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

Data File: VV023691.D

Acq On : 23 Nov 2021 23:16

Operator : SY/MD

Sample : VSTDCCC005EC

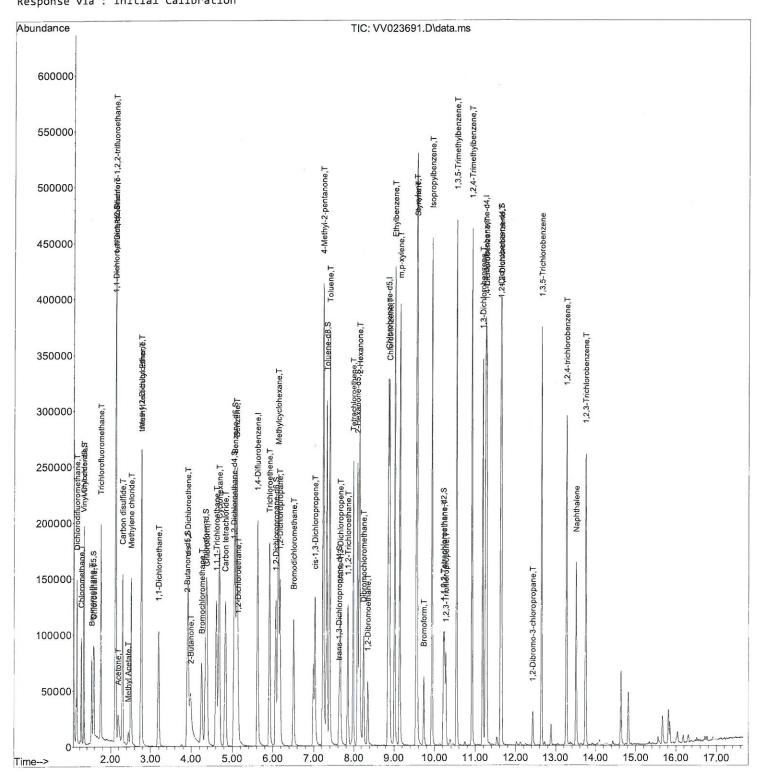
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 24 05:03:37 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
LabSampleId :
VSTDCCC005EC

#### **Manual IntegrationsAPPROVED**



#### Quantitation Report (Qedit)

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

Data File : VV023691.D

Acq On : 23 Nov 2021 23:16

Operator : SY/MD

Sample : VSTDCCC005EC

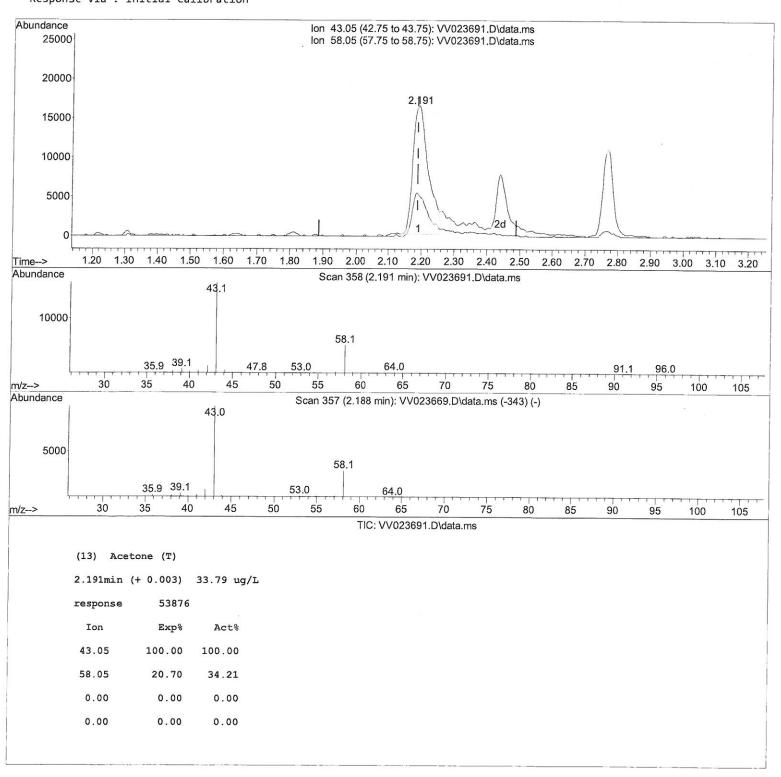
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 24 05:03:37 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
LabSampleId :
VSTDCCC005EC

### **Manual IntegrationsAPPROVED**



### Quantitation Report (Qedit)

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

Data File : VV023691.D

Acq On : 23 Nov 2021 23:16

Operator : SY/MD

Sample : VSTDCCC005EC

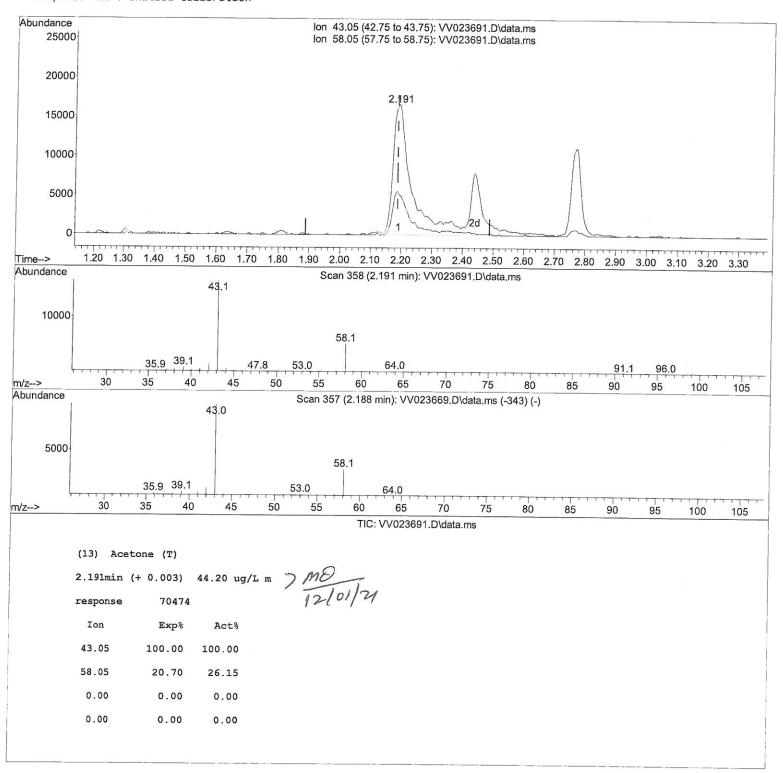
Misc : 25.0mL/MSVOA\_V/WATER
ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 24 05:03:37 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
LabSampleId :
VSTDCCC005EC

#### **Manual IntegrationsAPPROVED**



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

Data File : VV023691.D

Acq On : 23 Nov 2021 23:16
Operator : SY/MD
Sample : VSTDCCC005EC
Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 24 05:03:37 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 LabSampleId : VSTDCCC005EC

## **Manual IntegrationsAPPROVED**

Compound	R.T. C	Ion	Response	Conc Ur	its Dev(	Min)	
Internal Standards		0.00			741		
1) 1,4-Difluorobenzene	5.619		179550		ug/L	0.00	
28) Chlorobenzene-d5	8.854		173348		ug/L	0.00	
58) 1,4-Dichlorobenzene-d4	11.249	152	93778	5.000	ug/L	0.00	
System Monitoring Compounds							
4) Vinyl Chloride-d3	1.307	65	56915	3.861	ug/L	0.00	
Spiked Amount 5.000	Range 40 -		Recover				
7) Chloroethane-d5	1.568	69	45628	3.938	ug/L	0.00	
Spiked Amount 5.000	Range 65 -	130	Recover	'y =	78.800%		
11) 1,1-Dichloroethene-d2	2.111	63	103317	3.977	ug/L	0.00	
Spiked Amount 5.000	Range 60 -	125	Recover	ry =	79.600%		
20) 2-Butanone-d5	3.899	46	101066	57.037	ug/L	0.00	
Spiked Amount 50.000	Range 40 -				114.080%		
24) Chloroform-d	4.349	84	112168		0.0000	0.00	
Spiked Amount 5.000	Range 70 -		Recover				
26) 1,2-Dichloroethane-d4	5.034	65				0.00	
Spiked Amount 5.000	Range 70 -		Recover		88.800%		
32) Benzene-d6	5.053	84	218617		CONTRACTOR DESCRIPTION OF THE PARTY OF THE P	0.00	
Spiked Amount 5.000	Range 70 -	67	Recover	-		0 00	
36) 1,2-Dichloropropane-d6	6.069					0.00	*
Spiked Amount 5.000	Range 60 - 7.317	98	207315	y =		0 00	
41) Toluene-d8 Spiked Amount 5.000	Range 70 -		Recover		50 To 00 to 900 to	0.00	
43) trans-1,3-Dichloroprop		79		4.519		0.00	
Spiked Amount 5.000	Range 55 -		Recover		90.400%	0.00	
46) 2-Hexanone-d5	8.088	63	96310	THE RESERVE THE PARTY OF THE PA		0.00	
Spiked Amount 50.000	Range 45 -				108.640%	0.00	
56) 1,1,2,2-Tetrachloroeth		84				0.00	
	Range 65 -						
66) 1,2-Dichlorobenzene-d4						0.00	
	Range 80 -	120	Recover		94.200%		
Taugat Campunda					0	lua.	
Target Compounds 2) Dichlorodifluoromethane	1.130	85	74064	4.348	Qva]	100	
3) Chloromethane	1.243	50	61542	4.156	0.	98	
5) Vinyl chloride	1.310	62	65999	4.244		96	
6) Bromomethane	1.523	94	28279	3.207	0.	98	
8) Chloroethane		64	42062	4.268		97	
9) Trichlorofluoromethane		101	106566	4.205	1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-1-	99	
10) 1,1,2-Trichloro-1,2,2			53963	4.249		100	
12) 1,1-Dichloroethene	2.121	96	52192	4.339		95	-0
13) Acetone	2.191	43	70474m	44.200		2	ME
14) Carbon disulfide	2.298	76	172710	4.271		100	12/01/21
15) Methyl Acetate	2.439	43	16535	4.569		97	Communication (Communication)
16) Methylene chloride	2.510	84	61281	3.570	ug/L	97	
17) Methyl tert-butyl Ether	2.767	73	121250	4.918	ug/L	99	
18) trans-1,2-Dichloroethene		96	60609	4.423	-	97	
19) 1,1-Dichloroethane	3.191	63	103169	4.478		100	
21) 2-Butanone	3.986	43	97411	48.269	•	93	
22) cis-1,2-Dichloroethene	3.915	96	63294	4.817		98	
23) Bromochloromethane	4.252 1	.28	28442	4.612	ug/L	95	

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

Data File : VV023691.D

Acq On : 23 Nov 2021 23:16

Operator : SY/MD

Sample : VSTDCCC005EC
Misc : 25.0mL/MSVOA\_V/WATER ALS Vial : 31 Sample Multiplier: 1

Quant Time: Nov 24 05:03:37 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 LabSampleId : VSTDCCC005EC

# **Manual IntegrationsAPPROVED**

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)					
25) Chloroform	4.378	83	110461	4.304 ug/L	96					
27) 1,2-Dichloroethane	5.133	62	62206	4.559 ug/L	99					
29) 1,1,1-Trichloroethane	4.609	97	105691	4.666 ug/L	99					
30) Cyclohexane	4.680	56	94250	4.986 ug/L	99					
31) Carbon tetrachloride	4.828	117	97011	4.675 ug/L	100					
33) Benzene	5.101	78	249346	5.045 ug/L	100					
34) Trichloroethene	5.915	95	64152	4.846 ug/L	96					
35) Methylcyclohexane	6.133	83	103442	5.012 ug/L	98					
37) 1,2-Dichloropropane	6.172	63	54802	4.665 ug/L	99					
38) Bromodichloromethane	6.510	83	75102	4.710 ug/L	97					
39) cis-1,3-Dichloropropene	7.031	75	82564	4.939 ug/L	100					
40) 4-Methyl-2-pentanone	7.227	43	285968	50.995 ug/L	99					
42) Toluene	7.387	91	268512	5.010 ug/L	99					
44) trans-1,3-Dichloropropene	7.651	75	68717	4.891 ug/L	98					
45) 1,1,2-Trichloroethane	7.841	97	39505	4.857 ug/L	98					
47) Tetrachloroethene	7.976	164	56959	4.727 ug/L	98					
48) 2-Hexanone	8.140	43	201714	48.668 ug/L	99					
49) Dibromochloromethane	8.246	129	51198	4.593 ug/L	97					
50) 1,2-Dibromoethane	8.352	107	38287	4.827 ug/L	98					
51) Chlorobenzene	8.883	112	172499	4.855 ug/L	99					
52) Ethylbenzene	9.011	91	288812	5.155 ug/L	99					
53) m,p-xylene	9.140	106	112295	5.036 ug/L	97					
54) o-xylene	9.545	106	108510	5.117 ug/L	99					
55) Styrene	9.561	104	180892	5.063 ug/L	97					
57) 1,1,2,2-Tetrachloroethane	10.239	83	44593	4.927 ug/L	99					
59) Bromoform	9.731	173	28583	4.615 ug/L	99					
60) Isopropylbenzene	9.931	105	290909	5.195 ug/L	100					
61) 1,2,3-Trichloropropane	10.275	75	31500	4.736 ug/L	99					
62) 1,3,5-Trimethylbenzene	10.538	105	243726	5.231 ug/L	100					
63) 1,2,4-Trimethylbenzene	10.915	105	246887	5.358 ug/L	100					
64) 1,3-Dichlorobenzene	11.181	146	142286	4.970 ug/L	99					
65) 1,4-Dichlorobenzene	11.272	146	138567	4.816 ug/L	99					
67) 1,2-Dichlorobenzene	11.641	146	128748	4.914 ug/L	99					
68) 1,2-Dibromo-3-chloropr	12.429	75	6828	5.170 ug/L	89					
69) 1,3,5-Trichlorobenzene	12.644	180	112806	5.047 ug/L	99					
70) 1,2,4-trichlorobenzene	13.262	180	89919	5.185 ug/L	99					
71) Naphthalene	13.503	128	129584	5.552 ug/L	99					
72) 1,2,3-Trichlorobenzene	13.744	180	79080	5.262 ug/L	99					

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