

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
 Data File : VN070371.D
 Acq On : 30 Dec 2021 2:29
 Operator : JC/MD
 Sample : M5219-03
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 N48854-VOC

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Title : SW846 8260

Signal : TIC: VN070371.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.293	833	859	897	rBV2	130789	436812	10.08%	0.991%
2	4.588	950	969	970	rBV	95719	147530	3.41%	0.335%
3	6.498	1663	1681	1705	rVB5	28720	94327	2.18%	0.214%
4	7.691	2108	2126	2148	rBV5	85511	228069	5.27%	0.517%
5	8.021	2228	2249	2261	rBV2	643187	1552836	35.85%	3.523%
6	8.086	2262	2273	2305	rVB	1056809	2370587	54.73%	5.378%
7	8.440	2384	2405	2432	rBV2	869735	2255547	52.07%	5.117%
8	8.968	2582	2602	2625	rBV	1711085	3527080	81.42%	8.001%
9	9.628	2836	2848	2860	rBV2	27319	49077	1.13%	0.111%
10	10.331	3097	3110	3127	rBV5	25756	52815	1.22%	0.120%
11	10.443	3132	3152	3167	rBV	2347058	4331740	100.00%	9.826%
12	10.505	3167	3175	3199	rVB	175409	339846	7.85%	0.771%
13	11.090	3374	3393	3410	rBV2	74077	141700	3.27%	0.321%
14	11.744	3620	3637	3664	rBV	2015364	3743101	86.41%	8.491%
15	11.948	3703	3713	3725	rBV2	62323	111044	2.56%	0.252%
16	12.278	3828	3836	3856	rVB3	45567	84941	1.96%	0.193%
17	12.731	3989	4005	4028	rBV	1468105	2589514	59.78%	5.874%
18	13.055	4118	4126	4138	rVB8	47171	81177	1.87%	0.184%
19	13.369	4233	4243	4253	rBV	57324	89053	2.06%	0.202%
20	13.495	4280	4290	4303	rVB6	73950	127736	2.95%	0.290%
21	13.675	4342	4357	4374	rBV	1192893	2175986	50.23%	4.936%
22	13.836	4407	4417	4423	rBV2	133217	213944	4.94%	0.485%
23	13.876	4424	4432	4443	rVB	295570	470840	10.87%	1.068%
24	13.999	4467	4478	4489	rBV6	107703	200295	4.62%	0.454%
25	14.131	4519	4527	4533	rBV2	66747	105358	2.43%	0.239%
26	14.217	4548	4559	4572	rVV	814105	1236897	28.55%	2.806%
27	14.278	4575	4582	4589	rVV2	112612	170105	3.93%	0.386%
28	14.327	4589	4600	4614	rVB2	509808	887344	20.48%	2.013%
29	14.463	4646	4651	4659	rVB3	71235	89599	2.07%	0.203%
30	14.541	4662	4680	4689	rBV	784529	1445867	33.38%	3.280%
31	14.619	4704	4709	4718	rVB3	88820	109468	2.53%	0.248%
32	14.809	4770	4780	4790	rBV2	478679	799577	18.46%	1.814%
33	14.855	4792	4797	4807	rVB6	157338	213448	4.93%	0.484%
34	14.922	4807	4822	4836	rBV4	1653856	3807841	87.91%	8.638%
35	15.196	4903	4924	4933	rBV6	632668	1386956	32.02%	3.146%
36	15.308	4954	4966	4990	rVB3	636549	1569876	36.24%	3.561%

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37	15.504	5015	5039	5053	rBV	1486321	2950903	68.12%	6.694%
38	15.614	5071	5080	5090	rBV5	229469	404723	9.34%	0.918%
39	15.804	5136	5151	5168	rBV	626043	1297150	29.95%	2.943%
40	15.981	5206	5217	5223	rBV3	375179	701453	16.19%	1.591%
41	16.110	5255	5265	5277	rBV5	227445	404879	9.35%	0.918%
42	16.344	5338	5352	5366	rBV5	405123	903881	20.87%	2.050%
43	16.614	5445	5453	5462	rBV3	108168	181356	4.19%	0.411%

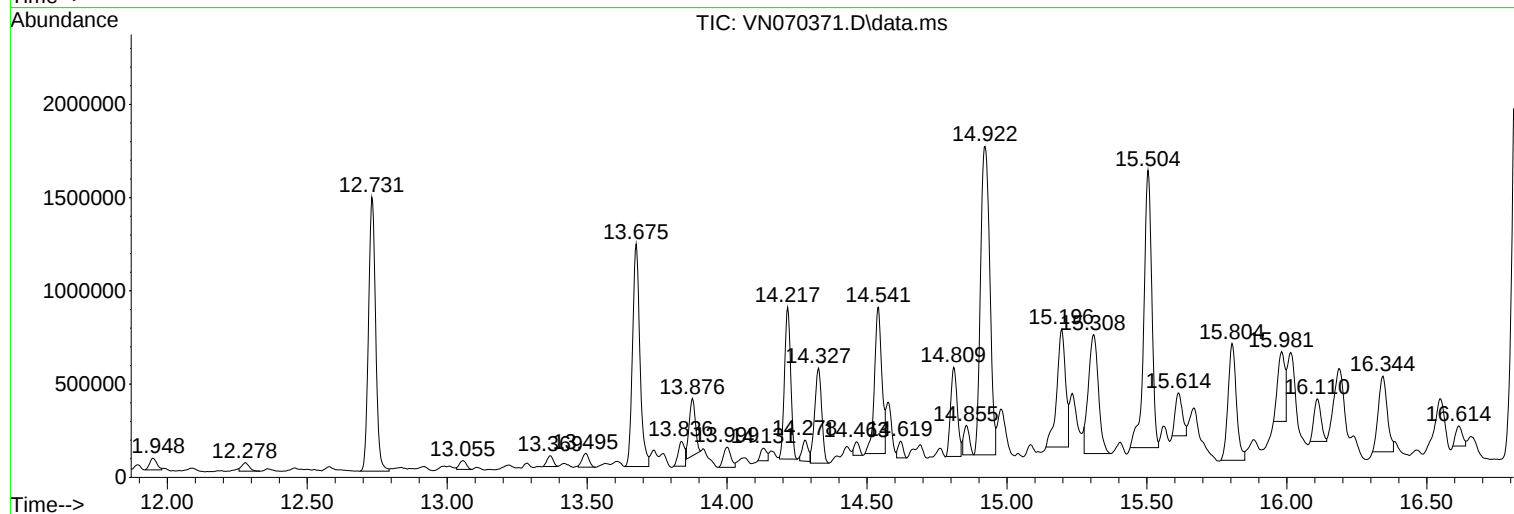
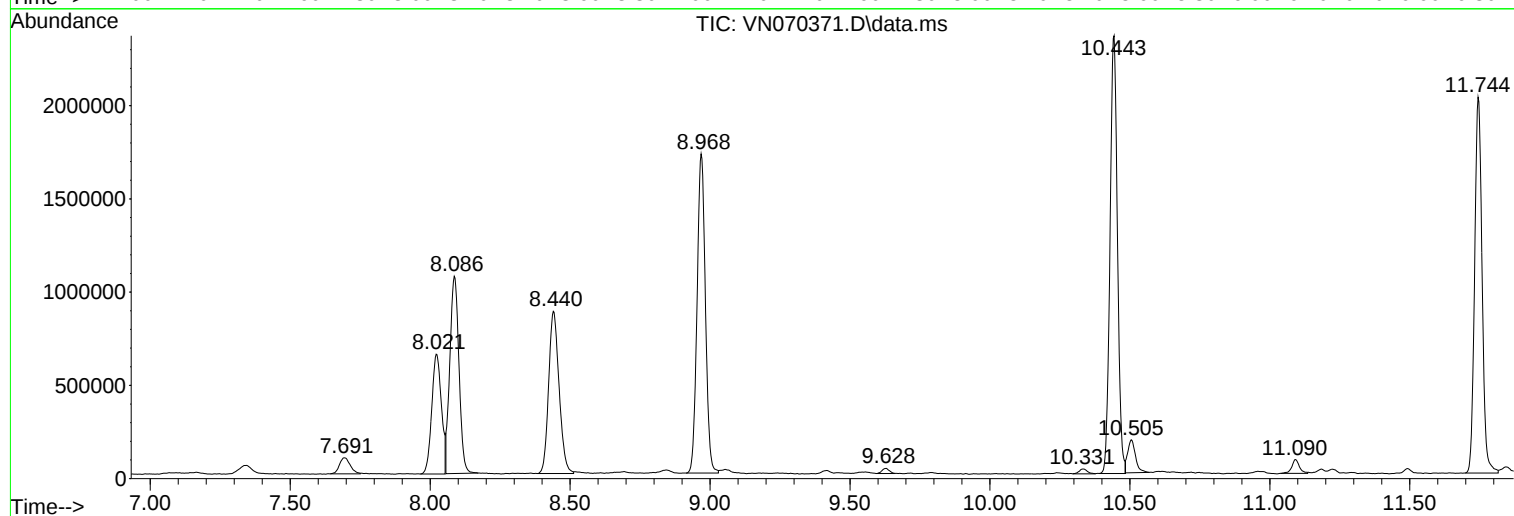
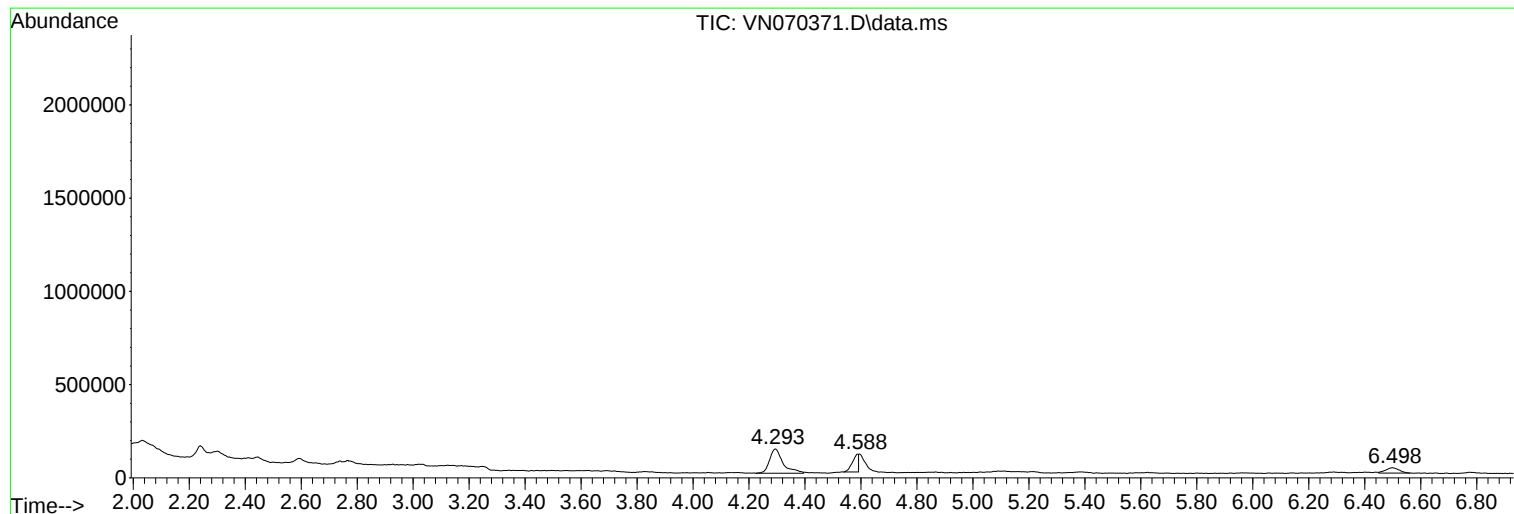
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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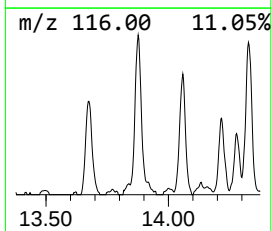
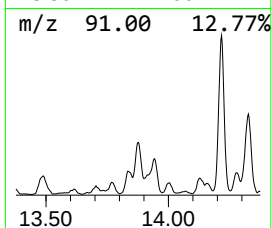
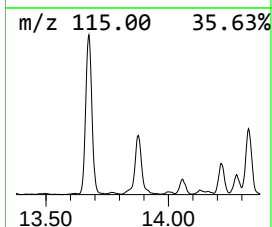
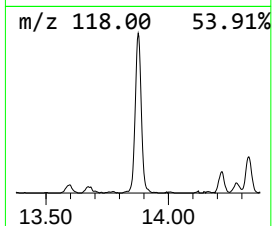
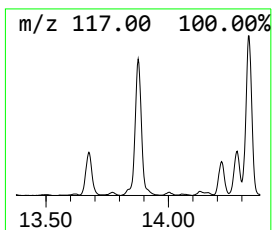
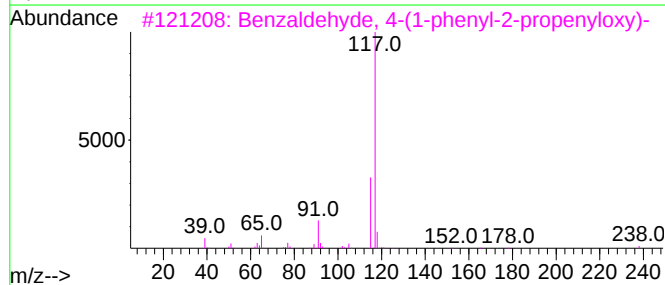
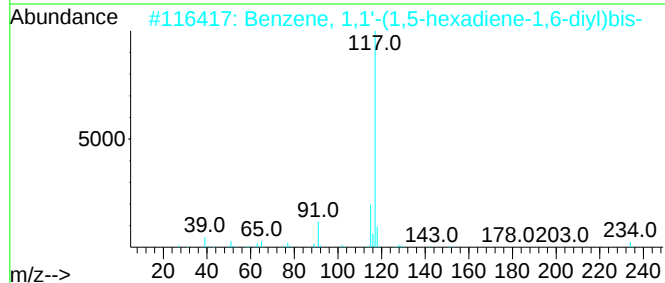
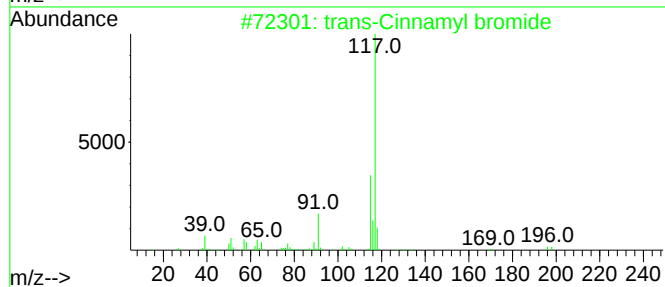
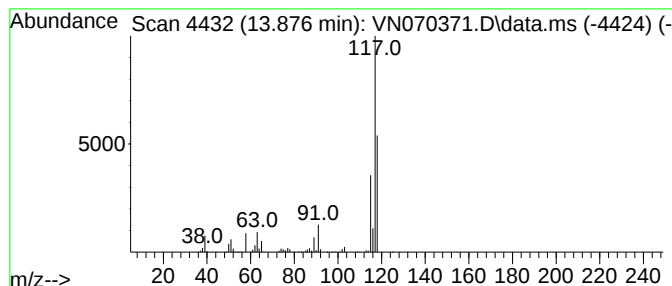
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 trans-Cinnamyl bromide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.876	10.82 ug/l	470840	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	53
2			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	50
3			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	50
4			Indole	117	C8H7N	000120-72-9	49
5			Benzene, (isocyanomethyl)-	117	C8H7N	010340-91-7	47



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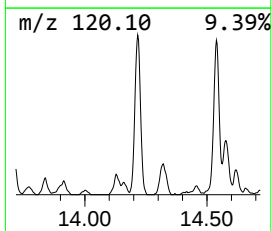
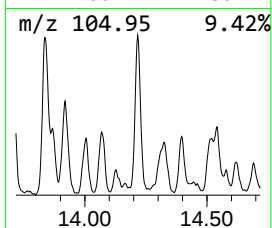
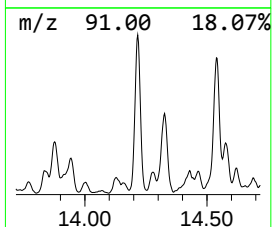
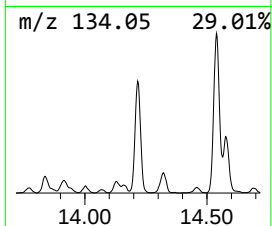
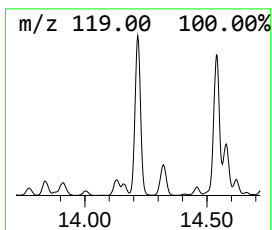
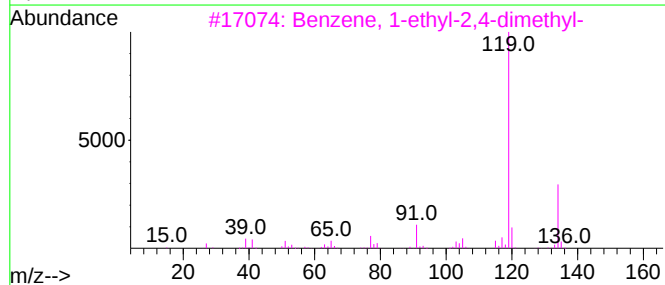
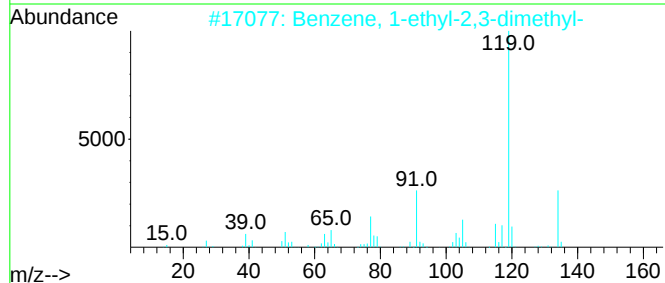
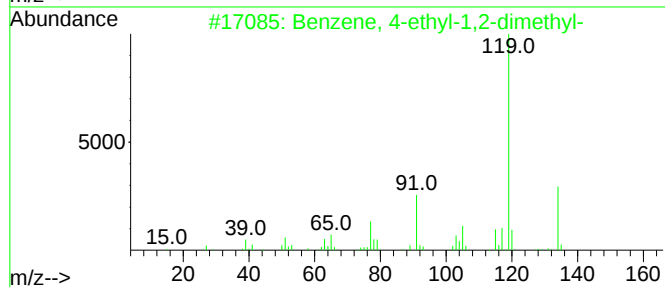
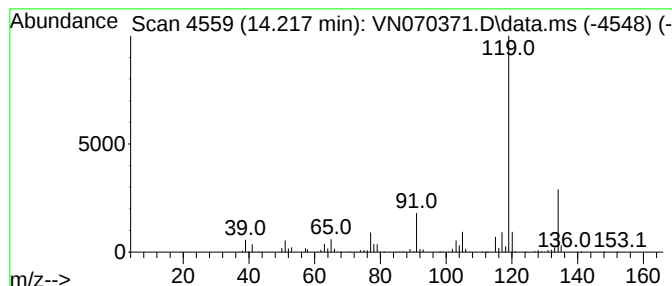
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.217	28.42 ug/l	1236900	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



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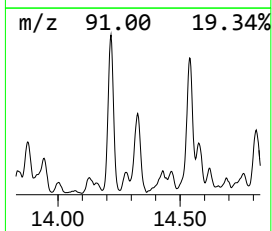
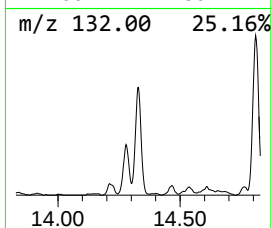
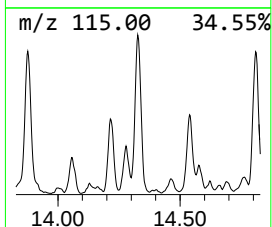
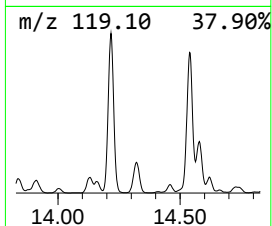
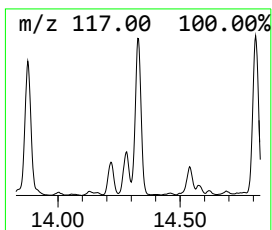
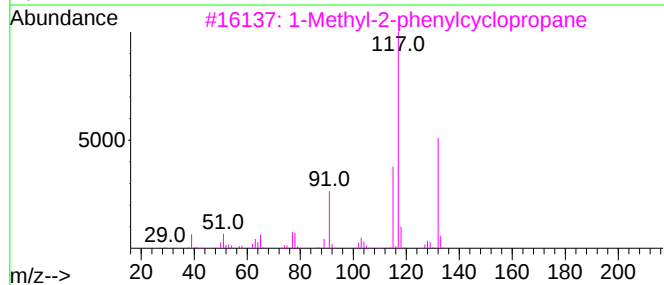
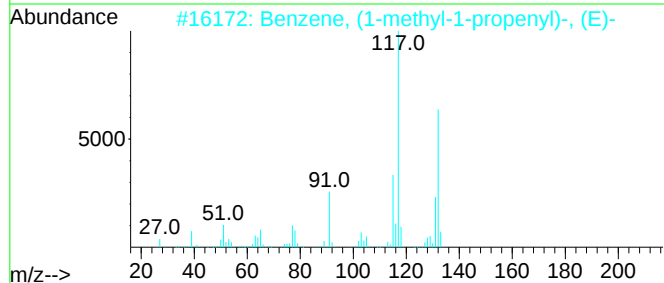
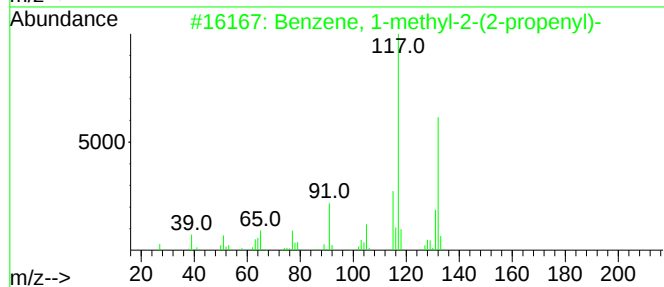
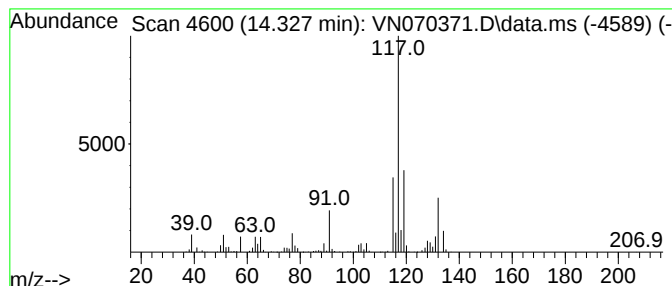
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-methyl-2-(2-prop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.327	20.39 ug/l	887344	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	72
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	70
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64
5			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64



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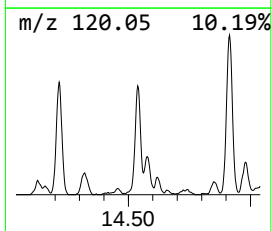
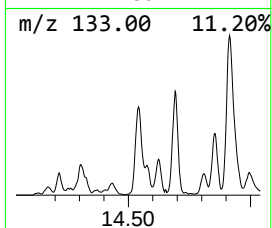
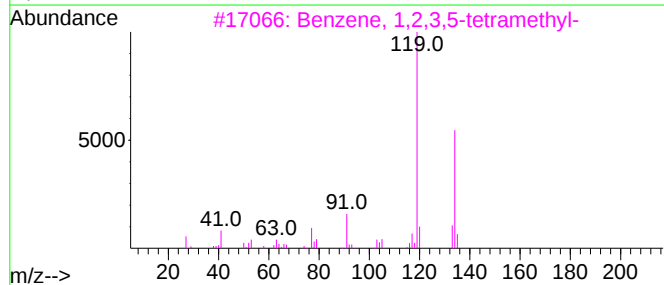
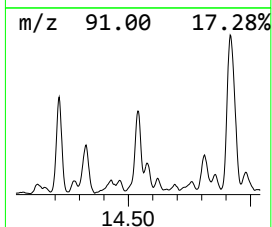
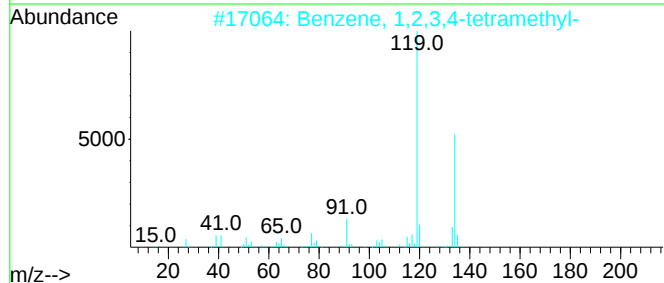
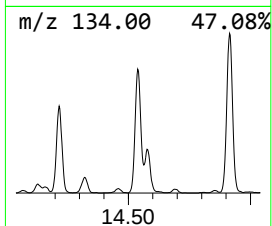
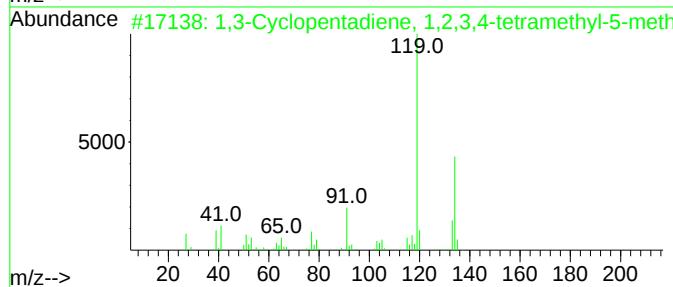
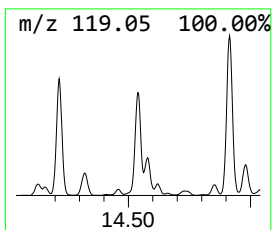
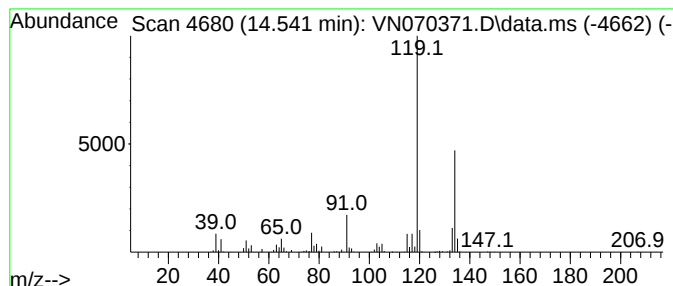
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.541	33.22 ug/l	1445870	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



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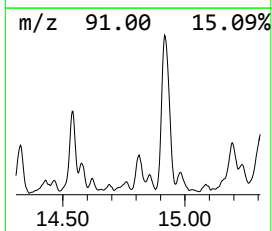
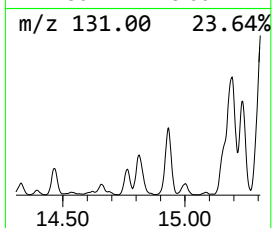
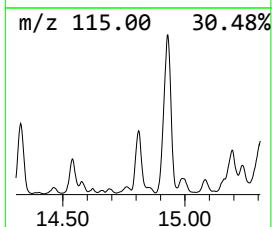
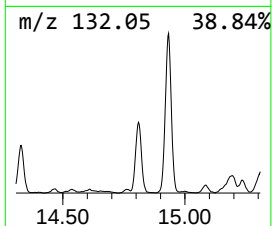
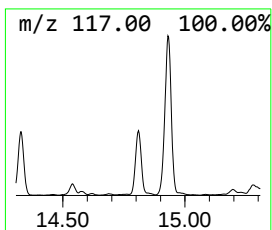
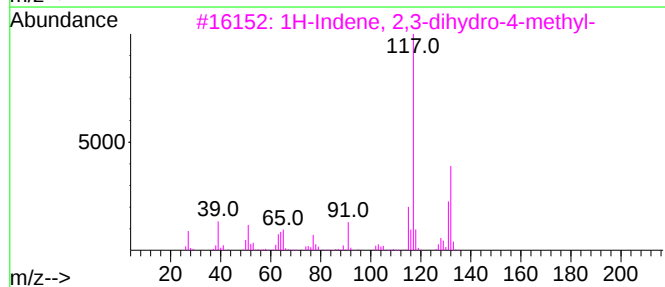
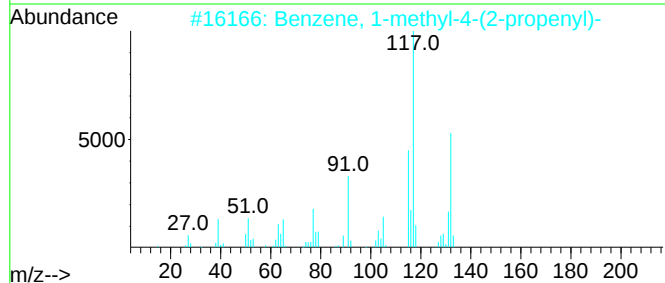
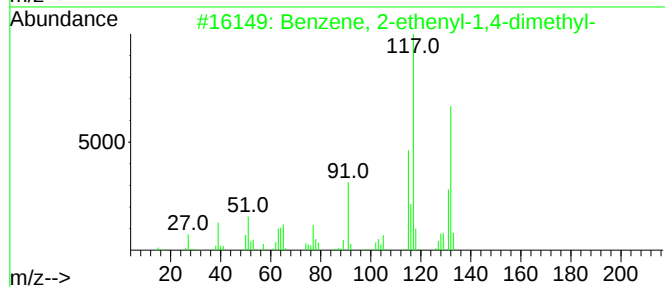
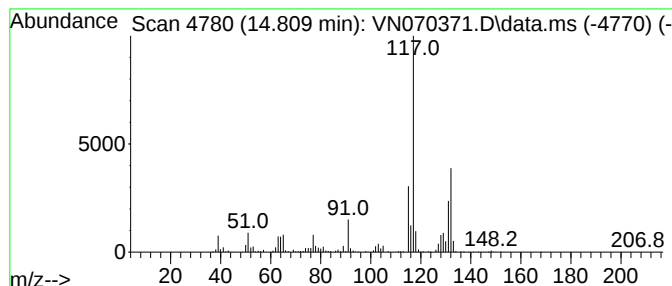
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.809	18.37 ug/l	799577	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	90
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

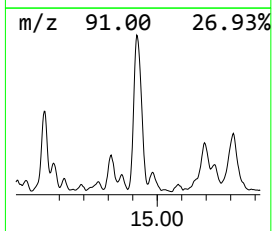
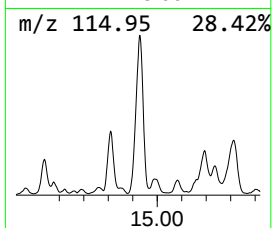
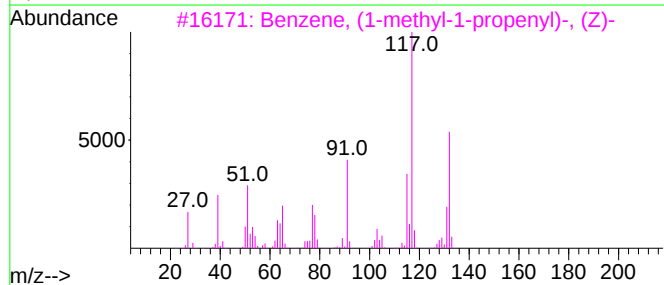
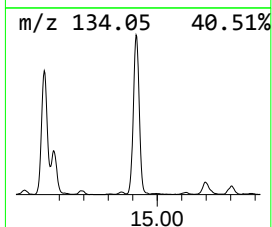
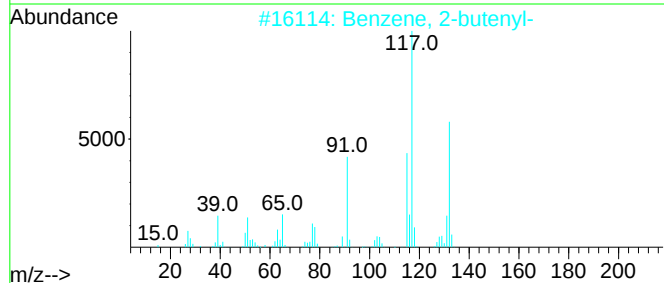
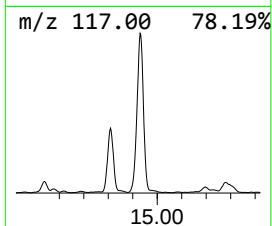
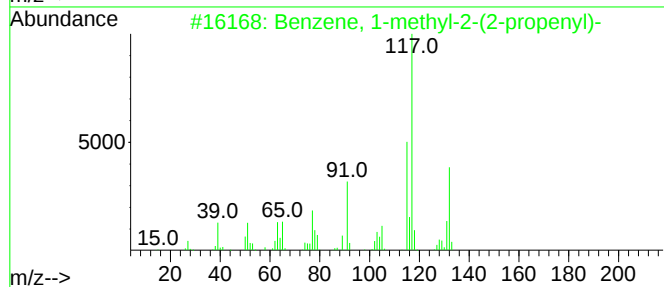
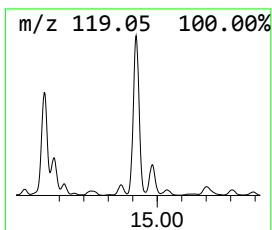
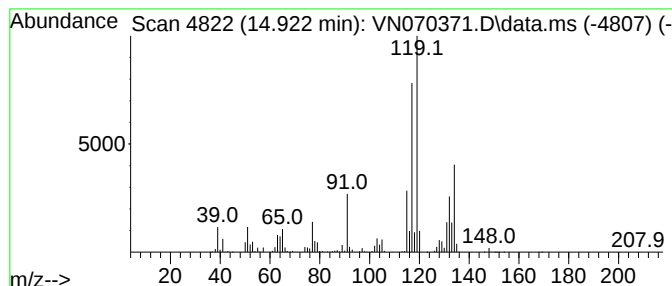
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-butenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.922	87.50 ug/l	3807840	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	50
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

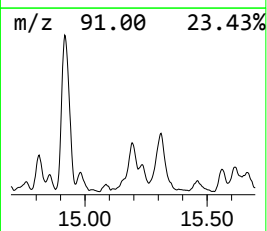
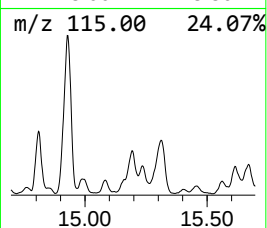
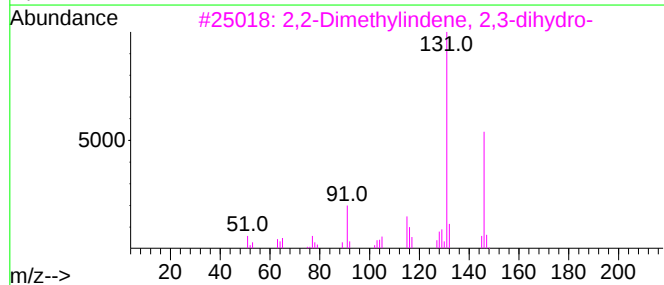
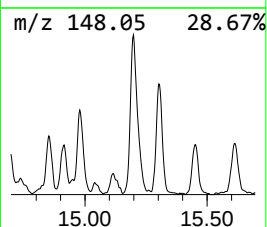
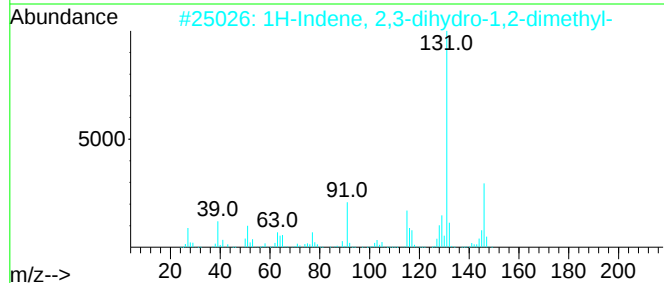
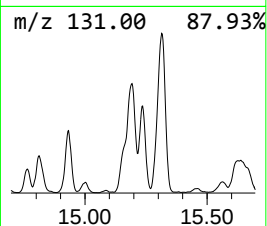
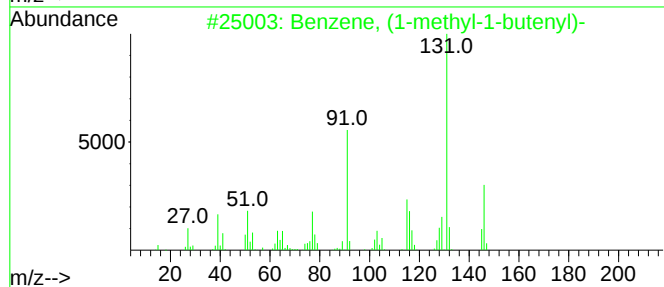
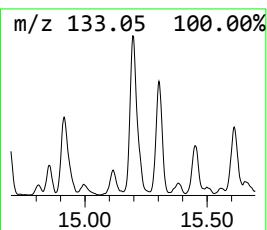
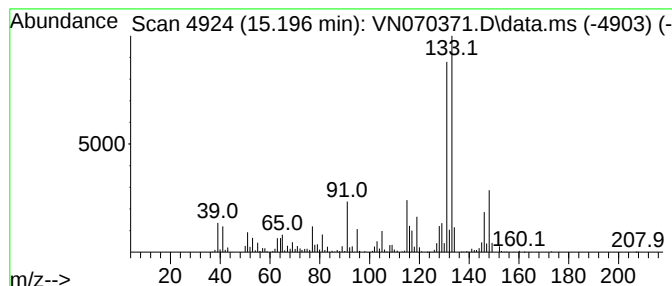
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, (1-methyl-1-butenyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.196	31.87 ug/l	1386960	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
5			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

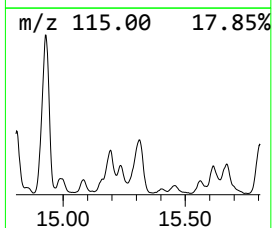
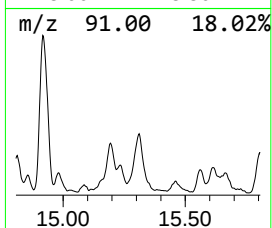
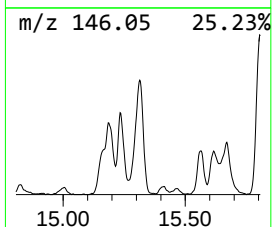
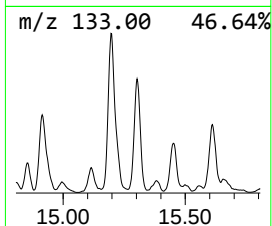
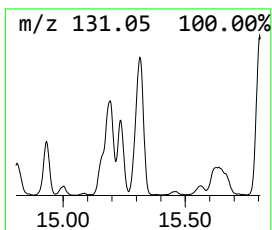
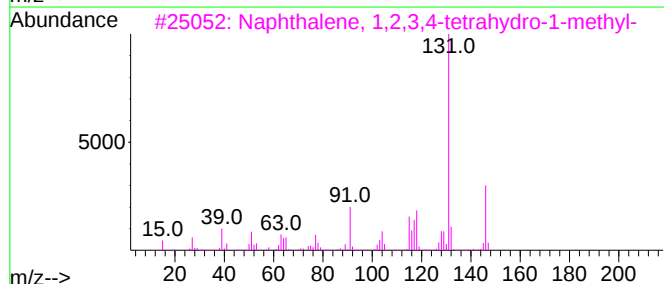
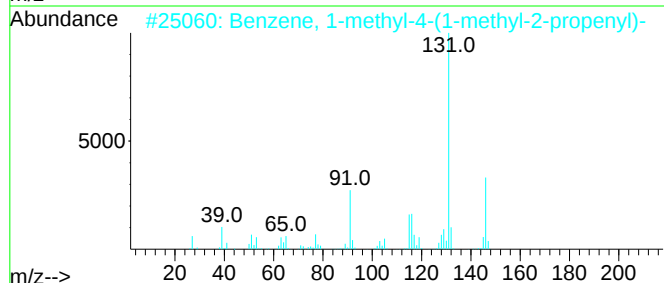
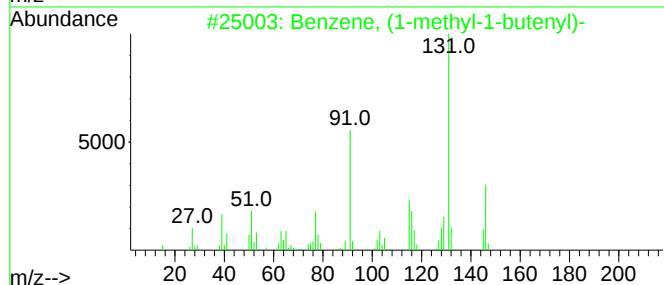
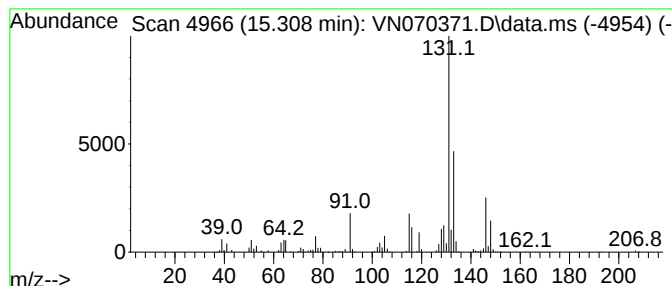
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-methyl-4-(1-meth... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.308	36.07 ug/l	1569880	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	83
2			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
4			1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	64
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

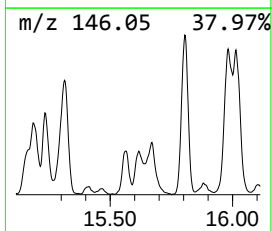
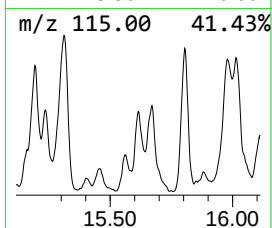
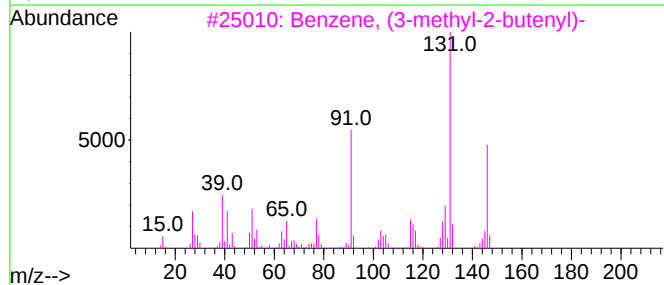
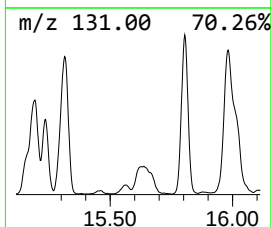
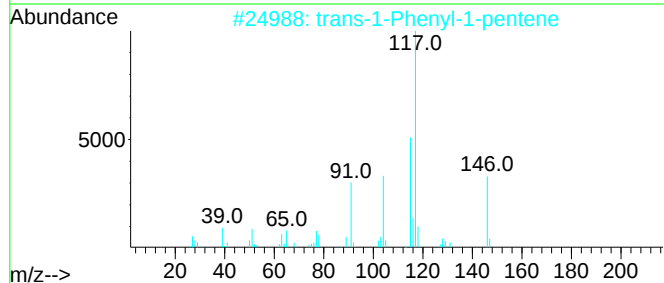
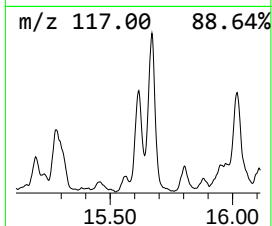
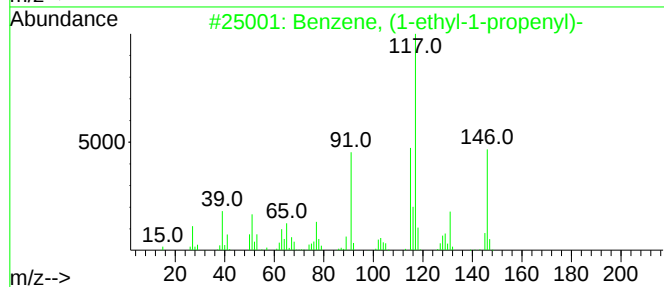
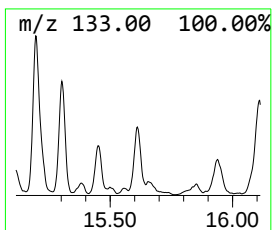
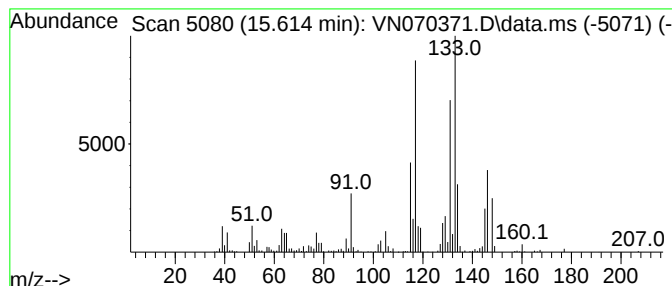
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown15.614 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.614	9.30 ug/l	404723	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	46
2			trans-1-Phenyl-1-pentene	146	C11H14	016002-93-0	42
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	35
4			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	30
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

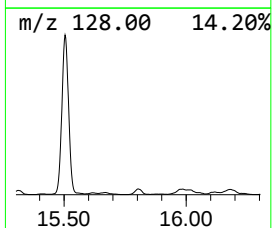
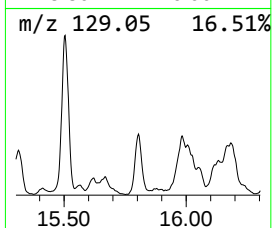
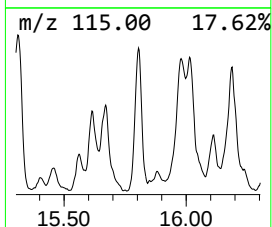
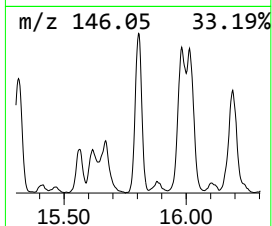
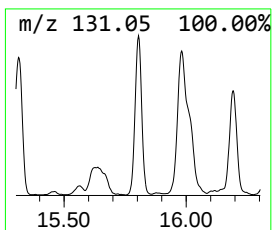
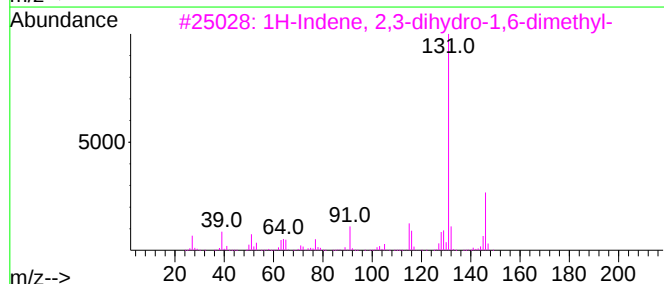
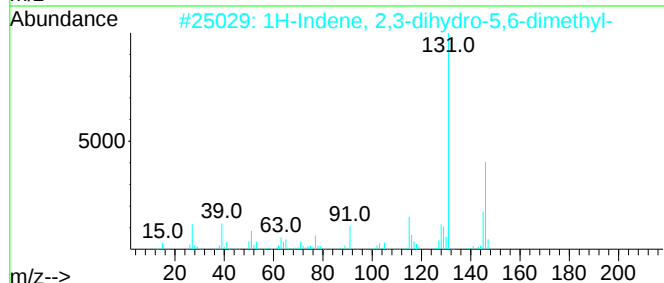
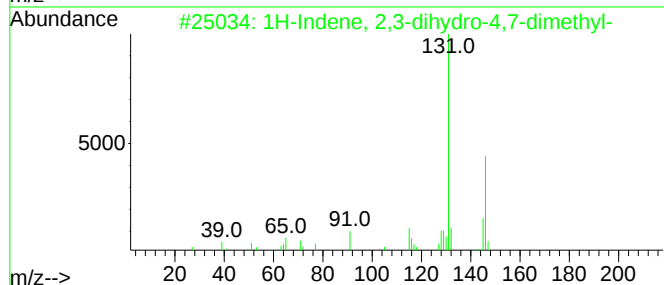
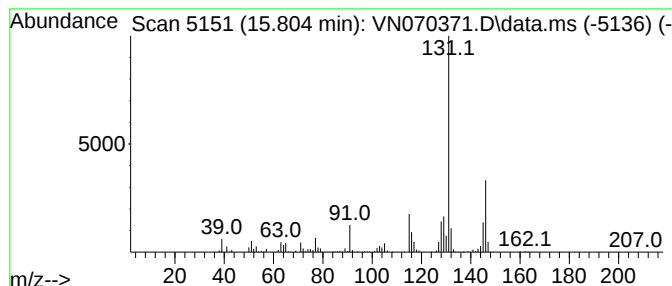
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.804	29.81 ug/l	1297150	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96		
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	94		
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94		
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91		
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

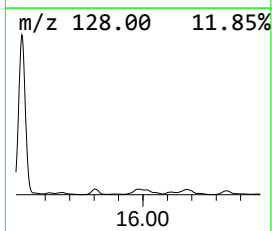
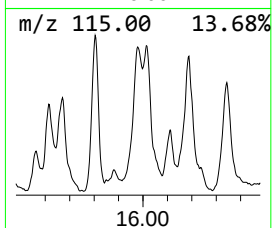
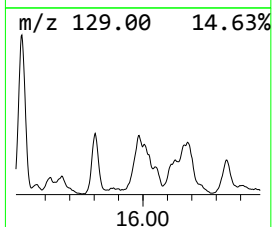
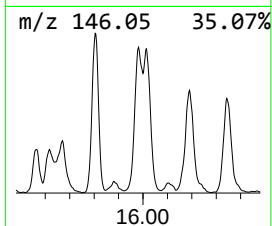
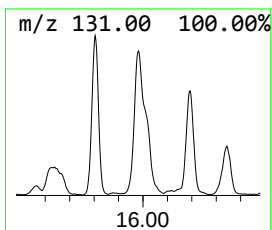
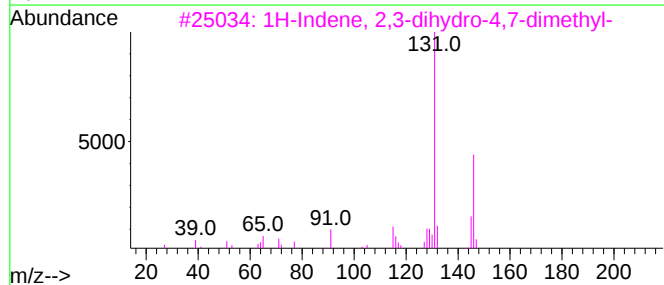
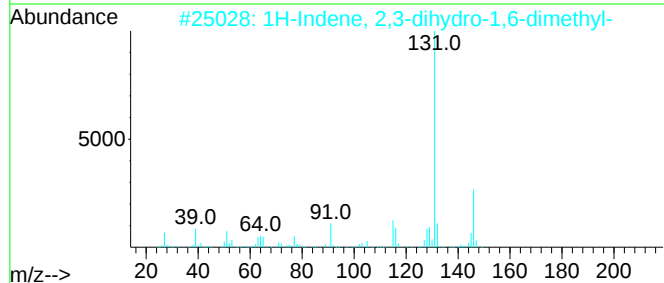
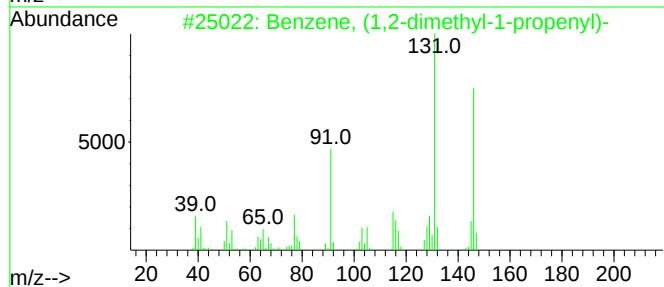
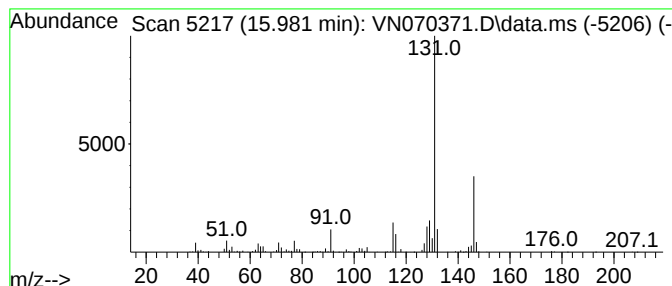
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.981	16.12 ug/l	701453	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
4			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

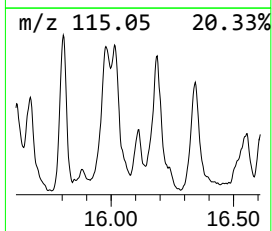
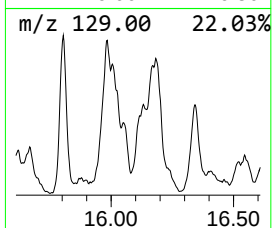
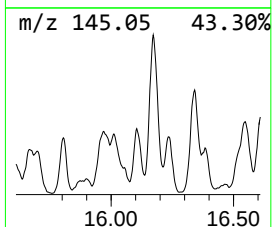
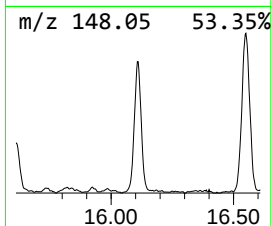
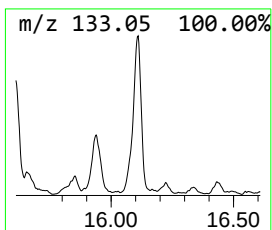
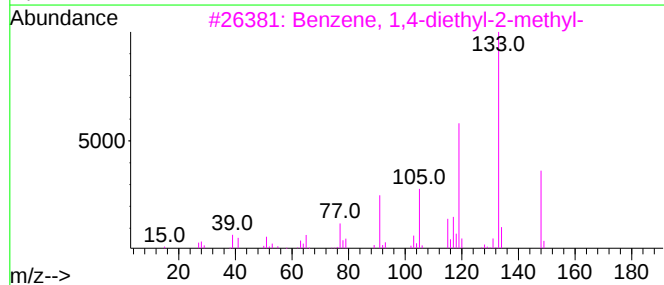
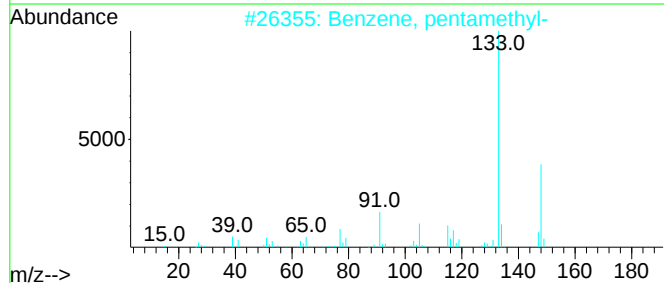
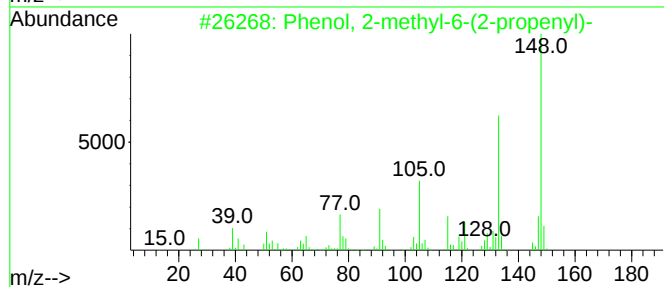
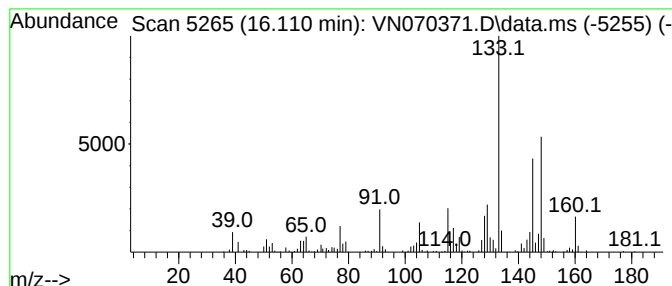
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Phenol, 2-methyl-6-(2-prope... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.110	9.30 ug/l	404879	1,4-Dichlorobenzene-d4	13.675

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenol, 2-methyl-6-(2-propenyl)-	148	C10H12O	003354-58-3	58
2		Benzene, pentamethyl-	148	C11H16	000700-12-9	55
3		Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	49
4		Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	46
5		3,4-Dimethylcumene	148	C11H16	1000370-34-1	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

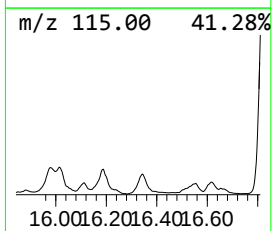
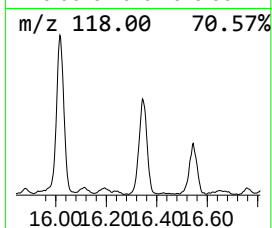
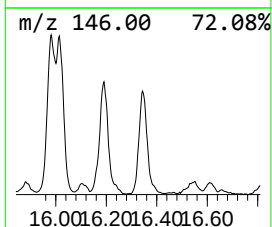
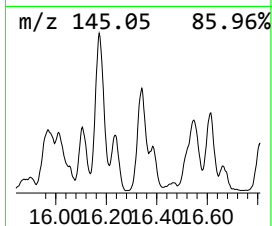
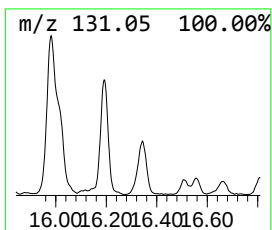
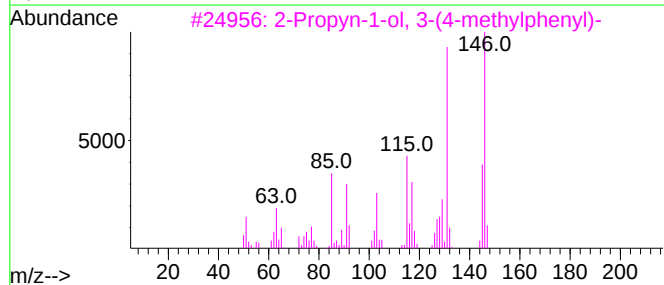
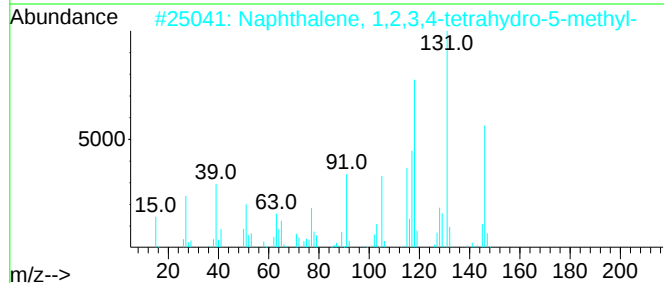
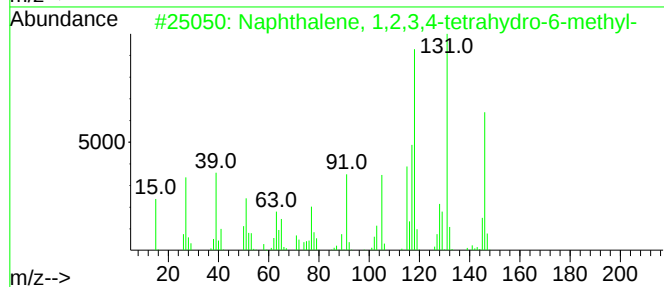
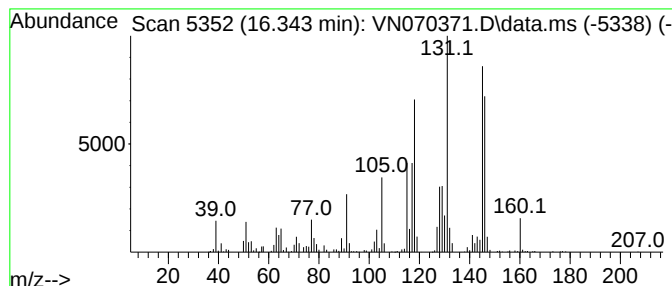
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.343	20.77 ug/l	903881	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	83
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	70
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	58
4			Benzene, 1-(1-methylethenyl)-2-(...	160	C12H16	005557-93-7	55
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
Data File : VN070371.D
Acq On : 30 Dec 2021 2:29
Operator : JC/MD
Sample : M5219-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 37 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
N48854-VOC

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
trans-Cinnamyl ...	13.876	10.8	ug/l	470840	4	13.675	2175990	50.0
Benzene, 4-ethy...	14.217	28.4	ug/l	1236900	4	13.675	2175990	50.0
Benzene, 1-meth...	14.327	20.4	ug/l	887344	4	13.675	2175990	50.0
1,3-Cyclopentad...	14.541	33.2	ug/l	1445870	4	13.675	2175990	50.0
Benzene, 2-ethe...	14.809	18.4	ug/l	799577	4	13.675	2175990	50.0
Benzene, 2-bute...	14.922	87.5	ug/l	3807840	4	13.675	2175990	50.0
Benzene, (1-met...	15.196	31.9	ug/l	1386960	4	13.675	2175990	50.0
Benzene, 1-meth...	15.308	36.1	ug/l	1569880	4	13.675	2175990	50.0
unknown15.614	15.614	9.3	ug/l	404723	4	13.675	2175990	50.0
1H-Indene, 2,3-...	15.804	29.8	ug/l	1297150	4	13.675	2175990	50.0
Benzene, (1,2-d...	15.981	16.1	ug/l	701453	4	13.675	2175990	50.0
Phenol, 2-methy...	16.110	9.3	ug/l	404879	4	13.675	2175990	50.0
Naphthalene, 1,...	16.343	20.8	ug/l	903881	4	13.675	2175990	50.0