Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW110821\

Data File : VW020806.D

Acq On : 08 Nov 2021 13:18

Operator : SY/VA Sample : VIBLK

Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

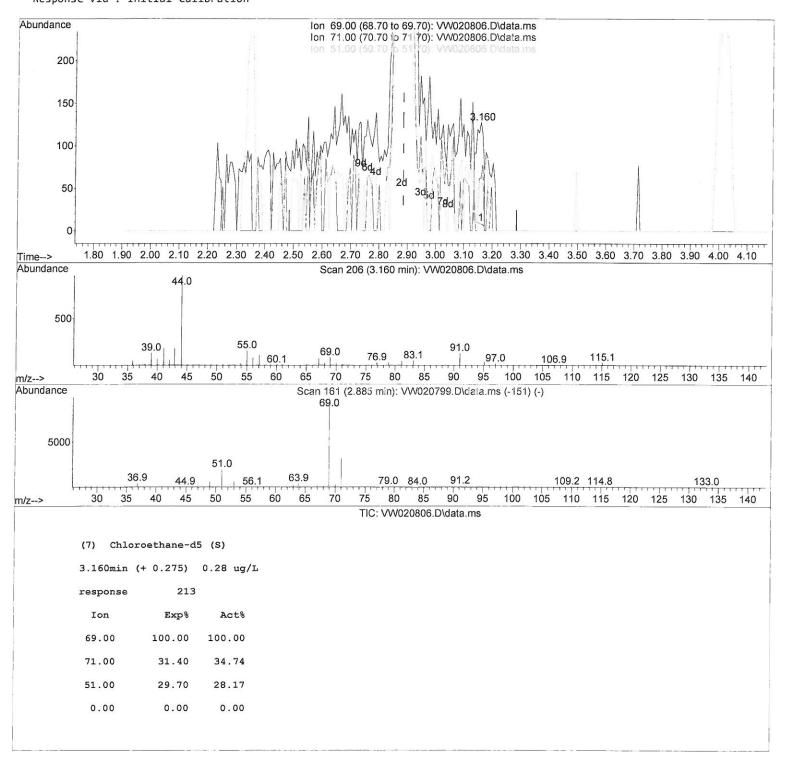
Quant Time: Nov 09 07:11:12 2021

 $\label{thm:constraint} Quant \ \mbox{Method} : \ Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM110621SMA.M$

Quant Title : SFAM01.0

QLast Update : Mon Nov 08 04:39:29 2021 Response via : Initial Calibration Instrument : MSVOA_W ClientSampleld : VIBLK479

Manual IntegrationsAPPROVED



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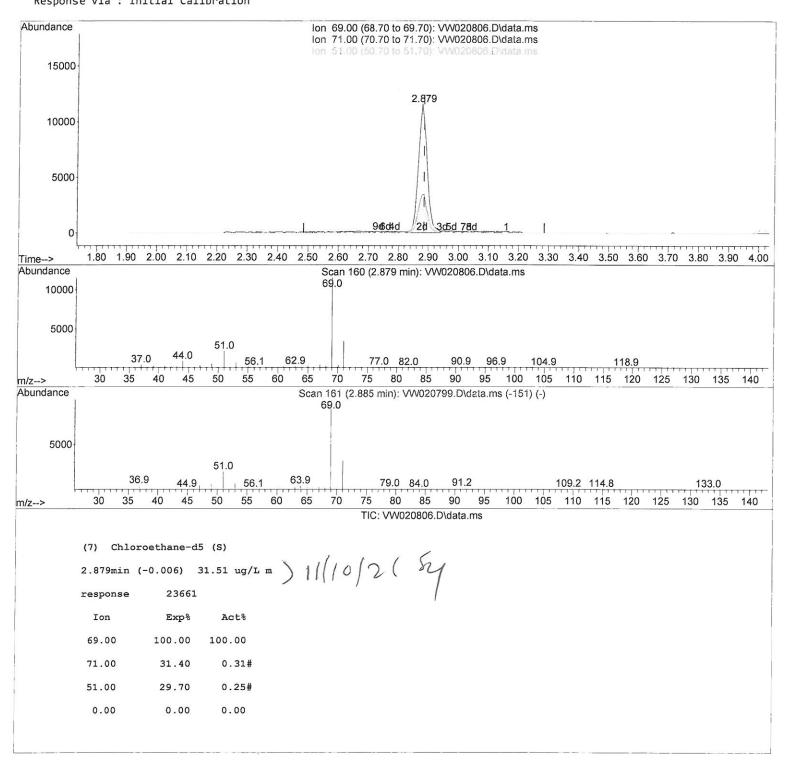
Quant Time: Nov 09 07:11:12 2021

Quant Method: Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM110621SMA.M

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 $\label{eq:Quant_Method} \textbf{Quant Method}: \textbf{Z:} \\ \textbf{Voasrv} \\ \textbf{HPCHEM1} \\ \textbf{MSVOA_W} \\ \textbf{Method} \\ \textbf{SFAMWLM110621SMA.M} \\ \\ \textbf{MSVOA_W} \\ \textbf{Method} \\ \textbf{MSVOA_W} \\ \textbf{Method} \\ \textbf{MSVOA_W} \\ \textbf{MSVOA_W} \\ \textbf{Method} \\ \textbf{MSVOA_W} \\ \textbf{MSVOA_W} \\ \textbf{Method} \\ \textbf{MSVOAW} \\ \textbf{Method} \\ \textbf{$

Quant Title : SFAM01.0

QLast Update: Mon Nov 08 04:39:29 2021 Response via : Initial Calibration

Instrument : MSVOA_W ClientSampleId : VIBLK479

Manual IntegrationsAPPROVED

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nesponse via i initial carro	acton		
Compound	R.T.	QIon	Response Conc Units Dev(Min)
Internal Standards			
1) 1,4-Difluorobenzene	8.848	114	140252 25.000 ug/L # 0.0
28) Chlorobenzene-d5	11.628	117	123977 25.000 ug/L 0.00
58) 1,4-Dichlorobenzene-d4	13.560	152	68537 25.000 ug/L 0.00
20, 2, 220.20.20.20.20.2	231300		25.000 ug/ 2
System Monitoring Compounds			
4) Vinyl Chloride-d3	2.349	65	38866 29.474 ug/L 0.00
Spiked Amount 25.000		- 150	Recovery = 117.880%
7) Chloroethane-d5	2.879	69	23661m 31.512 ug/L 0.00
Spiked Amount 25.000		- 150	Recovery = 126.040%
11) 1,1-Dichloroethene-d2	4.013	63	49442 17.863 ug/L 0.00
Spiked Amount 25.000	Range 45		Recovery = 71.440%
21) 2-Butanone-d5	7.080	46	
Spiked Amount 50.000		- 135	20346 46.459 ug/L 0.00 Recovery = 92.920%
24) Chloroform-d	7.647	84	•
Spiked Amount 25.000	Range 40		
26) 1,2-Dichloroethane-d4	0	65	Recovery = 98.440%
Spiked Amount 25.000	8.305		42311 25.504 ug/L 0.00
	0	- 130	Recovery = 102.000%
32) Benzene-d6	8.275	84	154588 27.598 ug/L 0.00
Spiked Amount 25.000		- 135	Recovery = 110.400%
36) 1,2-Dichloropropane-d6	9.274	67	45980 26.513 ug/L 0.00
Spiked Amount 25.000	0	- 120	Recovery = 106.040%
41) Toluene-d8	10.323	98	153895 28.240 ug/L 0.00
Spiked Amount 25.000	Range 30		Recovery = 112.960%
43) trans-1,3-Dichloroprop.		79	21934 25.500 ug/L 0.00
Spiked Amount 25.000	-	- 135	Recovery = 102.000%
47) 2-Hexanone-d5	10.920	63	17053 48.598 ug/L 0.00
Spiked Amount 25.000	O .	- 135	Recovery = 194.400%#
56) 1,1,2,2-Tetrachloroeth.		84	35742 24.868 ug/L 0.00
Spiked Amount 25.000	Range 45	- 120	Recovery = 99.480%
66) 1,2-Dichlorobenzene-d4	13.853	152	59805 27.766 ug/L 0.00
Spiked Amount 25.000	Range 75	- 120	Recovery = 111.080%
Target Compounds			Qvalue
20) cis-1,2-Dichloroethene	7.171	96	2639 1.313 ug/L # 62
34) Trichloroethene	9.091	95	2752 1.392 ug/L 88
42) Toluene	10.384	91	19360 2.425 ug/L 100

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

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Manual IntegrationsAPPROVED

