

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV080918\
 Data File : VV006964.D
 Acq On : 09 Aug 2018 21:47
 Operator : SY/MD
 Sample : J4399-02
 Misc : 25 mL/MSVOA V/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BC6R2

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR080918WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.029	29	43	68	rBV	3901206	4822328	100.00%	36.139%
2	1.318	124	133	145	rVB	122695	144903	3.00%	1.086%
3	1.585	209	216	229	rVB	102862	117370	2.43%	0.880%
4	2.131	377	386	397	rVV	183008	256398	5.32%	1.921%
5	2.209	400	410	429	rVB	34543	72915	1.51%	0.546%
6	3.228	719	727	747	rVB3	28873	64655	1.34%	0.485%
7	3.971	941	958	974	rBV	148079	424705	8.81%	3.183%
8	4.408	1077	1094	1126	rVB2	150967	377916	7.84%	2.832%
9	5.099	1291	1309	1320	rBV2	347878	824599	17.10%	6.180%
10	5.672	1471	1487	1503	rBV	287815	566579	11.75%	4.246%
11	5.922	1549	1565	1577	rBV4	37154	93667	1.94%	0.702%
12	6.125	1614	1628	1652	rBV	206451	425289	8.82%	3.187%
13	7.035	1899	1911	1929	rBV2	86564	154897	3.21%	1.161%
14	7.366	1999	2014	2028	rBV	323023	562766	11.67%	4.217%
15	7.671	2096	2109	2124	rBV2	50897	95697	1.98%	0.717%
16	8.147	2244	2257	2296	rBV2	373681	886932	18.39%	6.647%
17	8.903	2479	2492	2515	rBV2	411971	971912	20.15%	7.284%
18	9.060	2530	2541	2551	rBV	80622	127379	2.64%	0.955%
19	9.186	2569	2580	2598	rVV	122653	202313	4.20%	1.516%
20	9.282	2599	2610	2629	rVB3	28111	53594	1.11%	0.402%
21	9.594	2695	2707	2721	rBV	49683	83384	1.73%	0.625%
22	10.054	2837	2850	2863	rBV3	37600	63868	1.32%	0.479%
23	10.266	2902	2916	2931	rBV	150472	243419	5.05%	1.824%
24	10.395	2943	2956	2975	rBV4	56891	128267	2.66%	0.961%
25	11.302	3226	3238	3259	rBV	335230	610213	12.65%	4.573%
26	11.581	3309	3325	3338	rBV2	42212	86896	1.80%	0.651%
27	11.681	3344	3356	3384	rVB3	354562	784027	16.26%	5.876%
28	12.041	3457	3468	3481	rBV	57738	96922	2.01%	0.726%

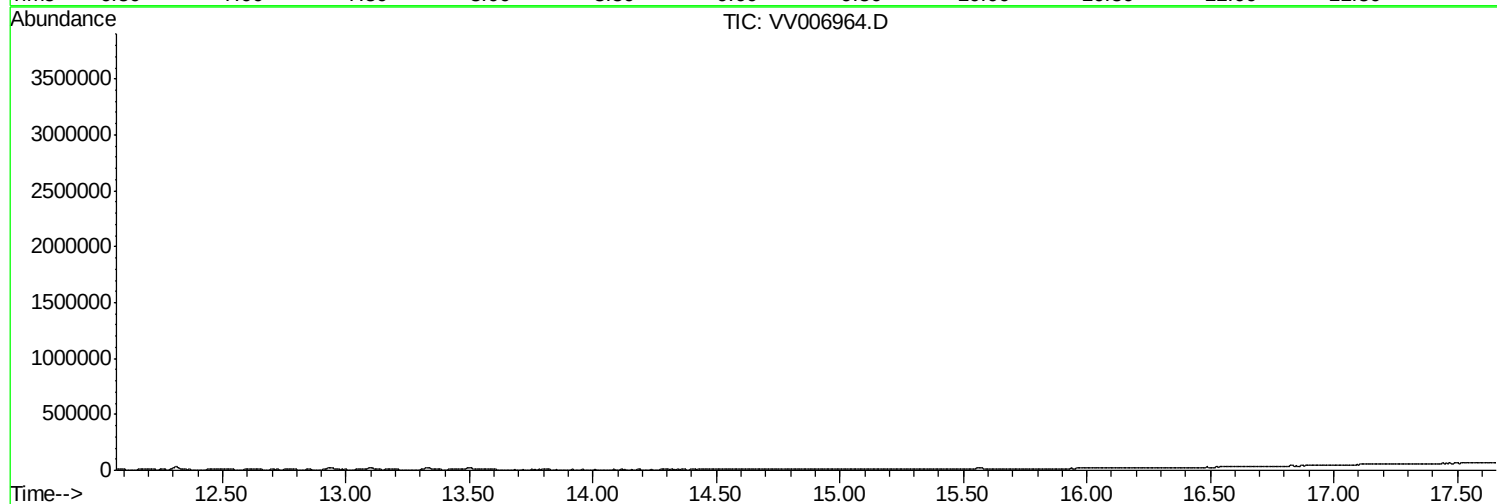
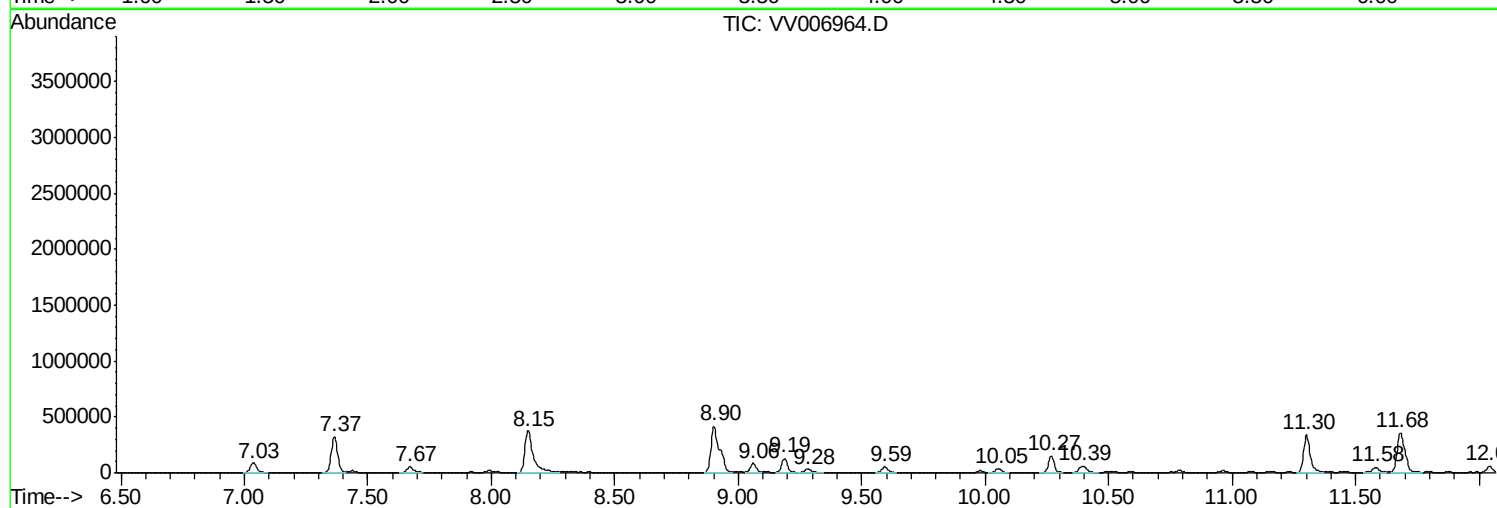
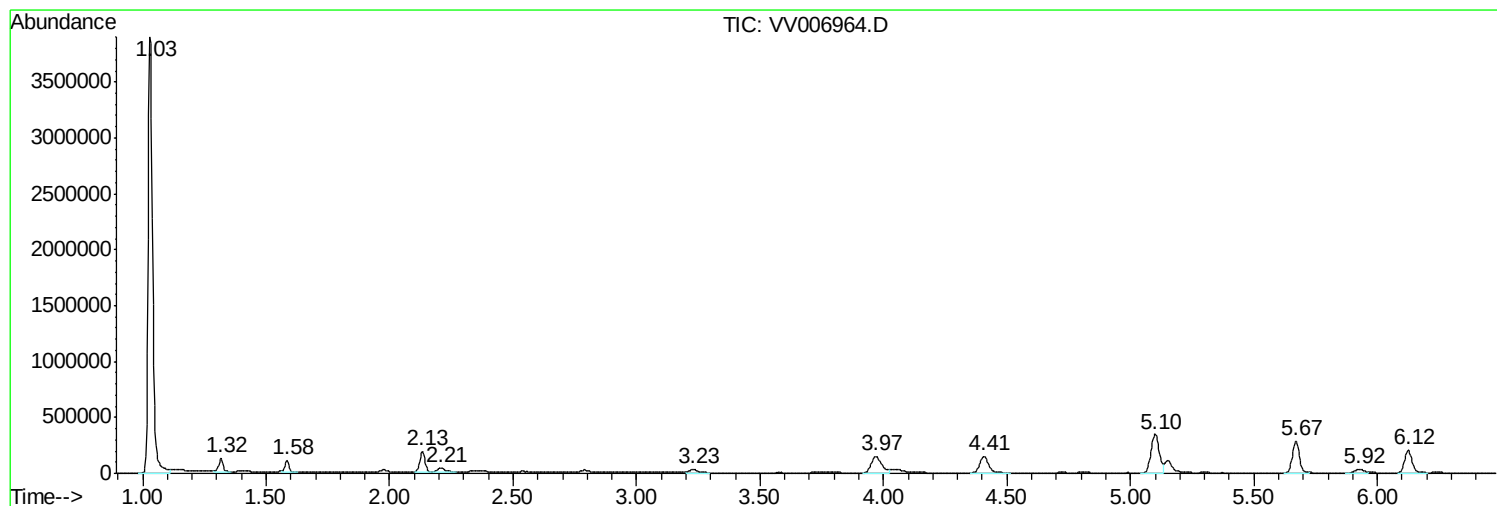
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ClientSampleId :
BC6R2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR080918WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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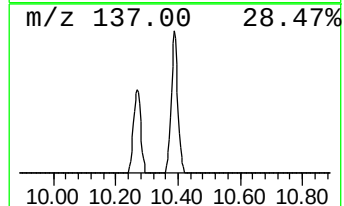
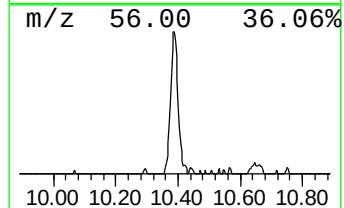
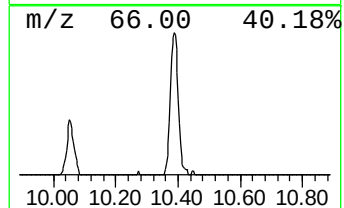
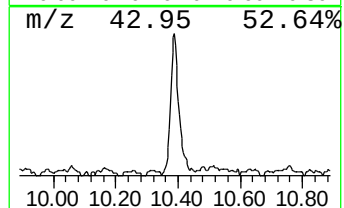
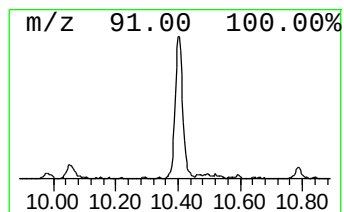
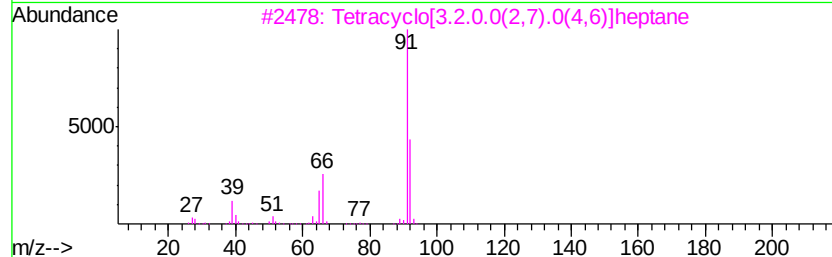
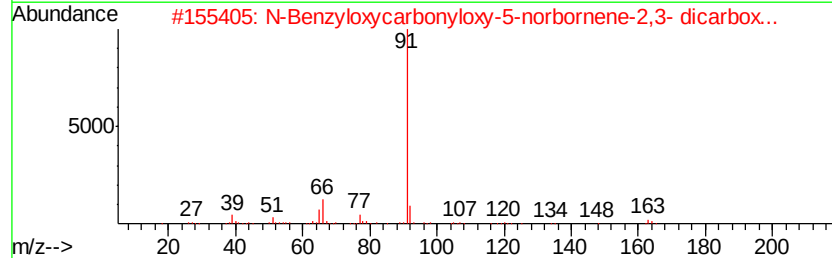
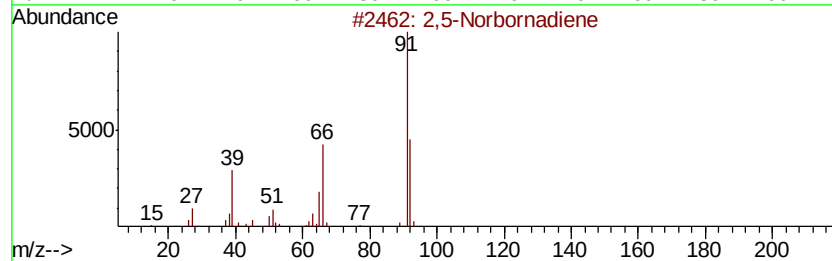
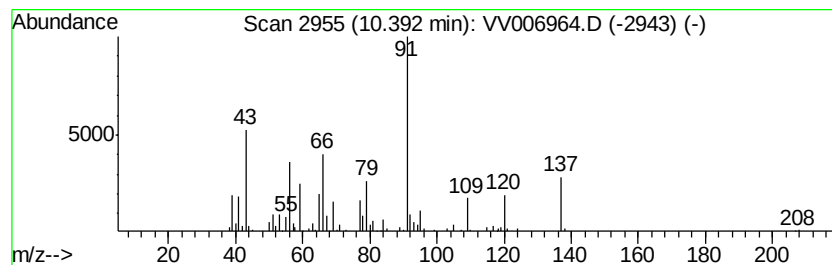
Instrument :
MSVOA_V
ClientSampleId :
BC6R2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR080918WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.39	1.05 ug/L	128267	1,4-Dichlorobenzene-d4	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Norbornadiene	92	C7H8	000121-46-0	38
2	N-Benzyloxycarbonyloxy-5-norborn...	313	C17H15N05	062210-73-5	37
3	Tetracyclo[3.2.0.0(2,7).0(4,6)]h...	92	C7H8	000278-06-8	32
4	Tetracyclo[3.2.0.0(2,7).0(4,6)]h...	92	C7H8	000278-06-8	32
5	2-(Hydroxymethyl)norbornadiene	122	C8H10O	067583-61-3	32



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Instrument :
MSVOA_V
ClientSampled :
BC6R2

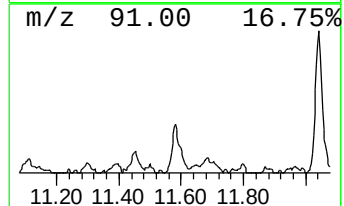
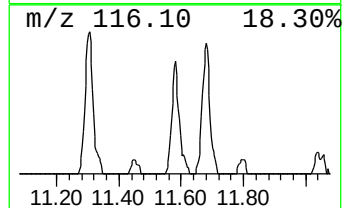
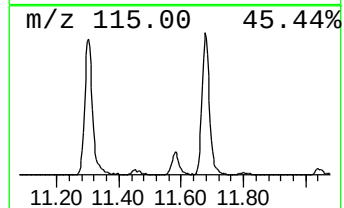
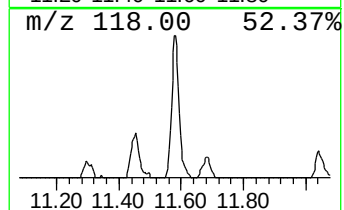
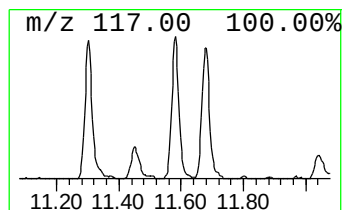
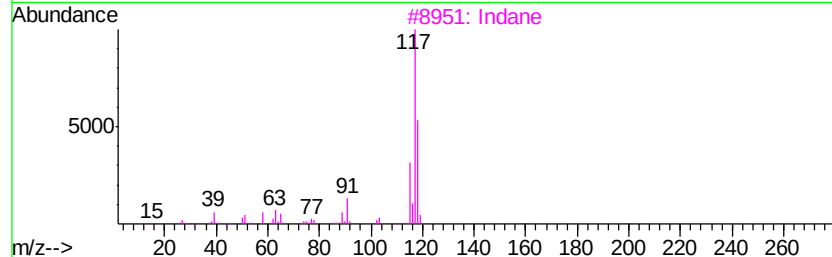
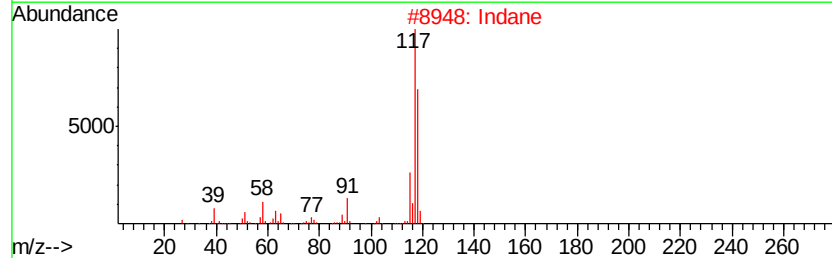
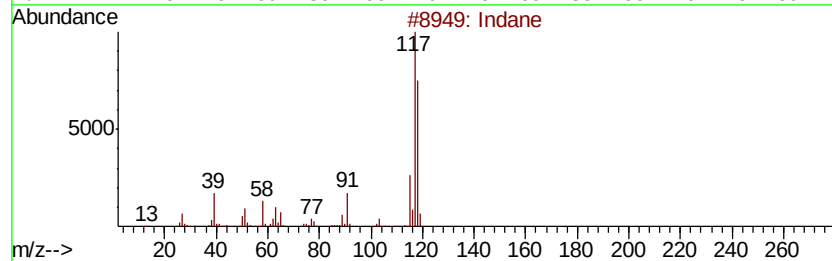
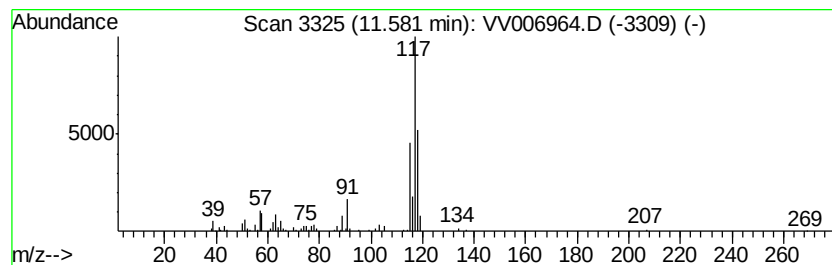
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.58	0.71 ug/L	86896	1,4-Dichlorobenzene-d4	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Indane		118	C9H10	000496-11-7	92
3	Indane		118	C9H10	000496-11-7	81
4	trans-Cinnamyl bromide		196	C9H9Br	026146-77-0	64
5	3-Bromo-1-phenyl-1-propene		196	C9H9Br	004392-24-9	59



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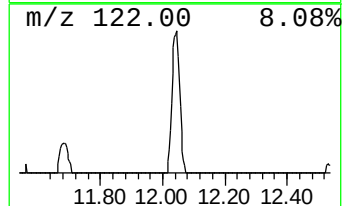
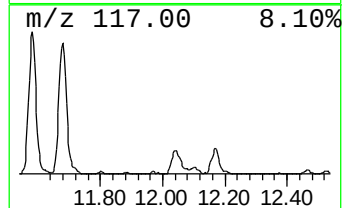
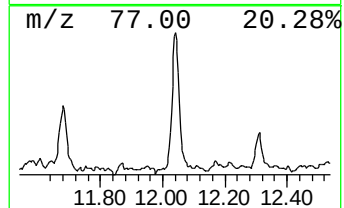
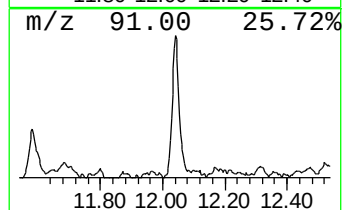
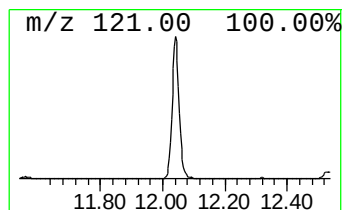
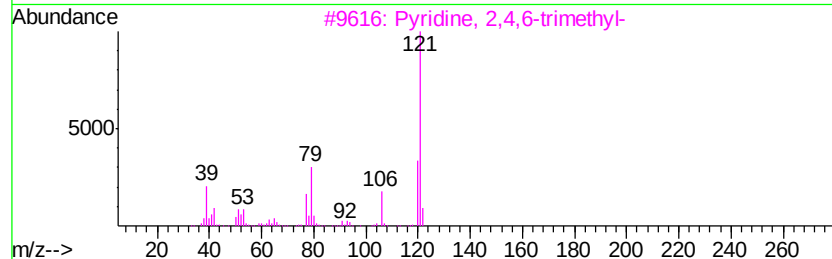
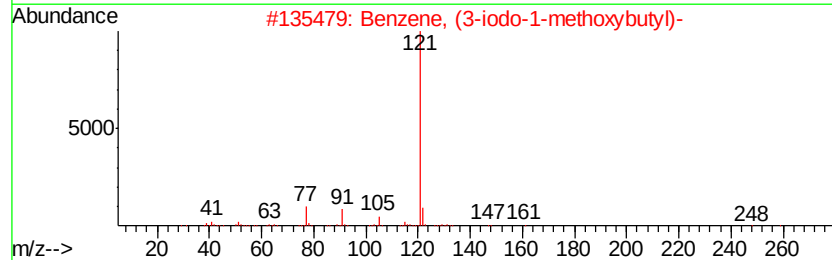
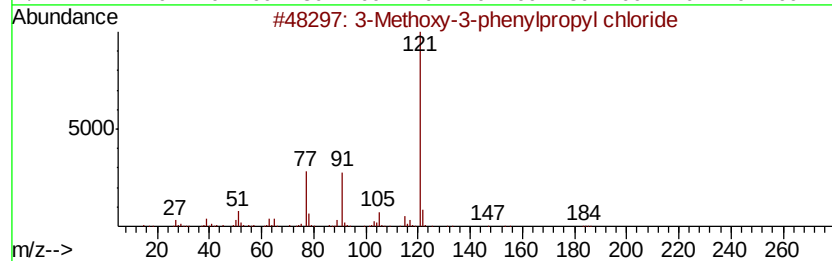
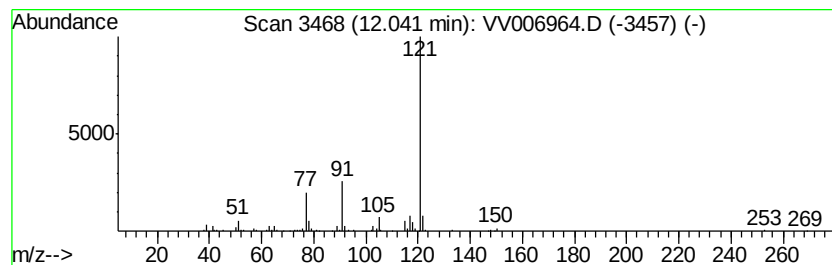
Instrument :
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ClientSampleId :
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 3-Methoxy-3-phenylpropyl ch... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.04	0.79 ug/L	96922	1,4-Dichlorobenzene-d4	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Methoxy-3-phenylpropyl chloride	184	C10H13ClO	1000370-35-5	80
2	Benzene, (3-iodo-1-methoxybutyl)-	290	C11H15IO	1000327-38-8	72
3	Pyridine, 2,4,6-trimethyl-	121	C8H11N	000108-75-8	64
4	Benzene, (3-iodo-1-methoxypropyl)-	276	C10H13IO	1000327-31-9	64
5	Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	64



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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	10.39	1.1	ug/L	128267	3	11.30	610213	5.0
Indane	11.58	0.7	ug/L	86896	3	11.30	610213	5.0
3-Methoxy-3-pheny...	12.04	0.8	ug/L	96922	3	11.30	610213	5.0