Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\

Data File : VW021040.D

Acq On : 02 Dec 2021 12:23

Operator : SY/VA Sample : M4886-12MS

Misc : 5.51g/10.0mL/MSVOA_W/SOIL ALS Vial : 11 Sample Multiplier: 1

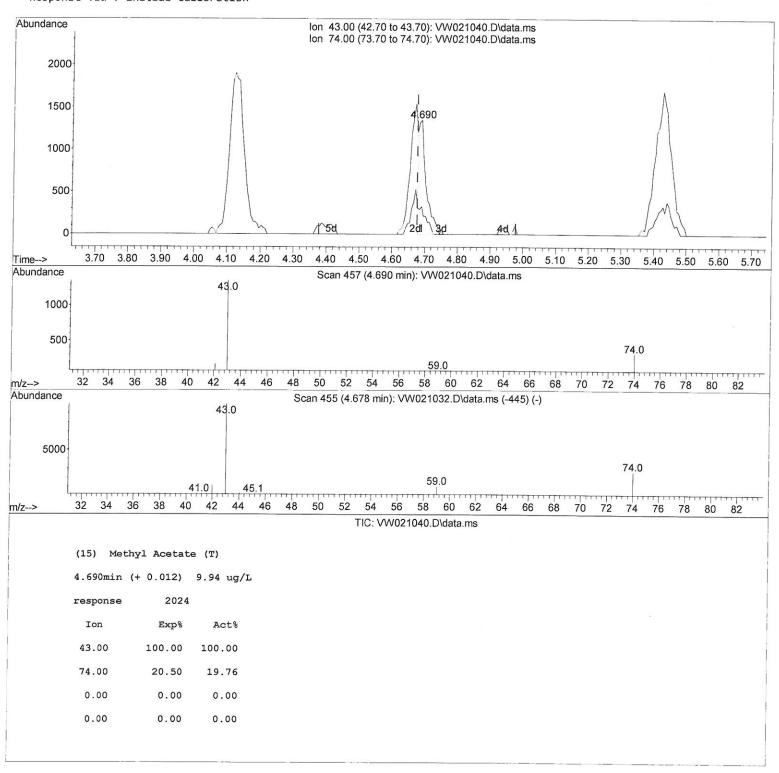
Quant Time: Dec 03 02:23:39 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M

Quant Title : SFAM01.0

QLast Update : Fri Dec 03 01:25:28 2021 Response via : Initial Calibration Instrument : MSVOA_W ClientSampleld : EX892MS

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\

Data File : VW021040.D

Acq On : 02 Dec 2021 12:23

Operator : SY/VA Sample : M4886-12MS

Misc : 5.51g/10.0mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

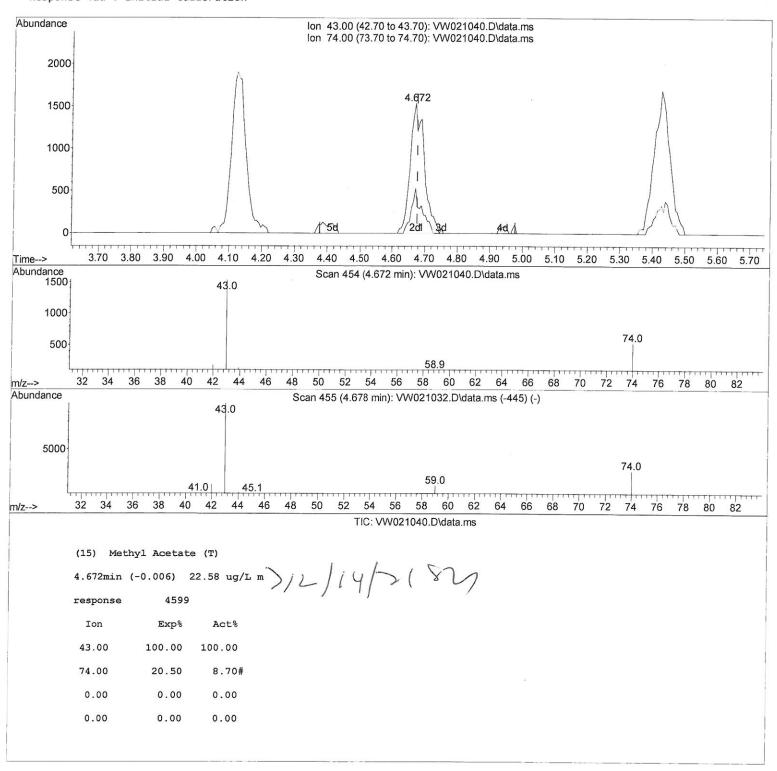
Quant Time: Dec 03 02:23:39 2021

 $\label{thm:power_power} Quant \ \ \mbox{Method} : \ \mbox{Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M}$

Quant Title : SFAM01.0

QLast Update : Fri Dec 03 01:25:28 2021 Response via : Initial Calibration Instrument : MSVOA_W ClientSampleId : EX892MS

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\

Data File : VW021040.D

Acq On : 02 Dec 2021 12:23

Operator : SY/VA

Sample : M4886-12MS Misc : 5.51g/10.0mL/MSVOA_W/SOIL ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 03 02:23:39 2021

 $\label{lem:quant_method} \mbox{Quant Method}: \mbox{Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M}$

Quant Title : SFAM01.0

QLast Update : Fri Dec 03 01:25:28 2021 Response via : Initial Calibration

Instrument : MSVOA_W ClientSampleId : EX892MS

Manual IntegrationsAPPROVED

Compound	R.T. Q	Ion Response	Conc Units Dev	(Min)	
Internal Standards					
<pre>Internal Standards 1) 1,4-Difluorobenzene</pre>	8.842 1	114 43943	25 000 110/1	# 0 00	
28) Chlorobenzene-d5		117 37626	25.000 ug/L 25.000 ug/L	# 0.00 0.00	
58) 1,4-Dichlorobenzene-d4		152 16844	25.000 ug/L	0.00	
	23,300	10044	25:000 ug/L	0.00	
System Monitoring Compounds					
4) Vinyl Chloride-d3	2.349	65 10288	20.184 ug/L	0.00	
Spiked Amount 25.000	Range 30 -	150 Recove		6	
7) Chloroethane-d5	2.885	69 10201	31.031 ug/L	0.00	
Spiked Amount 25.000	Range 30 -	150 Recove	ry = 124.120%	6	
11) 1,1-Dichloroethene-d2		63 22265	22.576 ug/L	0.00	
Spiked Amount 25.000	Range 45 -	110 Recove		6	
21) 2-Butanone-d5		46 5961	46.854 ug/L	0.00	
Spiked Amount 50.000	Range 20 -			ś	
24) Chloroform-d		84 25931	23.420 ug/L	0.00	
Spiked Amount 25.000	Range 40 -				
26) 1,2-Dichloroethane-d4 Spiked Amount 25.000		65 12781	22.742 ug/L	0.00	
32) Benzene-d6	Range 70 - : 8.275				
Spiked Amount 25.000	Range 20 - 1	84 48257	24.334 ug/L v = 97.320%	0.00	
36) 1,2-Dichloropropane-d6		135 Recover 67 13493	`y = 97.320% 23.894 ug/L		
Spiked Amount 25.000	Range 70 - 1			0.00	
41) Toluene-d8	-	98 46237	23.491 ug/L	0.00	
Spiked Amount 25.000	Range 30 - 3		y = 93.960%		
43) trans-1,3-Dichloroprop.		79 4951	18.549 ug/L	0.00	
Spiked Amount 25.000	Range 30 - 1				
47) 2-Hexanone-d5	127 August 2000 and 2	63 4573	48.827 ug/L	0.00	
Spiked Amount 50.000	Range 20 - 1				
56) 1,1,2,2-Tetrachloroeth.	12.689	34 11186	25.858 ug/L	0.00	
Spiked Amount 25.000	Range 45 - 1	120 Recover	y = 103.440%		
66) 1,2-Dichlorobenzene-d4	13.847 15	52 13671	22.772 ug/L	0.00	
Spiked Amount 25.000	Range 75 - 1	L20 Recover	y = 91.080%		
Tangat Compounds			<u>u</u>		
Target Compounds 2) Dichlorodifluoromethane	2.014 8	25 2070	Qva.		
3) Chloromethane		35 2070 50 8622	16.277 ug/L	99	
5) Vinyl chloride		52 16283	22.299 ug/L 24.348 ug/L	99	
6) Bromomethane		10283	29.172 ug/L	98 92	
8) Chloroethane		9157	32.569 ug/L	97	
9) Trichlorofluoromethane	3.257 10		45.539 ug/L	98	
10) 1,1,2-Trichloro-1,2,2			23.930 ug/L	91	
12) 1,1-Dichloroethene		6 11694	21.239 ug/L	82	
13) Acetone	4.123 4		50.445 ug/L		
14) Carbon disulfide	4.379 7	6 25463	17.100 ug/L	97 1111	21 64
15) Methyl Acetate	4.672 4	3 4599m	22.577 ug/L	12/17/	7
16) Methylene chloride	4.916 8	4 13434 /	21.920 ug/L	90	/
17) trans-1,2-Dichloroethene	5.428 9	6 12142	20.462 ug/L	81	
18) Methyl tert-butyl Ether	5.428 7	3 25721	28.877 ug/L	94	
19) 1,1-Dichloroethane	6.220 6		22.649 ug/L	95	
20) cis-1,2-Dichloroethene	7.171 9		21.481 ug/L	80	
22) 2-Butanone	7.171 4		44.457 ug/L	87	
23) Bromochloromethane	7.519 128	8 6508	22.259 ug/L #	68	

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\

Data File : VW021040.D

Acq On : 02 Dec 2021 12:23

Operator : SY/VA

Sample : M4886-12MS Misc : 5.51g/10.0mL/MSVOA_W/SOIL ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 03 02:23:39 2021

 $\label{eq:Quant_Method} \textbf{Quant Method}: \textbf{Z:}\\ \textbf{Voasrv}\\ \textbf{HPCHEM1}\\ \textbf{MSVOA_W}\\ \textbf{Method}\\ \textbf{SFAMWLM111521SMA.M}$

Quant Title : SFAM01.0

QLast Update : Fri Dec 03 01:25:28 2021 Response via : Initial Calibration

Instrument : MSVOA_W ClientSampleId : EX892MS

Manual IntegrationsAPPROVED

	Compound	R.T.	QIon	Response	Conc Units Dev(Min)
25)	Chloroform	7.677	83	24790	23.068 ug/L	100
27)	1,2-Dichloroethane	8.403	62	15518	23.847 ug/L	98
	Cyclohexane	7.958	56	18656	22.240 ug/L #	82
30)	1,1,1-Trichloroethane	7.872	97	24213	26.760 ug/L	94
31)	Carbon tetrachloride	8.067	117	20380	23.798 ug/L	97
33)	Benzene	8.323	78	51286	24.074 ug/L	100
	Trichloroethene	9.092	95	13908	23.340 ug/L	88
	Methylcyclohexane	9.335	83	21220	21.599 ug/L	86
37)	1,2-Dichloropropane	9.366	63	12031	24.266 ug/L #	96
	Bromodichloromethane	9.646	83	15947	22.412 ug/L #	94
39)	cis-1,3-Dichloropropene	10.073	75	17694	21.615 ug/L	100
40)	4-Methyl-2-pentanone	10.207	43	14572	49.379 ug/L #	93
	Toluene	10.390	91	58726	24.507 ug/L	96
44)	trans-1,3-Dichloropropene	10.604	75	14639	20.348 ug/L	99
45)	1,1,2-Trichloroethane	10.786	97	9656	25.041 ug/L	96
46)	Tetrachloroethene	10.866	164	11966	21.871 ug/L	92
	2-Hexanone	10.969	43	10322	50.496 ug/L #	98
49)	Dibromochloromethane	11.128	129	10877	20.460 ug/L	94
50)	1,2-Dibromoethane	11.238	107	8893	23.003 ug/L #	96
	Chlorobenzene	11.658	112	34863	21.875 ug/L	92
52) 1	Ethylbenzene	11.731	91	61715	23.242 ug/L	96
	m,p-Xylene	11.835	106	24896	22.965 ug/L	97
54)	o-Xylene	12.164	106	23860	22.917 ug/L	89
	Styrene	12.183	104	35346	20.445 ug/L	95
	1,1,2,2-Tetrachloroethane	12.713	83	10097	23.625 ug/L #	88
59) E	3romoform	12.347	173	5628	22.409 ug/L	98
	Isopropylbenzene	12.463	105	65036	28.348 ug/L	96
61) 1	1,2,3-Trichloropropane	12.768	75	7459	30.190 ug/L #	90
62) 1	l,3,5-Trimethylbenzene	12.938	105	50634	26.163 ug/L	92
63) 1	l,2,4-Trimethylbenzene	13.249	105	49891	26.128 ug/L	97
64) 1	l,3-Dichlorobenzene	13.493	146	23670	22.006 ug/L	95
65) 1	l,4-Dichlorobenzene	13.579	146	24185	22.290 ug/L	93
67) 1	,2-Dichlorobenzene	13.865	146	22940	24.203 ug/L	90
68) 1	,2-Dibromo-3-chloropr	14.481	75	1487	25.174 ug/L #	76
69) 1	.,3,5-Trichlorobenzene	14.627	180	15556	18.260 ug/L	96
70) 1	.,2,4-trichlorobenzene	15.127	180	10055	14.427 ug/L	94
	Taphthalene	15.365	128	21380	18.826 ug/L	100
72) 1	,2,3-Trichlorobenzene	15.554	180	8414	14.306 ug/L	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\

Data File : VW021040.D

Acq On : 02 Dec 2021 12:23

Operator : SY/VA Sample : M4886-12MS

Misc : 5.51g/10.0mL/MSVOA_W/SOIL
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 03 02:23:39 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M

Quant Title : SFAM01.0

QLast Update : Fri Dec 03 01:25:28 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

