Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\

Data File : VV023529.D

Acq On : 16 Nov 2021 13:12

Operator : SY/MD Sample : M4643-03MS

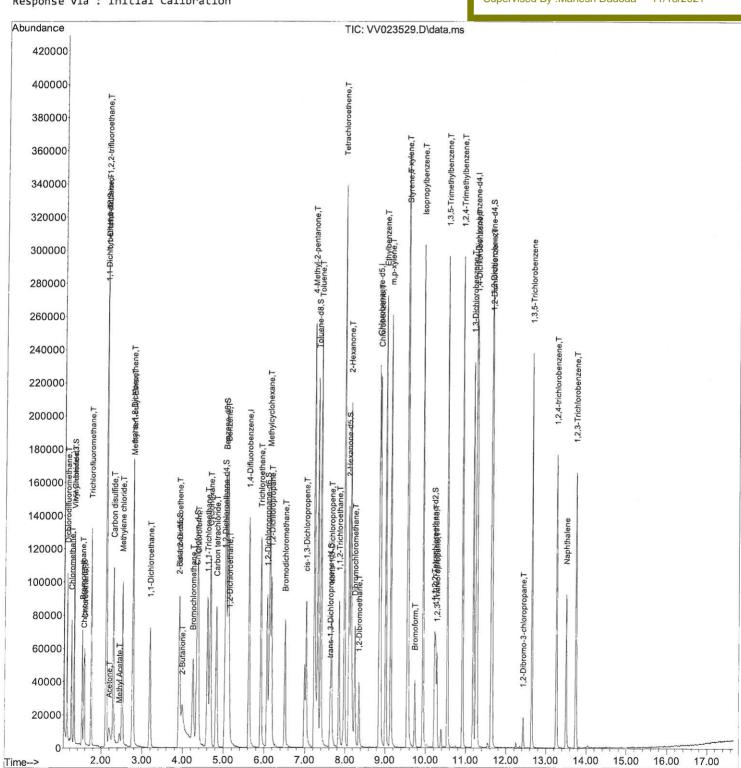
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 17 00:51:28 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 00:48:57 2021 Response via : Initial Calibration Instrument:
MSVOA_V
ClientSampleId:
GB8.J9MS

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV111621\

Data File : VV023529.D

Acq On : 16 Nov 2021 13:12

Operator : SY/MD Sample : M4643-03MS

Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

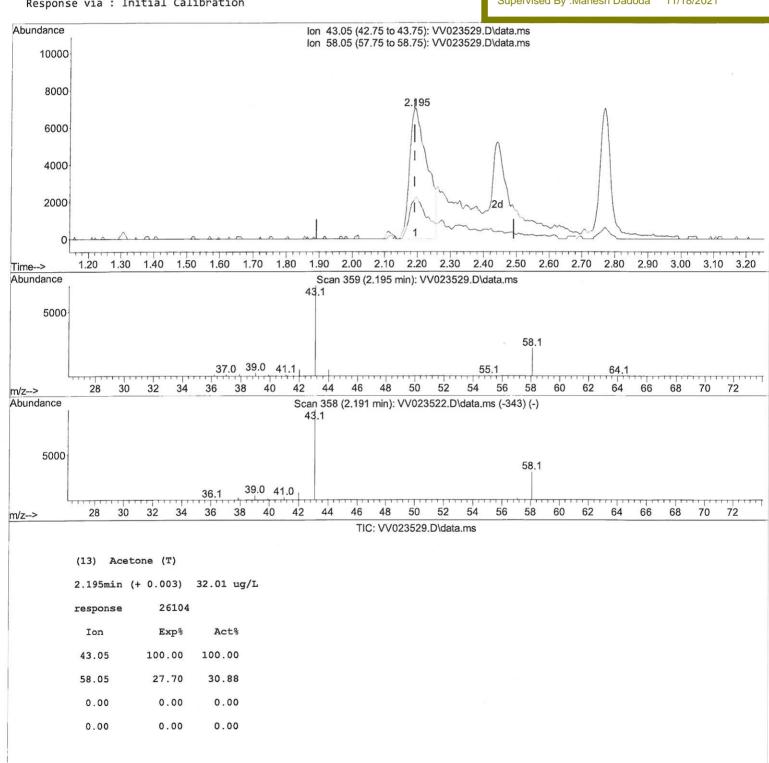
Quant Time: Nov 17 00:51:28 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0

QLast Update : Wed Nov 17 00:48:57 2021 Response via : Initial Calibration Instrument: MSVOA_V ClientSampleId: GB8J9MS

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\

Data File: VV023529.D

Acq On : 16 Nov 2021 13:12

Operator : SY/MD Sample : M4643-03MS

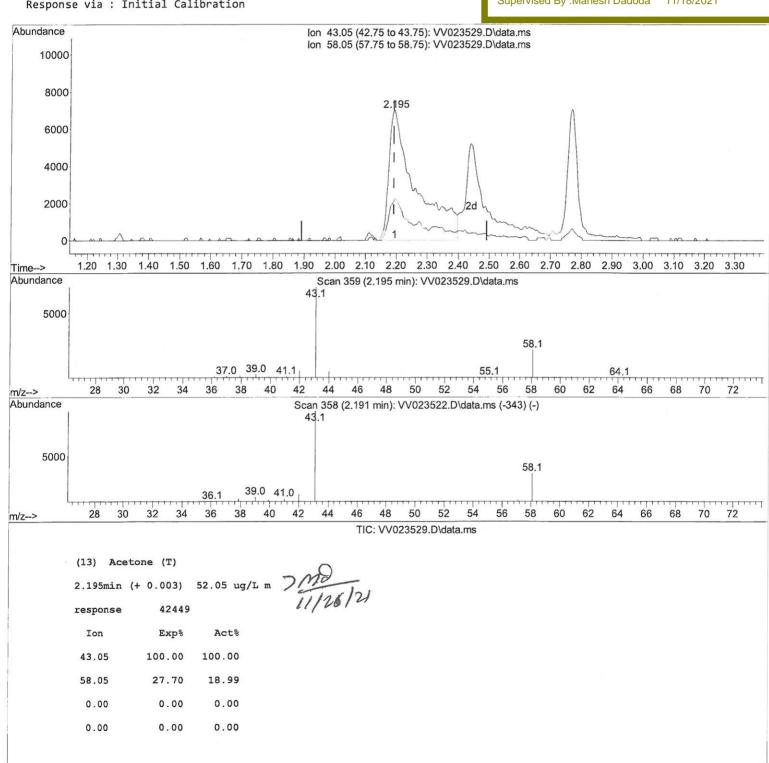
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 17 00:51:28 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 00:48:57 2021 Response via : Initial Calibration Instrument : MSVOA_V ClientSampleId :

Manual IntegrationsAPPROVED



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\

Data File: VV023529.D

Acq On : 16 Nov 2021 13:12 Operator : SY/MD

Sample : M4643-03MS Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 17 00:51:28 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update: Wed Nov 17 00:48:57 2021 Response via: Initial Calibration

Instrument: MSVOA_V ClientSampleId: GB8J9MS

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Un	its Dev(Min)	
Internal Standards							
1) 1,4-Difluorobenzene	5.619	11/	123674	5.000	ua/I	0.00	
28) Chlorobenzene-d5	8.854		122669	5.000		0.00	
58) 1,4-Dichlorobenzene-d4	11.249		67661	5.000	_	0.00	
56) 134-Dichiol Obenzene u+	11.243	132	07001	3.000	ug/ L	0.00	
System Monitoring Compounds							
4) Vinyl Chloride-d3	1.307	65	34583	4.464	ug/L	0.00	
Spiked Amount 5.000	Range 40	- 130	Recover		89.200%		
7) Chloroethane-d5	1.568	69	28434	4.503	ug/L	0.00	
Spiked Amount 5.000	Range 65		Recover	-	90.000%		
<pre>11) 1,1-Dichloroethene-d2</pre>	2.111		69046	4.760		0.00	
Spiked Amount 5.000	Range 60		Recover	•	95.200%		
20) 2-Butanone-d5	3.908	46	54248	40.642		0.00	
Spiked Amount 50.000	Range 40		Recover	A. The second second second	81.280%		
24) Chloroform-d	4.349	84	74702	4.524	10.00	0.00	
Spiked Amount 5.000	Range 70		Recover		90.400%	0 00	
26) 1,2-Dichloroethane-d4	5.034	65	36245	4.882		0.00	
Spiked Amount 5.000	Range 70		Recover		97.600%	0 00	
<pre>32) Benzene-d6 Spiked Amount 5.000</pre>	5.053	84 125		4.720	94.400%	0.00	
Spiked Amount 5.000 36) 1,2-Dichloropropane-d6	Range 70 6.069	67	Recover 42946	17:0		0.00	
Spiked Amount 5.000	Range 60		Recover	4.635	92.800%	0.00	
41) Toluene-d8	7.317	98	144027	4.883		0.00	
Spiked Amount 5.000	Range 70		Recover		97.600%	0.00	
43) trans-1,3-Dichloroprop.		79	18031	5.132		0.00	
Spiked Amount 5.000	Range 55		Recover		.02.600%		
46) 2-Hexanone-d5	8.091	63	63059	48.784		0.00	
Spiked Amount 50.000	Range 45		Recover		97.560%		
56) 1,1,2,2-Tetrachloroeth.			31999	4.803	ug/L	0.00	
Spiked Amount 5.000	Range 65		Recover		96.000%		
66) 1,2-Dichlorobenzene-d4	11.625	152	53133	4.716	ug/L	0.00	
Spiked Amount 5.000	Range 80	- 120	Recover	·y =	94.400%		
T					0		
Target Compounds	1 120	0E	E2E72	4 250	Qva]		
2) Dichlorodifluoromethane3) Chloromethane	1.130	85 50	52572 47077	4.359 4.591		100 99	
5) Vinyl chloride	1.240	62	48333	4.720		100	
6) Bromomethane	1.523	94	27371	4.182		96	
8) Chloroethane	1.584	64	29381	4.972		97	
9) Trichlorofluoromethane	1.754	101	75723	4.921		98	
10) 1,1,2-Trichloro-1,2,2		101	39837	5.143	-	98	
12) 1,1-Dichloroethene	2.121	96	35529	4.817		89	.0
13) Acetone	2.195	43	42449m	52.047		7	man
14) Carbon disulfide	2.294	76	118776	4.268		100	41126/21
15) Methyl Acetate	2.442	43	11019	4.774	ug/L	91	S 100
16) Methylene chloride	2.510	84	41488	3.855	ug/L	95	
17) Methyl tert-butyl Ether	2.770	73	79914	4.922	ug/L	96	
18) trans-1,2-Dichloroethene		96	41771	4.607		93	
19) 1,1-Dichloroethane	3.191	63	72457	4.734		98	
21) 2-Butanone	3.989	43	49302	37.389	0	100	
22) cis-1,2-Dichloroethene	3.915	96	42016		ug/L #	91	
23) Bromochloromethane	4.249	128	19438	4.831	ug/L #	79	

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111621\

Data File : VV023529.D

Acq On : 16 Nov 2021 13:12 Operator : SY/MD Sample : M4643-03MS Misc : 25.0mL/MSVOA_V/WATER ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 17 00:51:28 2021

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR110421WMA.M

Quant Title : TRACE VOA SFAM1.0 QLast Update : Wed Nov 17 00:48:57 2021 Response via : Initial Calibration

Instrument: MSVOA_V ClientSampleId: GB8J9MS

Manual IntegrationsAPPROVED

		**************************************				. Indiana de Managaria			
Co	mpound	R.T.	QIon	Response	Conc Un	its Dev	/(Min)		
25) Chlo	noform	4.378	83	81584	5.000	/I	99		
	Dichloroethane	5.133	62	42686	4.918	•	99		
	1-Trichloroethane	4.613	97	71957	4.830	•	99		
30) Cycl		4.680	56	59979	4.493	0.	98		
, ,	on tetrachloride	4.831	117	65435	4.493	O.	97		
33) Benz		5.101	78	165209	4.819	•	100		
	nloroethene	5.915	95	43556	4.819		96		
	/lcyclohexane	6.133	83	65565	4.777		95		
	Dichloropropane	6.175	63	38700	4.835		97		
, , , , , , , , , , , , , , , , , , , ,	odichloromethane	6.510	83	52662	4.833	0.	98		
			75		4.909	•			
	1,3-Dichloropropene	7.030		56143		0.	100		
	thyl-2-pentanone	7.230	43	204631	55.122	0.	99		
42) Tolu		7.387	91	185799	5.067		97		
	s-1,3-Dichloropropene	7.654	75	46831	4.902	-	99		
	2-Trichloroethane	7.841	97	28373	4.933	•	97		
	chloroethene	7.976	164	78220	9.899	0.	98		
48) 2-He		8.143	43	145116	55.787	•	97		
,	omochloromethane	8.246	129	36694	5.035	G.	97		
,	ibromoethane	8.352	107	26784	5.025	0	99		
51) Chlor		8.882	112	119079	4.885	0.	99		
52) Ethy		9.014	91	185995	4.809	-	100		
53) m,p->		9.140	106	75264	4.958		98		
54) o-xy		9.545	106	71397	5.014	-	100		
55) Styre		9.561	104	123863	5.077		97		
	,2-Tetrachloroethane	10.243	83	30294	4.808		96		
59) Bromo		9.731	173	19466		ug/L #			
	opylbenzene	9.931	105	190904	4.917	0.	99		
	-Trichloropropane	10.275	75	21571	4.799	0.	95		
	-Trimethylbenzene	10.538	105	156006	4.846	•	99		
	-Trimethylbenzene	10.915	105	157731	4.922		98		
	ichlorobenzene	11.181	146	96561	4.867	0.	98		
	ichlorobenzene	11.271	146	96766	4.776		98		
	ichlorobenzene	11.644	146	89422	5.037		99		
	ibromo-3-chloropr	12.429	75	4750	4.960		99		
	-Trichlorobenzene	12.644	180	72316	4.656		97		
	-trichlorobenzene	13.262	180	54818	4.407	0.	99		
71) Napht		13.503	128	74204	4.046		99		
72) 1,2,3	-Trichlorobenzene	13.744	180	50025	4.596	ug/L	99		

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