Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\

Data File : VY006948.D

Acq On : 04 Dec 2021 17:37

Operator : SY/MD Sample : M4884-18MS

Misc : 5.68g/10.0mL/MSVOA_Y/SOIL ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 05 06:08:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYLM120321SMA.M

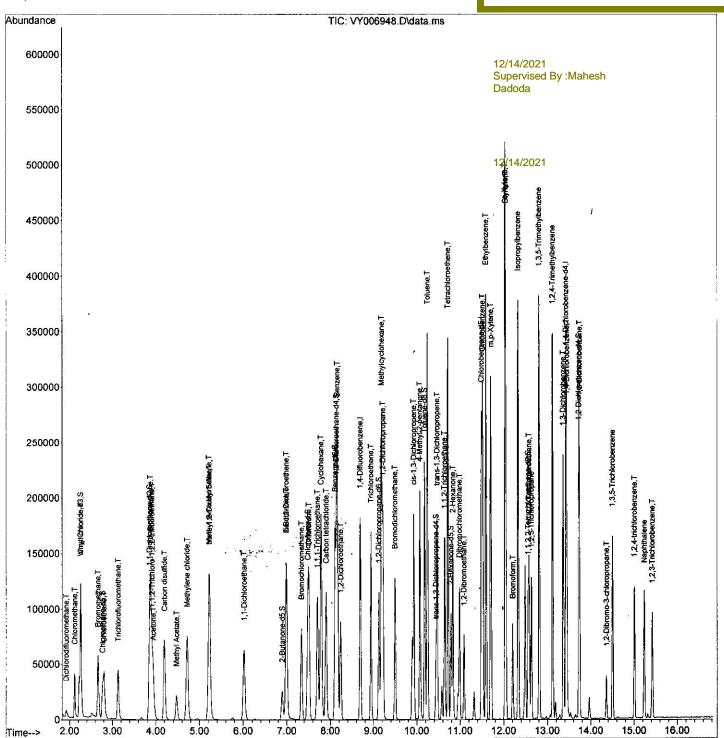
Quant Title : VOC Analysis

QLast Update : Sat Dec 04 02:26:29 2021 Response via : Initial Calibration Instrument : MSVOA_Y ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :John

Carlone



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\

Data File : VY006948.D

Acq On : 04 Dec 2021 17:37

Operator : SY/MD : M4884-18MS Sample

Misc : 5.68g/10.0mL/MSVOA_Y/SOIL ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 05 06:08:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYLM120321SMA.M

Quant Title : VOC Analysis

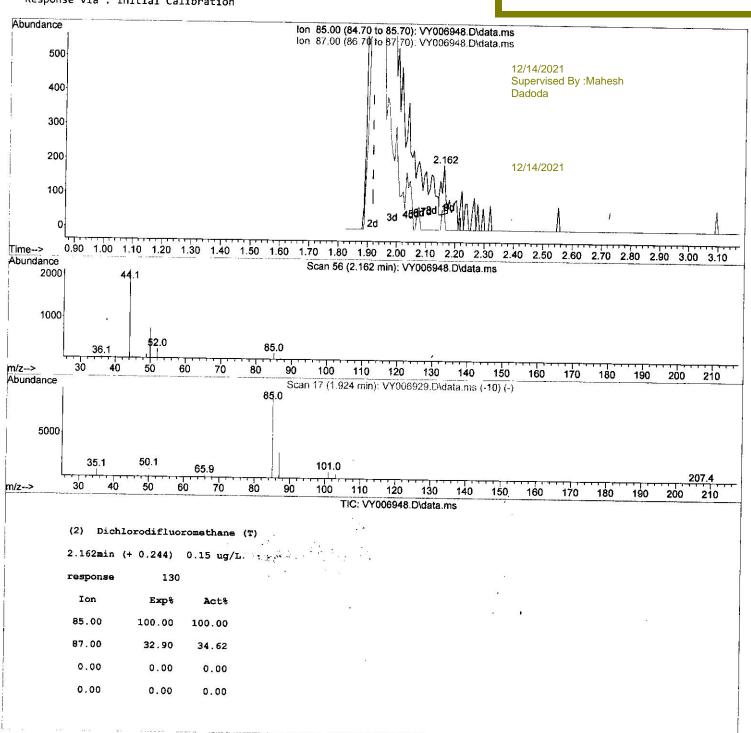
QLast Update : Sat Dec 04 02:26:29 2021 Response via : Initial Calibration

Instrument: MSVOA_Y ClientSampleId:

Manual IntegrationsAPPROVED

Reviewed By :John

Carlone



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\

Data File: VY006948.D

Acq On : 04 Dec 2021 17:37

Operator : SY/MD Sample : M4884-18MS

Misc : 5.68g/10.0mL/MSVOA_Y/SOIL ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 05 06:08:51 2021

 $\label{thm:policy} Quant \ \mbox{Method}: \ Z:\voasrv\HPCHEM1\MSVOA_Y\mbox{method}s\SFAMYLM120321SMA.M$

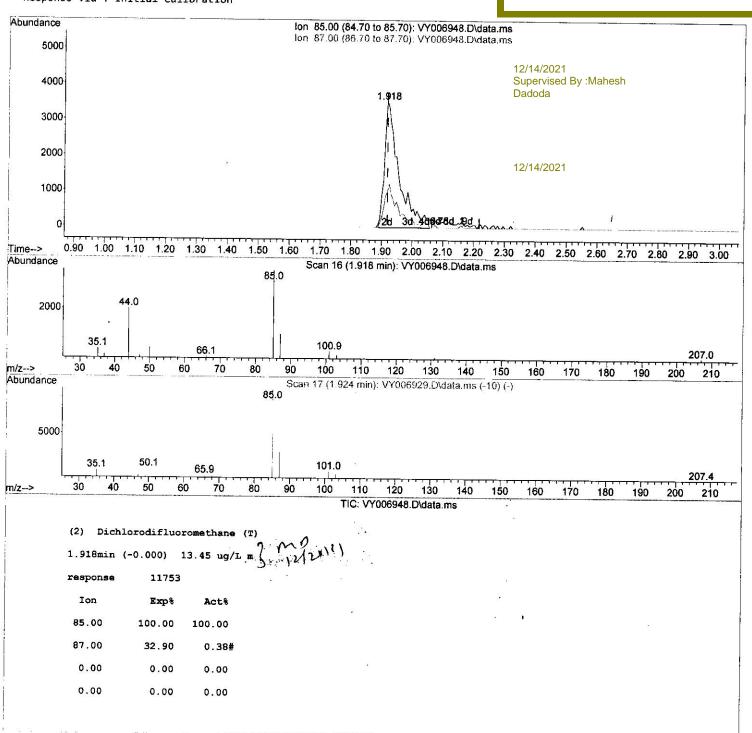
Quant Title : VOC Analysis

QLast Update : Sat Dec 04 02:26:29 2021 Response via : Initial Calibration Instrument : MSVOA_Y ClientSampleId :

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Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\

Data File : VY006948.D Acq On : 04 Dec 2021 17:37

Operator : SY/MD Sample : M4884-18MS

Misc : 5.68g/10.0mL/MSVOA_Y/SOIL ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 05 06:08:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYLM120321SMA.M

Quant Title : VOC Analysis QLast Update : Sat Dec 04 02:26:29 2021 Response via : Initial Calibration

Instrument: MSVOA_Y ClientSampleId: EW9K3MS

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Reviewed By :John

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Compound	R.T. QIon	Response Conc Units Dev(Min)	
Internal Standards				
1) 1,4-Difluorobenzene	8.691 114	138447 25.000 ug/L	0.00	40/44/0004
28) Chlorobenzene-d5	11.489 117	126149 25.000 ug/L	0.00	12/14/2021 Supervised By :Mahesh
58) 1,4-Dichlorobenzene-d4	13.428 152	59852 25.000 ug/L	0.00	Dadoda
		•		
System Monitoring Compounds	3.053 48		V.10 D.0790AD	
4) Vinyl Chloride-d3 Spiked Amount 25.000	2.253 65	55710 31.244 ug/L	0.00	
Spiked Amount 25.000 (7) Chloroethane-d5	Range 30 - 150 2.771 69	Recovery = 124.960%		
	Range 30 - 150	39017 27.894 ug/L Recovery = 111.560%	0.00	12/14/2021
11) 1,1-Dichloroethene-d2	3.857 63	86043 23.211 ug/L	0.00	
	Range 45 - 110	Recovery = 92.840%		
21) 2-Butanone-d5	6.893 46	45999 66.792 ug/L	0.00	r .
Spiked Amount 50.000	Range 20 - 135	Recovery = 133.580%		
24) Chloroform-d	7.484 84	81056 23.493 ug/L	0.00	
	Range 40 - 150	Recovery = 93.960%		
26) 1,2-Dichloroethane-d4	8.142 65	52353 25.279 ug/L	0.00	
	Range 70 - 130	Recovery = 101.120%		
32) Benzene-d6	8.112 84	159466 23.797 ug/L	0.00	
Spiked Amount 25.000 F 36) 1,2-Dichloropropane-d6	Range 20 - 135	Recovery = 95.200%		
e :	9.124 67	52443 24.909 ug/L	0.00	
41) Toluene-d8	Range 70 - 120 10.179 98	Recovery = 99.640% 145560 23.193 ug/L	0.00	
	Range 30 - 130	145560 23.193 ug/L Recovery = 92.760%	0.00	
43) trans-1,3-Dichloroprop		25880 25.413 ug/L	0.00	
	Range 30 - 135	Recovery = 101.640%	0.00	
47) 2-Hexanone-d5	10.788 63	30984 63.585 ug/L	0.00	
Spiked Amount 50.000 F	Range 20 - 135	Recovery = 127.160%		
56) 1,1,2,2-Tetrachloroeth	12.562 84	47067 26.833 ug/L	0.00	
Spiked Amount 25.000 F	Range 45 - 120	Recovery = 107.320%		
66) 1,2-Dichlorobenzene-d4	13.721 152	40528 20.764 ug/L	0.00	
Spiked Amount 25.000 F	Range 75 - 120	Recovery = 83.040%		
Target Compounds		1 Qva	lue	.0. 10
Dichlorodifluoromethane	1.918 85	11753m 13.446 ug/L	LUC	12/20/21
Chloromethane	2.119 50	37763 22.508 ug/L	99	1210
Vinyl chloride	2.259 62	79924 :32.340 ug/L	97	
6) Bromomethane	2.662 94	44005 29.458 ug/L	100	
8) Chloroethane	2.802 . 64	43835 29.893 ug/L	100	
9) Trichlorofluoromethane	3.131 101	46341 25.820 ug/L	98	
10) 1,1,2-Trichloro-1,2,212) 1,1-Dichloroethene		42980 24.702 ug/L	92	•
13) Acetone	3.875 96 3.942 43	43292 23.107 ug/L	89	•
14) Carbon disulfide	3.942 43 4.198 76	35882 56.850 ug/L 143755 23.275 ug/L	97	j.
15) Methyl Acetate	4.472 43	43687 31.749 ug/L	100	
16) Methylene chloride	4.716 84	51952 20.925 ug/L	97 83	
17) trans-1,2-Dichloroethene	5.222 96	45377 22.715 ug/L	95	
18) Methyl tert-butyl Ether	5.222 73	95627 26.536 ug/L	98	
19) 1,1-Dichloroethane	6.021 63	93980 25.013 ug/L	99	
20) cis-1,2-Dichloroethene	6.984 96	54246 25.324 ug/L	96	
22) 2-Butanone	6.984 43	67534 76.489 ug/L	91	
23) Bromochloromethane	7.338 128	26051 26.651 ug/L #	84	

7.508 83

100231 26.480 ug/L

98

25) Chloroform

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY120421\

Data File : VY006948.D

Acq On : 04 Dec 2021 17:37

Operator : SY/MD

Sample : M4884-18MS

Misc : 5.68g/10.0mL/MSVOA_Y/SOIL

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 05 06:08:51 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\SFAMYLM120321SMA.M

Quant Title : VOC Analysis QLast Update : Sat Dec 04 02:26:29 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
27) 1,2-Dichloroethane	8.240	62	78035	28.449 ug/L	98
29) Cyclohexane	7.789	56	106927	30.412 ug/L	92
30) 1,1,1-Trichloroethane	7.703	97	88185	27.480 ug/L	98
31) Carbon tetrachloride	7.905	117	81071	26.477 ug/L	100
33) Benzene	8.161	78	225364	26.572 ug/L	100
34) Trichloroethene	8.941	95	54674	26.172 ug/L	90
35) Methylcyclohexane	9.185	83	104471	27.909 ug/L	95
37) 1,2-Dichloropropane	9.215	63	63932	29.407 ug/L	98
38) Bromodichloromethane	9.496	83	78999	28.046 ug/L	99
39) cis-1,3-Dichloropropene	9.929	75	94873	27.799 ug/L	100
40) 4-Methyl-2-pentanone	10.069	43	147867	77.979 ug/L	96
42) Toluene	10.246	91	235207	25.995 ug/L	100
44) trans-1,3-Dichloropropene	10.465	75	92316	28.946 ug/L	96
45) 1,1,2-Trichloroethane	10.642	97	46588	29.085 ug/L	92
46) Tetrachloroethene	10.721	164	69391	39.648 ug/L	97
48) 2-Hexanone	10.831	43	103743	82.903 ug/L	97
49) Dibromochloromethane	10.983	129	54338	27.875 ug/L	92
50) 1,2-Dibromoethane	11.087	107	45507	30.019 ug/L	98
51) Chlorobenzene	11.514	112	135872	24.622 ug/L	95
52) Ethylbenzene	11.593	91	255196	26.239 ug/L	99
53) m,p-Xylene	11.703	106	90168	24.508 ug/L	96
54) o-Xylene	12.032	106	84870	24.522 ug/L	98
55) Styrene	12.044	104	147284	24.323 ug/L	94
57) 1,1,2,2-Tetrachloroethane	12.581	83	61705	31.927 ug/L	92
59) Bromoform	12.209	173	34326	28.121 ug/L	98
60) Isopropylbenzene	12.331	105	228701	26.112 ug/L	99
61) 1,2,3-Trichloropropane	12.636	75	49393	33.951 ug/L	97
62) 1,3,5-Trimethylbenzene	12.812	105	190867	26.504 ug/L	100
63) 1,2,4-Trimethylbenzene	13.123	105	183703	26.080 ug/L	99
64) 1,3-Dichlorobenzene	13.361	146	93796	23.795 ug/L	95
65) 1,4-Dichlorobenzene	13.446	146	94874	23.620 ug/L	98
67) 1,2-Dichlorobenzene	13.739	146	83399	23.628 ug/L	96
68) 1,2-Dibromo-3-chloropr	14.361	75	9357	29.870 ug/L	91
69) 1,3,5-Trichlorobenzene	14.501	180	44807	16.703 ug/L	99
70) 1,2,4-trichlorobenzene	15.007	180	32995	15.250 ug/L	99
71) Naphthalene	15.239	128	89698	20.690 ug/L	99
72) 1,2,3-Trichlorobenzene	15.422	180	29035	. 15.882 ug/L	98
***************************************			;		

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

Instrument: MSVOA_Y
ClientSampleId: EW9K3MS

Manual IntegrationsAPPROVED

Reviewed By :John Carlone

12/14/2021

Supervised By: Mahesh Dadoda

12/14/2021