegrationsAPPROVED**



#### Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV112321\

Data File: VV023684.D

Acq On : 23 Nov 2021 20:30

Operator : SY/MD Sample : M4723-14

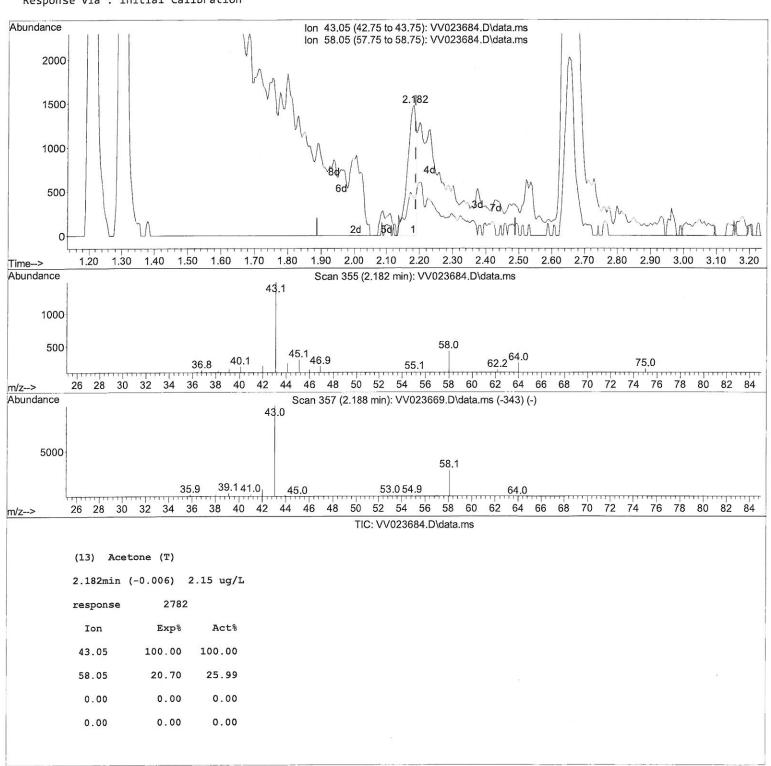
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Nov 24 05:01:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR112321WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 24 04:42:45 2021 Response via : Initial Calibration Instrument : MSVOA\_V ClientSampleId : C0G43

# **Manual Integrations APPROVED**



### Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV112321\

Data File: VV023684.D

Acq On : 23 Nov 2021 20:30

Operator : SY/MD Sample : M4723-14

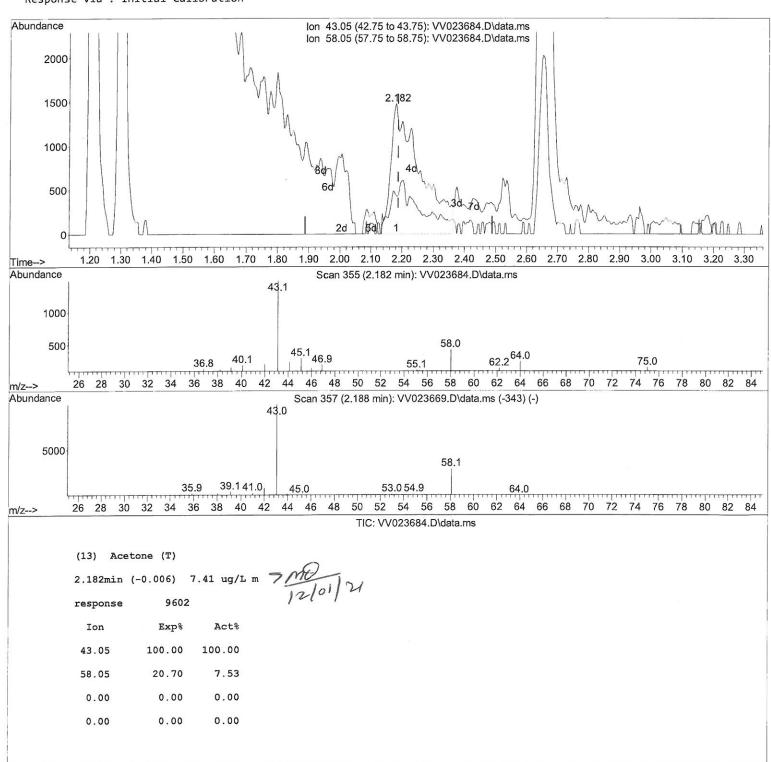
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Nov 24 05:01:51 2021

 $\label{thm:local_var_var} \mbox{Quant Method}: \mbox{Z:\voasrv\HPCHEM1\MSVOA\_v\Method\SFAMVTR112321WMA.M}$ 

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 24 04:42:45 2021 Response via : Initial Calibration Instrument : MSVOA\_V ClientSampleId : C0G43

# **Manual IntegrationsAPPROVED**



Data Path : Z:\voasrv\HPCHEM1\MSVOA\_V\Data\VV112321\

Data File : VV023684.D

Acq On : 23 Nov 2021 20:30

Operator : SY/MD Sample : M4723-14

Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Nov 24 05:01:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_V\Method\SFAMVTR112321WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 24 04:42:45 2021 Response via : Initial Calibration Instrument: MSVOA\_V ClientSampleId: C0G43

# **Manual IntegrationsAPPROVED**

Compound	R.T.	QIon	Response	Conc Un	its Dev(	Min)	
Internal Standards							
<ol> <li>1,4-Difluorobenzene</li> </ol>	5.613	114	145886	5.000	ug/L	0.00	
28) Chlorobenzene-d5	8.850	117	144113	5.000	ug/L	0.00	1
58) 1,4-Dichlorobenzene-d4	11.249	152	72923	5.000	ug/L	0.00	ĺ
System Monitoring Compounds							
<ol><li>Vinyl Chloride-d3</li></ol>	1.301	65	52024	4.344	ug/L	0.00	
Spiked Amount 5.000	Range 40	- 130	Recover	ry =	86.800%		
7) Chloroethane-d5	1.561	69	42922	4.559	ug/L	0.00	
Spiked Amount 5.000	Range 65	- 130	Recover	-y =	91.200%		
<pre>11) 1,1-Dichloroethene-d2</pre>	2.101	63	69663	3.301	ug/L	0.00	
Spiked Amount 5.000	Range 60	- 125	Recover	'y =	66.000%		
20) 2-Butanone-d5	3.883	46	96394	66.954	ug/L	-0.02	
Spiked Amount 50.000	Range 40	- 130	Recover	·y =	133.900%	#	
24) Chloroform-d	4.343	84	95213	4.566	ug/L	0.00	
Spiked Amount 5.000	Range 70	- 125	Recover	'y =	91.400%		
26) 1,2-Dichloroethane-d4	5.027	65	46362	4.759	ug/L	0.00	
Spiked Amount 5.000	Range 70	- 130	Recover	·y =	95.200%		
32) Benzene-d6	5.047	84	186268	4.745	ug/L	0.00	
Spiked Amount 5.000	Range 70	- 125	Recover	·y =	94.800%		
36) 1,2-Dichloropropane-d6	6.066	67	54381	4.941	ug/L	0.00	
	Range 60	- 140	Recover	y =	98.800%		
41) Toluene-d8	7.313	98	169554	4.623	ug/L	0.00	
and the second control of the second control	Range 70	- 130	Recover		92.400%		
43) trans-1,3-Dichloroprop	. 7.622	79	20222	4.558	ug/L	0.00	
	Range 55	- 130	Recover	y =	91.200%		
46) 2-Hexanone-d5	8.088	63	80536	54.640	ug/L	0.00	
A STATE OF THE STA	Range 45	- 130	Recover		109.280%		
56) 1,1,2,2-Tetrachloroeth	250	84		4.900		0.00	
•	Range 65	- 120	Recover		98.000%		
66) 1,2-Dichlorobenzene-d4	11.622		68439	5.308		0.00	
	Range 80				106.200%		
Target Compounds					Qva:	lue	
5) Vinyl chloride	1.304	62	17563	1.390	77	88	0
13) Acetone	2.182	43	9602m	7.412	ug/L	7	MOTO
14) Carbon disulfide	2.291	76	3521	0.107		100	12/01/4
18) trans-1,2-Dichloroethene		96	3075	0.276		86	, ,
22) cis-1,2-Dichloroethene	3.905	96	141871	13.288		99	
30) Cyclohexane	4.674	56	2817		ug/L #	88	
34) Trichloroethene	5.908	95	236454	21.484		97	
35) Methylcyclohexane	6.124	83	1737		ug/L #	76	
47) Tetrachloroethene	7.973	164	262791	26.233		98	
54) o-xylene	9.545	106	3432	0.195		92	
60) Isopropylbenzene	9.934	105	16008	0.368		99	

<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed