Quantitation Report (QT Reviewed)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025133.D

Acq On : 11 Nov 2021 16:06

Operator : JC/MD Sample : VSTDICV050

Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 12 04:59:20 2021

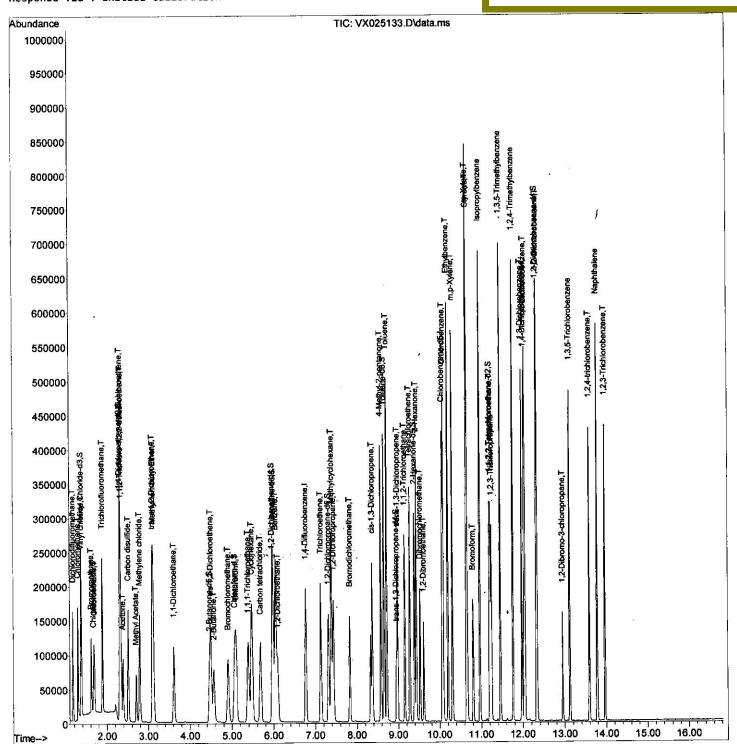
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Nov 11 16:29:23 2021 Response via : Initial Calibration Instrument : MSVOA_X ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 11/12/2021 Supervised By :Mahesh Dadoda 11/12/2021



SFAMXLM111121WMA.M Fri Nov 12 05:01:11 2021

Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025133.D

: 11 Nov 2021 16:06 Acq On

: JC/MD Operator : VSTDICV050 Sample

: 5.0mL/MSVOA_X/WATER Misc ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 12 04:59:20 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

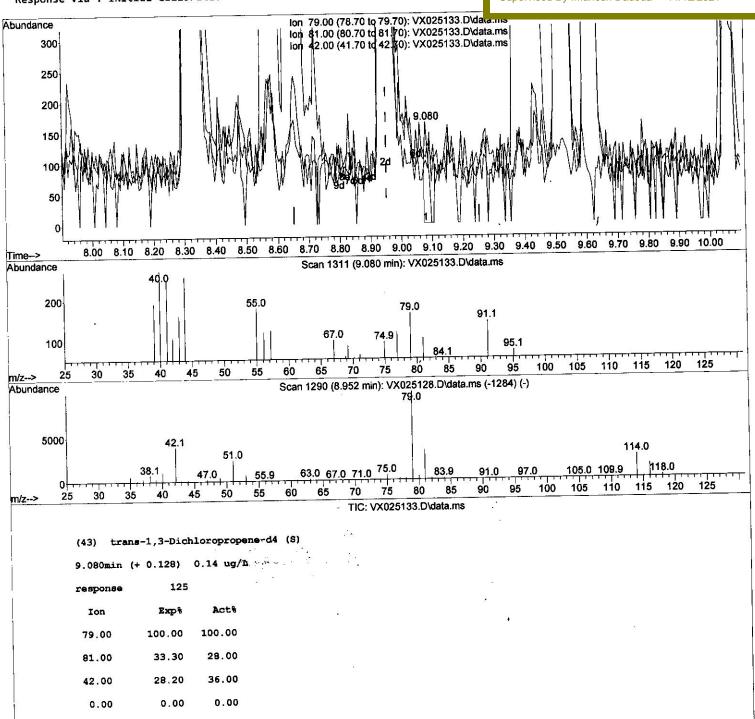
Quant Title : VOC Analysis

QLast Update : Thu Nov 11 16:29:23 2021 Response via : Initial Calibration

Instrument: MSVOA_X ClientSampleId:

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Reviewed By :John Carlone 11/12/2021 Supervised By: Mahesh Dadoda 11/12/2021



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025133.D

: 11 Nov 2021 16:06 Acq On

: JC/MD Operator : VSTDICV050 Sample

: 5.0mL/MSVOA_X/WATER Misc Sample Multiplier: 1 : 9 ALS Vial

Quant Time: Nov 12 04:59:20 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

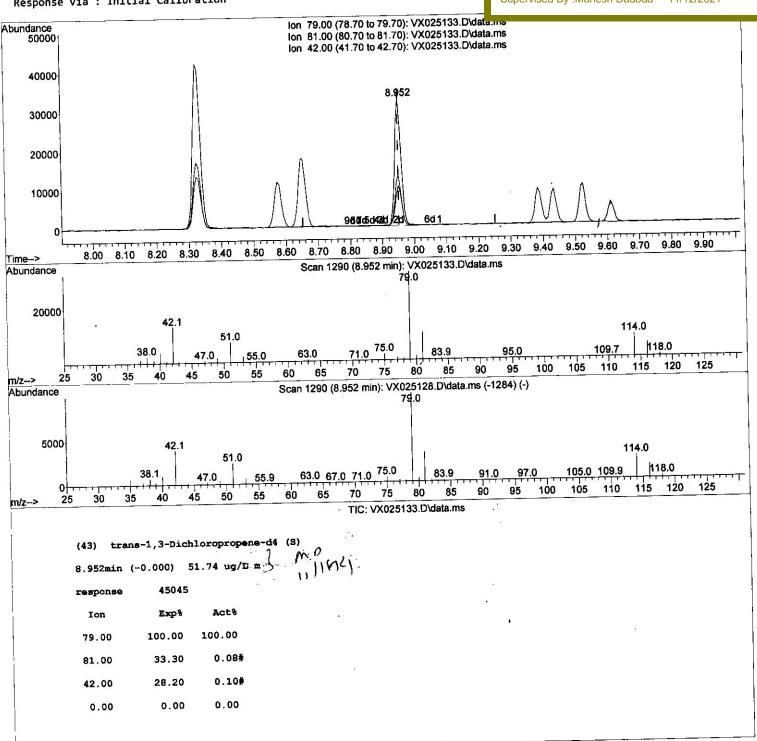
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QLast Update : Thu Nov 11 16:29:23 2021 Response via : Initial Calibration

Instrument: MSVOA_X ClientSampleId:

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Operator : JC/MD : VSTDICV050 Sample

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Quant Time: Nov 12 04:59:20 2021

Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M Quant Method

Quant Ti

QLast Up Response

lethod : T:/Aogsta/ubcutut/	IZAON_V	Auc cu	Ja (DI MINELIN			
itle : VOC Analysis						
pdate : Thu Nov 11 16:29:23	2021					
e via : Initial Calibration						
Compound	R.T.	QIon	Response	Conc Units D	Dev(Min)	
nal Standards	6.763	114	211029	50.000 ug/l	L 0.06	ð

Internal Standards					_		200	
1) 1,4-Difluorobe	enzene	6.	763	114	211029	50.000	1.00	0.00
28) Chlorobenzene	-d5	10.	055	117	192427	50.000	100000000000000000000000000000000000000	0.00
58) 1.4-Dichlorobe		12.	024	152	101323	50.000	ug/L	0.00
System Monitoring					15 A (Contract Contract)			
4) Vinyl Chloride	e-d3		368	65	72258	50.683		0.00
Spiked Amount	50.000	Range	60		Recover		101.360%	
7) Chloroethane-	d5	1.	666	69	46853	57.663		0.00
Spiked Amount	50.000	Range	70	- 130	Recover		115.320%	2.2
11) 1,1-Dichloroe	thene-d2	2.	307	63	122533	49.972		0.00
Spiked Amount	50.000	Range	60	- 125	Recover	٠y =	99.940%	
21) 2-Butanone-d5			459	46	113710	105.544	ug/L	0.00
Spiked Amount	100.000	Range	40	- 130	Recover	^y =	105.540%	
24) Chloroform-d			.062	84	128524	51.244	ug/L	0.00
Spiked Amount	50.000	Range	70	- 125	Recover	ry =	102.480%	
26) 1,2-Dichloroe		0.000 0.000 0.000	.958	65	77285	50.837	ug/L	0.00
Spiked Amount	50.000	Range		- 125	Recove	ry ≖	101.680%	
32) Benzene-d6	301005		.977	84	266768	50.794	ug/L	0.00
Spiked Amount	50.000	Range		- 125	Recove	rv =	101.580%	
36) 1,2-Dichlorop			.312	67	81600	50,885		0.00
	50.000	Range			Recove	rv =	101.760%	
Spiked Amount	30.000		.653		255001	50.843		0.00
41) Toluene-d8	FO 000	Range			Recove			
Spiked Amount	50.000		.952		7	51.749		0.00
43) trans-1,3-Dic					Recové	3		PER ENGERORIA
Spiked Amount	50.000	Range	.385		90090	104.619		0.00
47) 2-Hexanone-d5		0 -2 7		_	Recove			1000 BENEV
Spiked Amount	100.000			- 130 84	116986		g ug/L	0.00
56) 1,1,2,2-Tetra	ichloroeth.		. 195				101.300%	
Spiked Amount	50.000			- 120			101.300% 3 ug/L	0.00
66) 1,2-Dichlorob			.323		101067		100.640%	
Spiked Amount	50.000	Range	80	- 120	Recove	ry =	100.040	•

Qvalue Target Compounds 98 -83399 50.566 ug/L 85 1.167 2) Dichlorodifluoromethane 89 49.284 ug/L 1.288 50 87999 Chloromethane 91993 49.961 ug/L 100 62 5) Vinyl chloride 1.374 96 1.612 94 1.685 64 39345 50830 55.685 ug/L 1.612 6) Bromomethane 97 54.974 ug/L 8) Chloroethane 100 50.039 ug/L 1.886 101 134488 9) Trichlorofluoromethane 50.314 ug/L 96 70073 2.331 10) 1,1,2-Trichloro-1,2,2-... 49.511 ug/L 87 66617 96 12) 1,1-Dichloroethene 2.319 98 99965 98.915 ug/L 2.386 43 13) Acetone 48.204 ug/L 99 199839 2.514 76 14) Carbon disulfide 81094 49.286 ug/L # 82 2.703 43 15) Methyl Acetate 81 2.788 72382 48.630 ug/L # 16) Methylene chloride 90 71284 48.783 ug/L 17) trans-1,2-Dichloroethene 3.093 96 90 226950 49.689 ug/L # 3.111 73 18) Methyl tert-butyl Ether 95 49.609 ug/L 19) 1,1-Dichloroethane 3.611 63 122622 98 49.339 ug/L # 96 79770 20) cis-1,2-Dichloroethene 4.489 84 100.755 ug/L 4.562 43 135543 22) 2-Butanone 41215 80 49.514 ug/L # 23) Bromochloromethane 4.904 128

Instrument: MSVOA_X ClientSampleId: VICV636

Manual IntegrationsAPPROVED

Reviewed By :John Carlone 11/12/2021 Supervised By: Mahesh Dadoda 11/12/2021

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111121\

Data File : VX025133.D

Acq On : 11 Nov 2021 16:06
Operator : JC/MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_X/WATER

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 12 04:59:20 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM111121WMA.M

Quant Title : VOC Analysis

QLast Update : Thu Nov 11 16:29:23 2021 Response via : Initial Calibration

Compound	R.T. (lon	Response	Conc Units Dev(M:	in)
25) Chloroform	5.093	83	125462	49.522 ug/L	99
27) 1,2-Dichloroethane	6.092	62	91966	49.509 ug/L #	87
29) Cyclohexane	5.471	56	122203	49.708 ug/L	86
30) 1,1,1-Trichloroethane	5.385	97	118522	50.123 ug/L #	94
31) Carbon tetrachloride	5.678	117	104807	50.196 ug/L	100
33) Benzene	6.044	78	292991	49.559 ug/L	100
34) Trichloroethene	7.129	95	76308	49.393 ug/L	83
35) Methylcyclohexane	7.385	83	130210	49.842 ug/L	92
37) 1,2-Dichloropropane	7.434	63	73018	49.303 ug/L	100
38) Bromodichloromethane	7.824	83	98275	49.035 ug/L	97
39) cis-1,3-Dichloropropene	8.366	75	120481	49.544 ug/L	99
40) 4-Methyl-2-pentanone	8.574	43	235892	100.221 ug/L #	84
42) Toluene	8.720	91	318130	49.542 ug/L	97
44) trans-1,3-Dichloropropene	8.976	75	117566	49.676 ug/L	98
45) 1,1,2-Trichloroethane	9.153	97	72636	48.570 ug/L	99
46) Tetrachloroethene	9.275	164	64800	50.273 ug/L	89
48) 2-Hexanone	9.433	43	192519	99.787 ug/L #	85
49) Dibromochloromethane	9.525	129	83882	49.136 ug/L	100
50) 1,2-Dibromoethane	9.610	107	79206	49.489 ug/L #	99
51) Chlorobenzene	10.080	112	206196	49.975 ug/L	97
52) Ethylbenzene	10.195	91	344548	49.917 ug/L	93
53) m,p-Xylene	10.305	106	140009	50.170 ug/L	79
54) o-Xylene	10.647	106	135777	49.305 ug/L	82
55) Styrene	10.659	104	231349	49.474 ug/L	82
57) 1,1,2,2-Tetrachloroethane	11.213	83	114978	48.425 ug/L	98
59) Bromoform	10.799	173	65162	47.587 ug/L #	95
60) Isopropylbenzene	10.964	105	353993	48.859 ug/L	95
61) 1,2,3-Trichloropropane	11.244	75	91032	47.713 ug/L	95
62) 1,3,5-Trimethylbenzene	11.451	105	302406	49.195 ug/L	89
63) 1,2,4-Trimethylbenzene	11.756	105	302190	48.999 ug/L #	87
64) 1,3-Dichlorobenzene	11.969	146	161437	48.880 ug/L	97
65) 1,4-Dichlorobenzene	12.043	146	161022	48.798 ug/L	95
67) 1,2-Dichlorobenzene	12.335	146	159680	48.678 ug/L	94
68) 1,2-Dibromo-3-chloropr	12.945	75	27049	49.019 ug/L #	61
69) 1,3,5-Trichlorobenzene	13.116	180	119843	50.156 ug/L	97
70) 1,2,4-trichlorobenzene	13.591		105344	50.543 ug/L	97
71) Naphthalene	13.780		368477		99
72) 1,2,3-Trichlorobenzene	13.963	180	105373	50.997 ug/L	94
		11111			

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

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ClientSampleId: VICV636

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