

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
 Data File : VY015451.D
 Acq On : 06 Sep 2023 15:44
 Operator : SY/MD
 Sample : 04299-05
 Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SB-03-9.5-10.0

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_Y\methods\82Y082923S.M

Title : SW846 8260

Signal : TIC: VY015451.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.149	1021	1038	1054	rBV	4009177	9062236	10.13%	2.152%
2	8.679	1117	1125	1142	rBV	442920	903500	1.01%	0.215%
3	10.173	1362	1370	1375	rBV	793225	1376181	1.54%	0.327%
4	11.483	1578	1585	1592	rBV	647046	1061673	1.19%	0.252%
5	11.587	1595	1602	1612	rVV	21602386	33753645	37.71%	8.015%
6	11.697	1612	1620	1630	rVB	11766834	21375220	23.88%	5.075%
7	12.026	1667	1674	1685	rVB	13011202	19477779	21.76%	4.625%
8	12.325	1717	1723	1729	rBV	1834248	2762332	3.09%	0.656%
9	12.733	1784	1790	1792	rBV	4047356	6006174	6.71%	1.426%
10	12.764	1792	1795	1799	rVV	5389442	7712104	8.62%	1.831%
11	12.806	1799	1802	1809	rVB	3077047	4499928	5.03%	1.068%
12	12.971	1821	1829	1837	rBV	1822198	2940285	3.29%	0.698%
13	13.117	1843	1853	1858	rBV	7147076	11350222	12.68%	2.695%
14	13.166	1858	1861	1867	rVB	1288860	1776967	1.99%	0.422%
15	13.312	1878	1885	1891	rBV2	788086	1555996	1.74%	0.369%
16	13.453	1904	1908	1912	rVV	3153654	4820538	5.39%	1.145%
17	13.495	1912	1915	1924	rVB	1990001	2917716	3.26%	0.693%
18	13.629	1924	1937	1949	rBV	36519872	72995463	81.56%	17.333%
19	13.806	1960	1966	1973	rVB	1021837	1694831	1.89%	0.402%
20	13.971	1987	1993	1997	rBV	1472052	2202926	2.46%	0.523%
21	14.080	2005	2011	2012	rBV2	717339	1029679	1.15%	0.244%
22	14.105	2012	2015	2020	rVV	1239653	1878145	2.10%	0.446%
23	14.330	2049	2052	2057	rBV	736300	997866	1.11%	0.237%
24	14.562	2084	2090	2096	rBV	4536261	6731344	7.52%	1.598%
25	14.678	2102	2109	2114	rBV3	1259009	2496882	2.79%	0.593%
26	14.745	2114	2120	2128	rVV	5359112	8550404	9.55%	2.030%
27	14.830	2128	2134	2140	rVV	11872374	18339501	20.49%	4.355%
28	14.928	2140	2150	2158	rVV2	1260398	2643147	2.95%	0.628%
29	15.245	2182	2202	2210	rBV3	36344204	89497125	100.00%	21.251%
30	15.330	2210	2216	2225	rVB	1579726	2722655	3.04%	0.646%
31	15.562	2250	2254	2266	rVB	308699	966006	1.08%	0.229%
32	15.708	2272	2278	2283	rBV	630747	1241931	1.39%	0.295%
33	15.830	2291	2298	2312	rVB	663465	1920118	2.15%	0.456%
34	16.293	2366	2374	2386	rVV	21958967	43407268	48.50%	10.307%
35	16.489	2398	2406	2416	rVB	14061779	28478838	31.82%	6.762%

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Instrument :
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ClientSampleId :
SB-03-9.5-10.0

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_Y\methods\82Y082923S.M
Title : SW846 8260

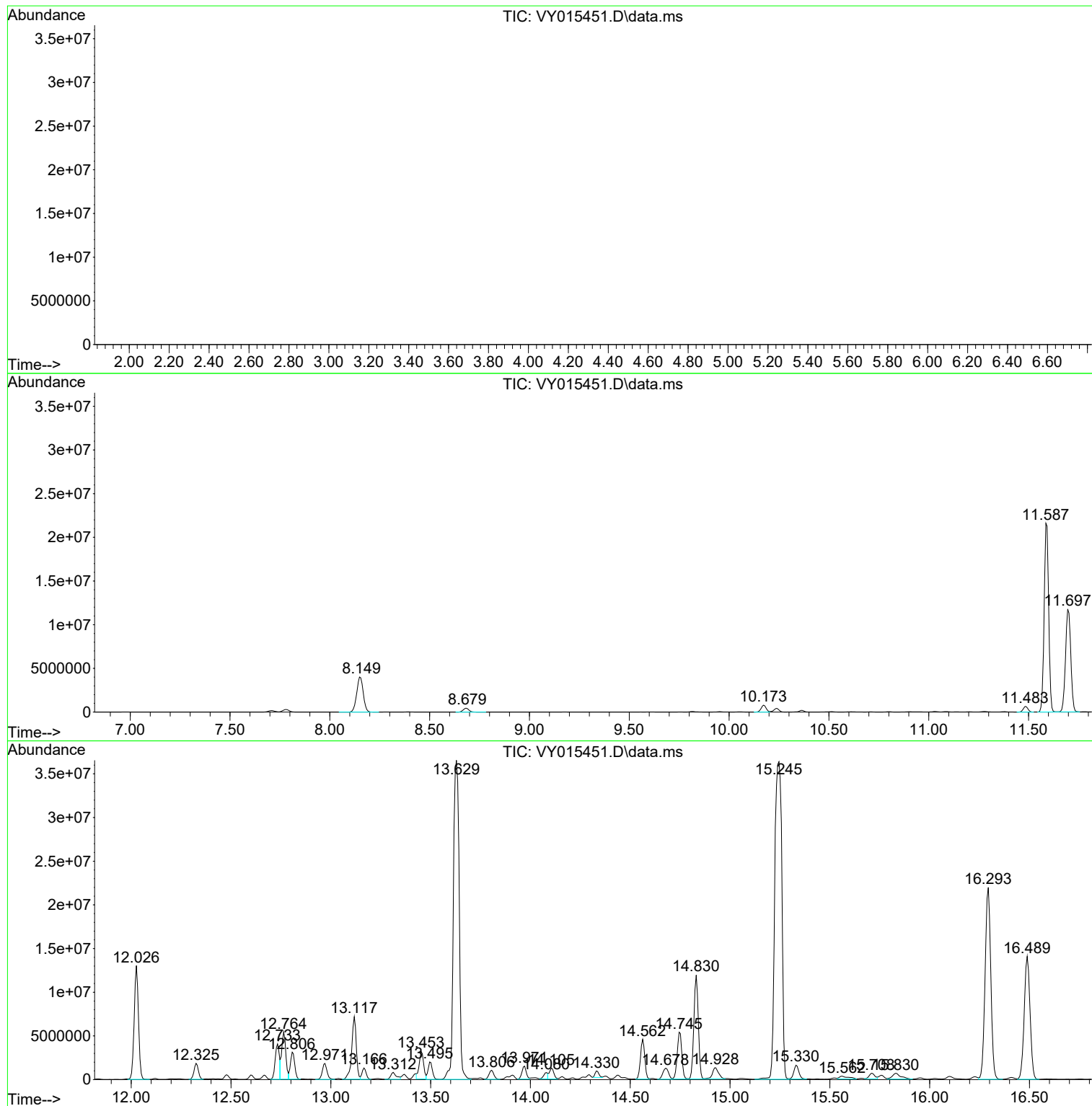
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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



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SB-03-9.5-10.0

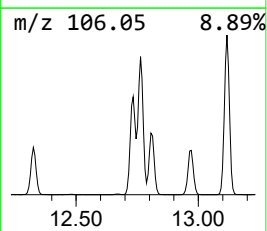
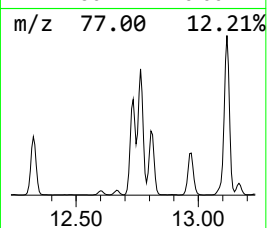
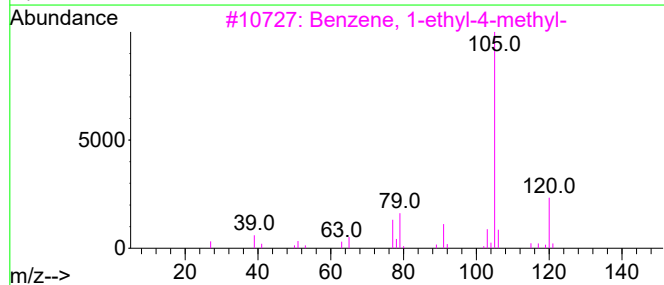
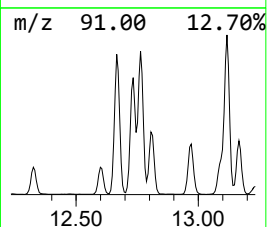
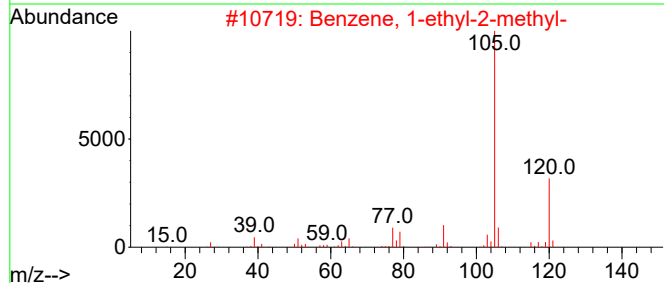
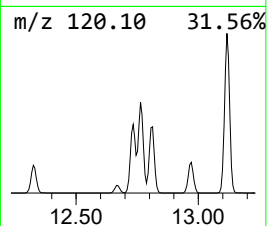
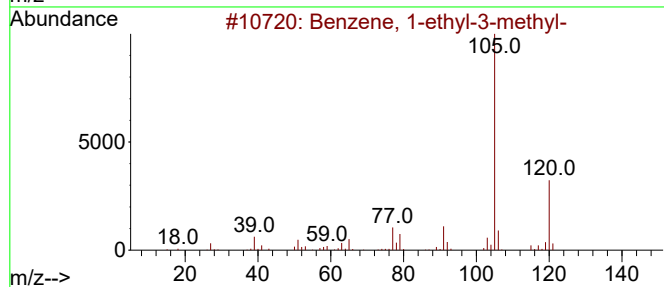
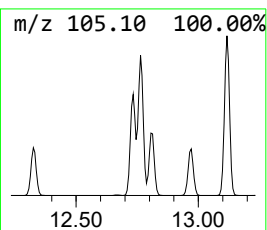
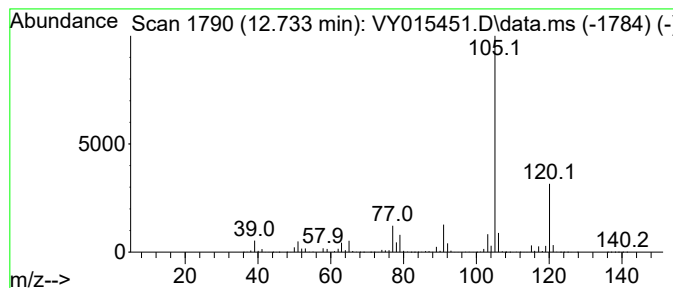
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	62.30 ug/l	6006170	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

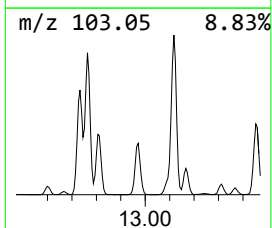
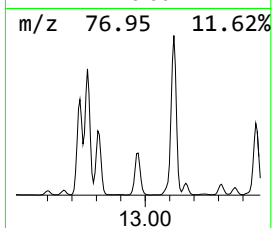
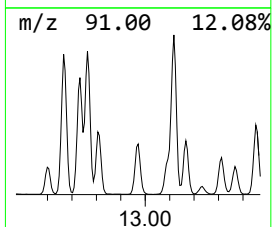
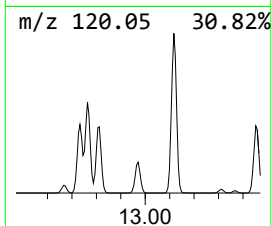
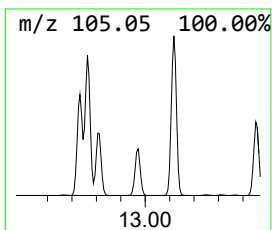
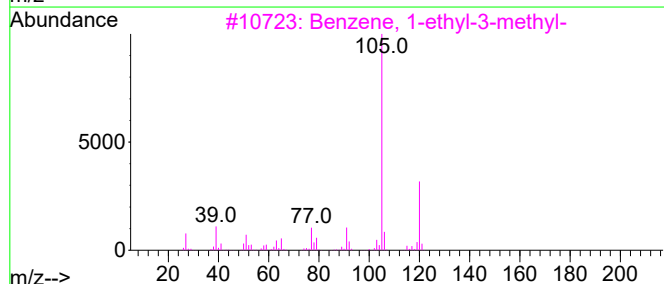
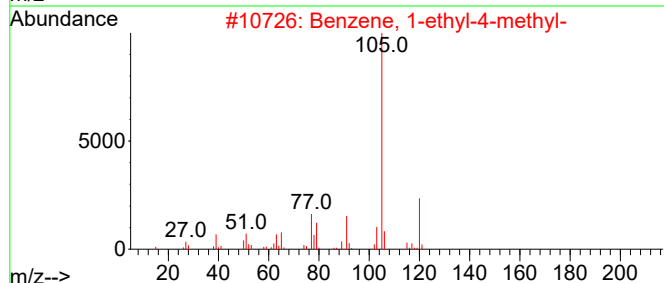
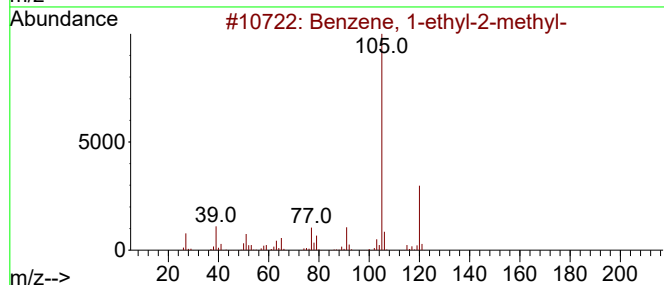
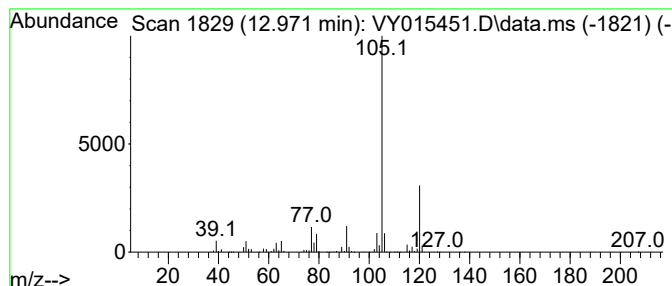
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	30.50 ug/l	2940290	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



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Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

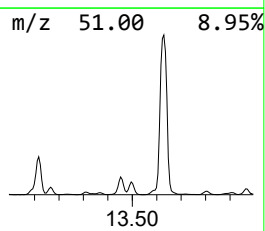
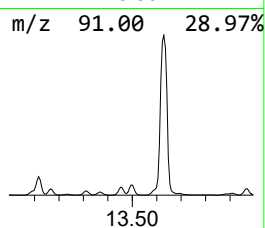
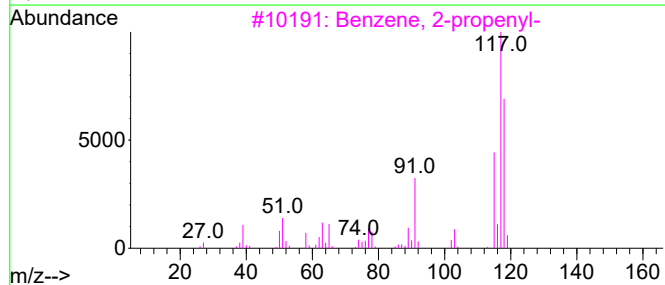
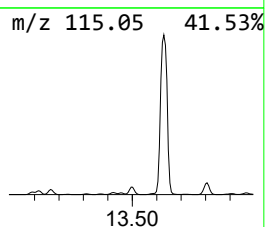
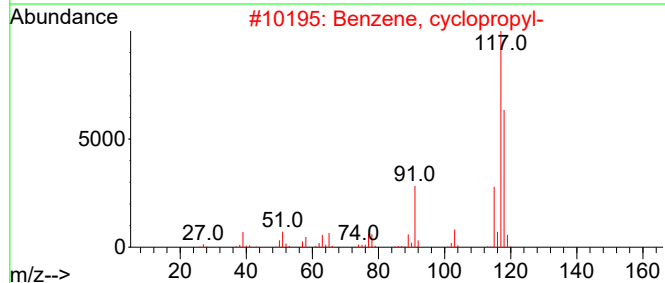
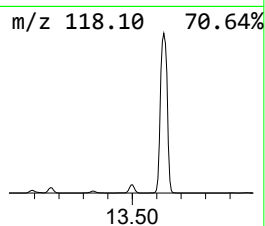
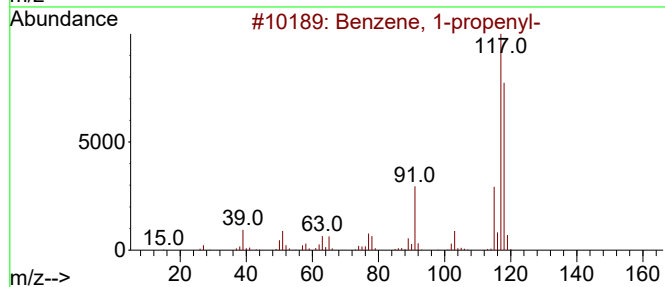
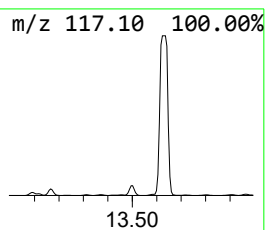
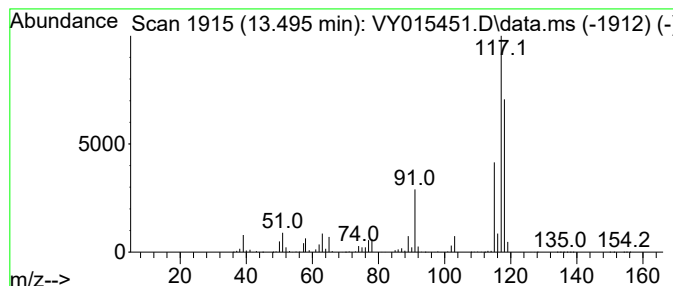
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-propenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.495	30.26 ug/l	2917720	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	95
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	93
4			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	89
5			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	80



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ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

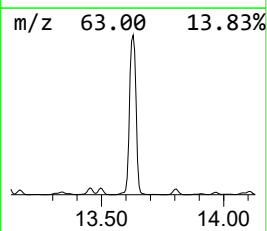
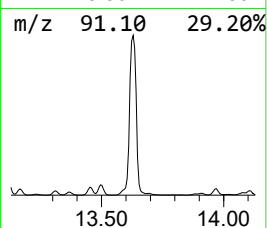
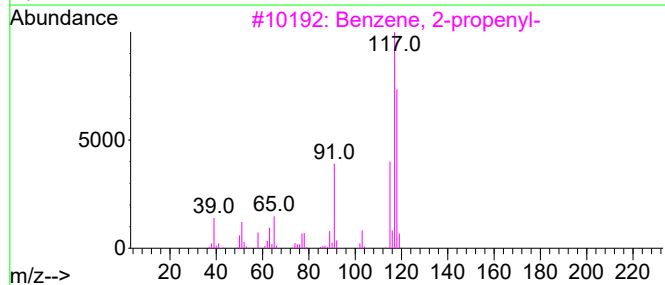
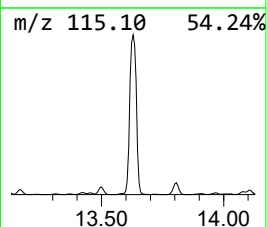
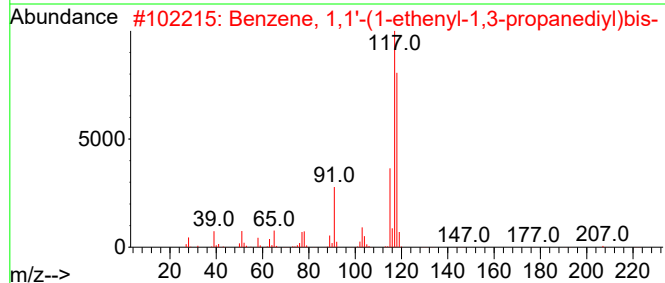
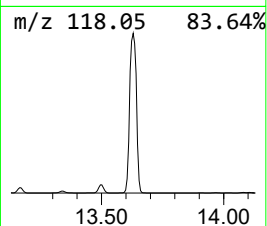
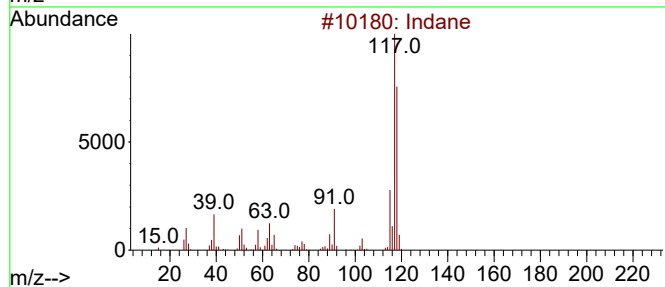
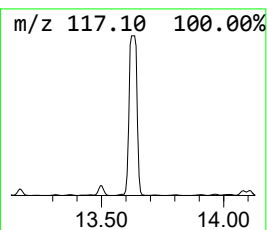
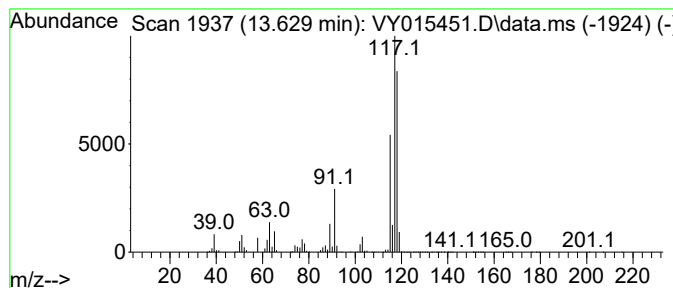
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.629	757.13 ug/l	72995500	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	90
2			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	90
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	81



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Instrument :
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ClientSampleId :
SB-03-9.5-10.0

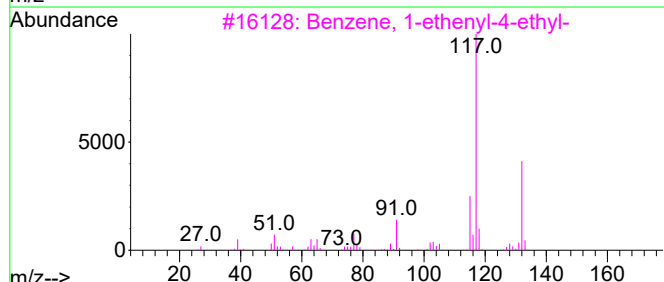
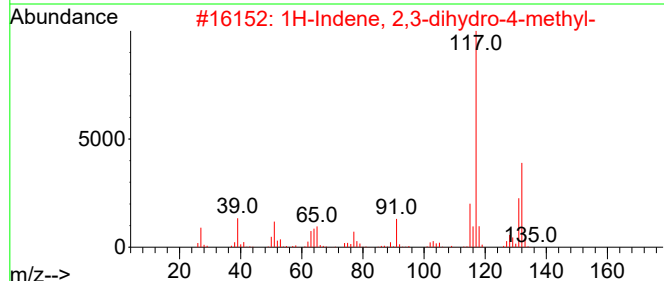
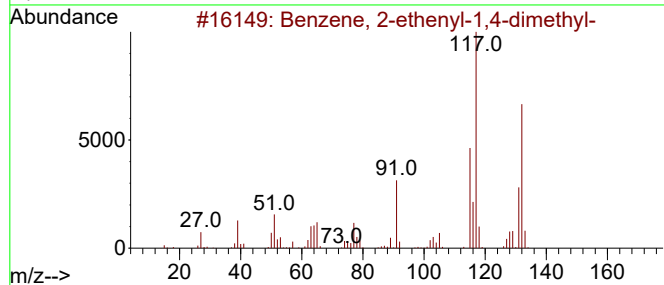
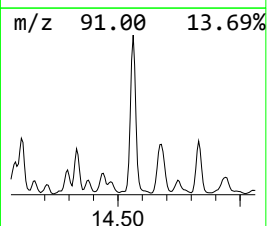
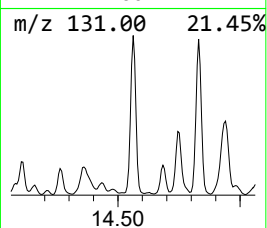
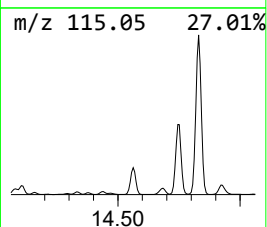
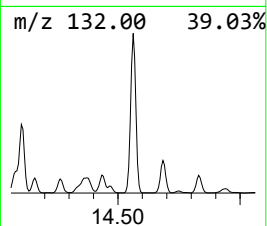
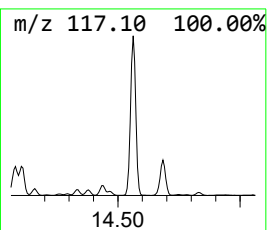
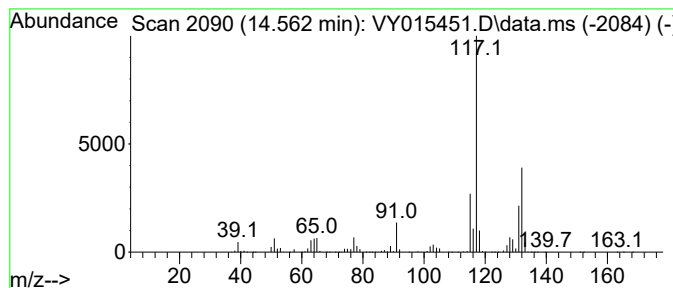
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.562	69.82 ug/l	6731340	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
Data File : VY015451.D
Acq On : 06 Sep 2023 15:44
Operator : SY/MD
Sample : 04299-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

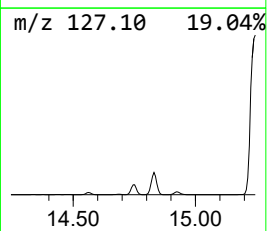
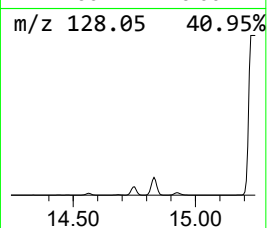
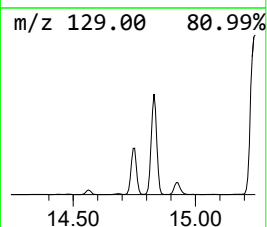
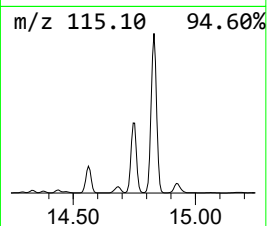
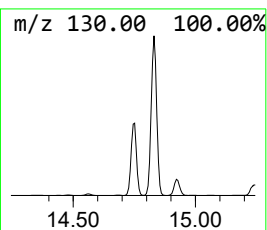
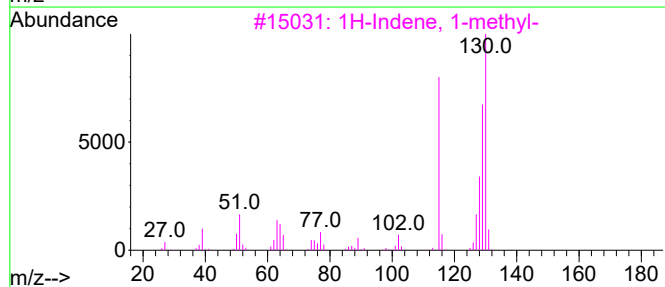
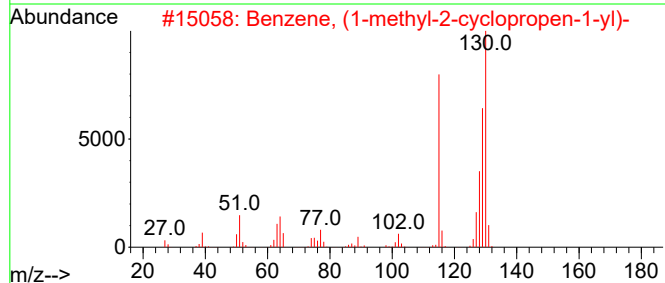
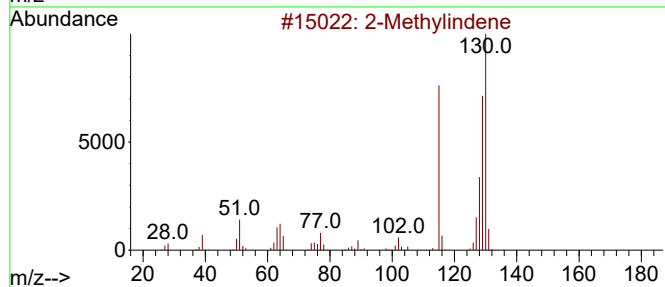
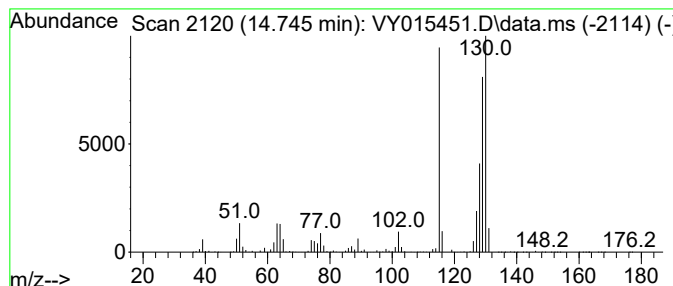
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1-methyl-2-cyclop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.745	88.69 ug/l	8550400	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene			130	C10H10	002177-47-1	97
2	Benzene, (1-methyl-2-cyclopropen...			130	C10H10	065051-83-4	97
3	1H-Indene, 1-methyl-			130	C10H10	000767-59-9	96
4	1,4-Dihydronaphthalene			130	C10H10	000612-17-9	95
5	1H-Indene, 3-methyl-			130	C10H10	000767-60-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
Data File : VY015451.D
Acq On : 06 Sep 2023 15:44
Operator : SY/MD
Sample : 04299-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

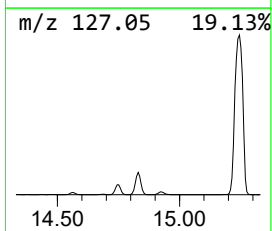
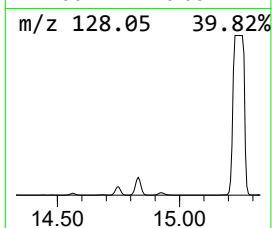
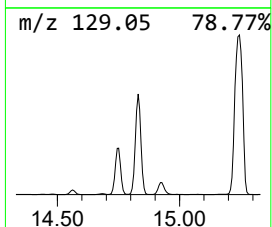
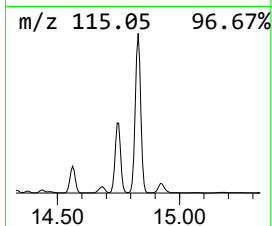
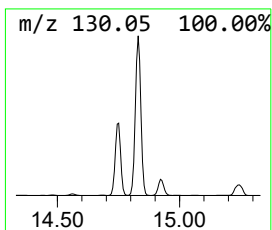
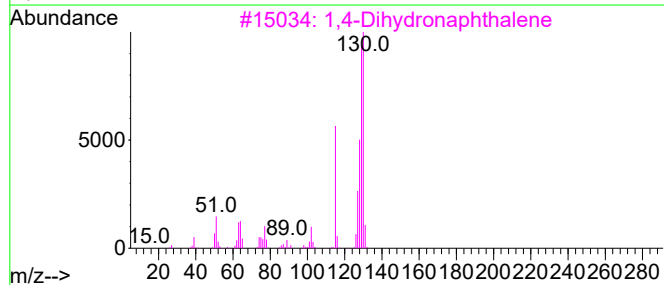
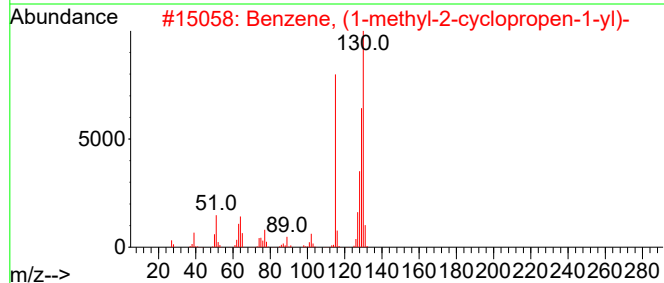
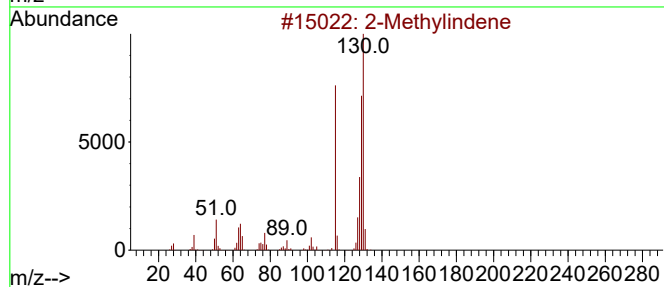
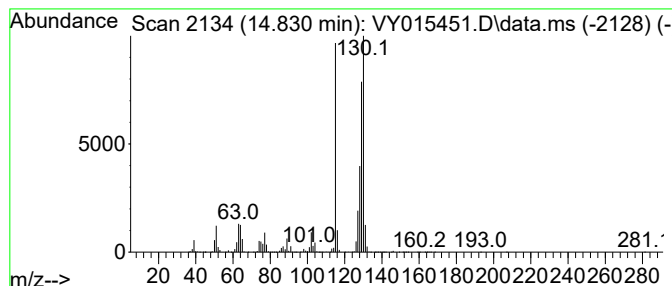
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Quant Title : SW846 8260

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

Peak Number 7 Benzene,1-methyl-1,2-propad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.830	190.22 ug/l	18339500	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene			130	C10H10	002177-47-1	97
2	Benzene, (1-methyl-2-cyclopropen...			130	C10H10	065051-83-4	96
3	1,4-Dihydronaphthalene			130	C10H10	000612-17-9	95
4	Benzene,1-methyl-1,2-propadienyl-			130	C10H10	022433-39-2	94
5	1H-Indene, 1-methyl-			130	C10H10	000767-59-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
Data File : VY015451.D
Acq On : 06 Sep 2023 15:44
Operator : SY/MD
Sample : 04299-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

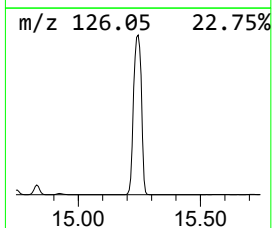
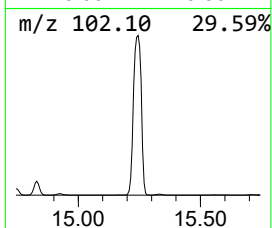
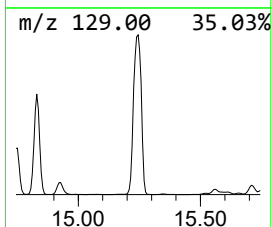
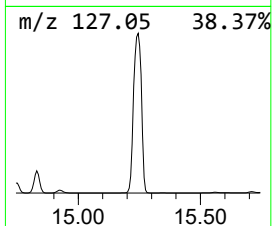
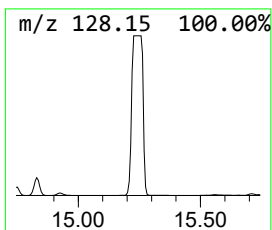
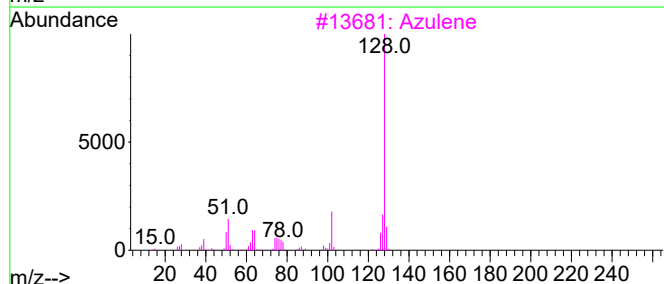
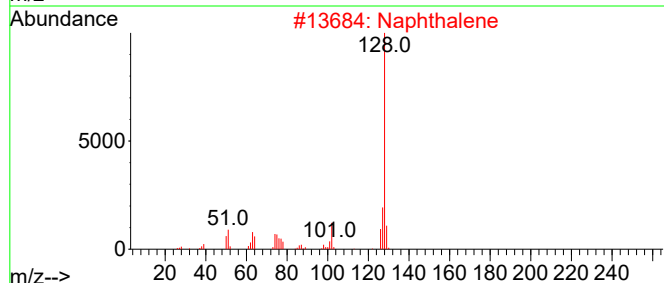
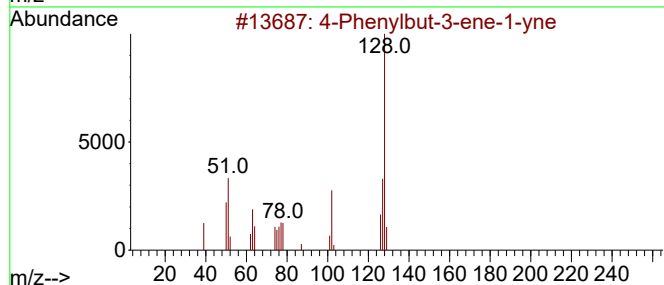
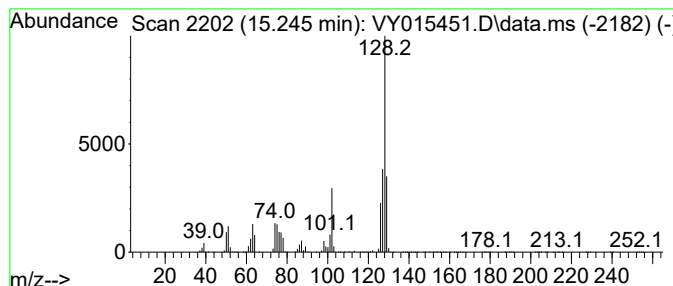
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Quant Title : SW846 8260

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

Peak Number 8 4-Phenylbut-3-ene-1-yne Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.245	928.29 ug/l	89497100	1,4-Dichlorobenzene-d4	13.422

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Phenylbut-3-ene-1-yne	128	C10H8	1000222-12-0	76
2		Naphthalene	128	C10H8	000091-20-3	70
3		Azulene	128	C10H8	000275-51-4	70
4		[4.2.2]Propella-2,4,7,9-tetraene	128	C10H8	088090-34-0	52
5		1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
Data File : VY015451.D
Acq On : 06 Sep 2023 15:44
Operator : SY/MD
Sample : 04299-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

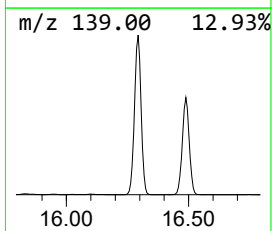
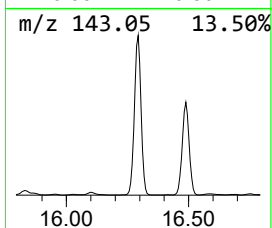
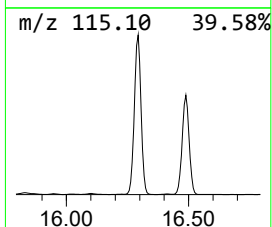
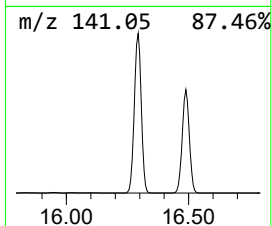
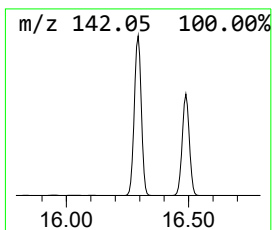
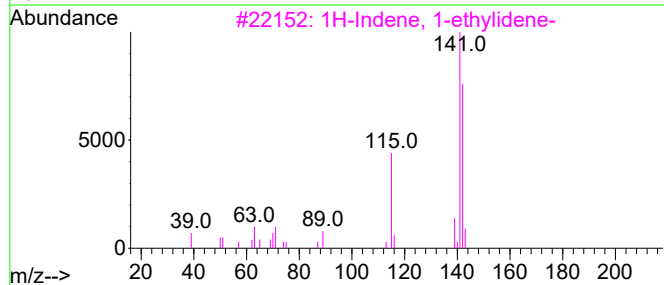
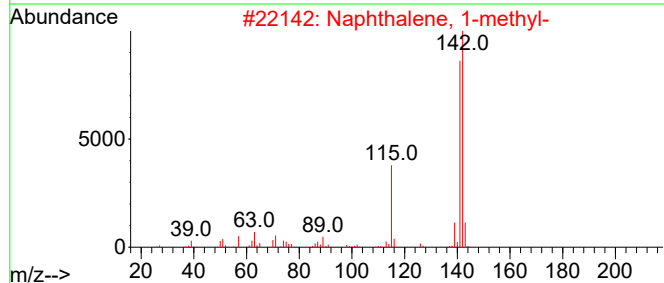
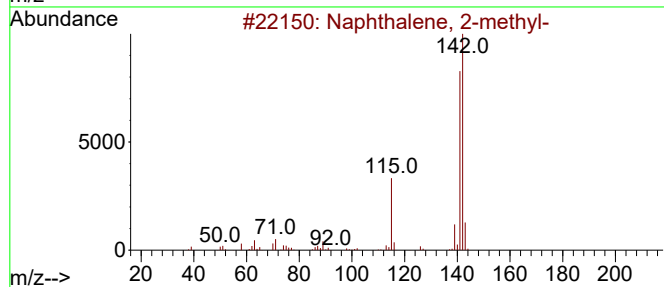
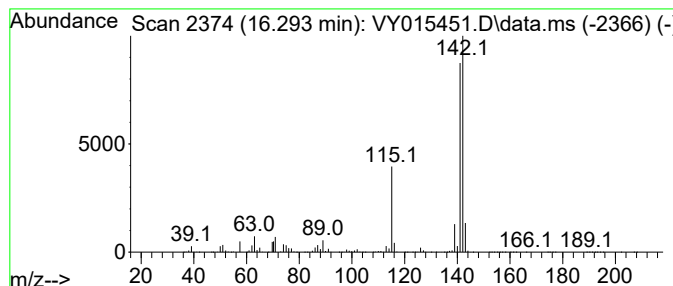
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.294	450.23 ug/l	43407300	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	94
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
Data File : VY015451.D
Acq On : 06 Sep 2023 15:44
Operator : SY/MD
Sample : 04299-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

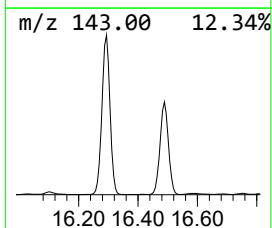
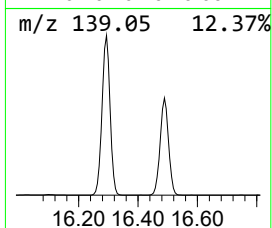
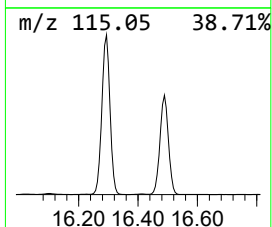
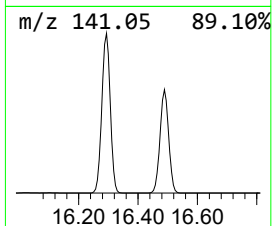
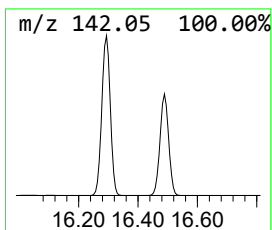
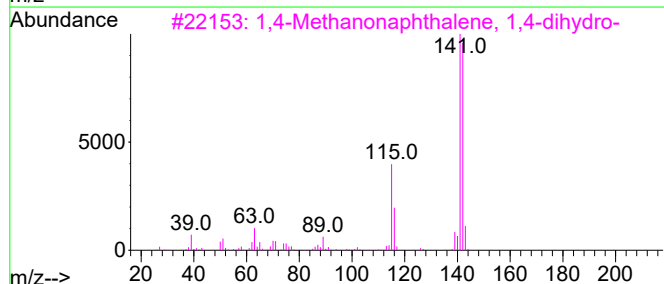
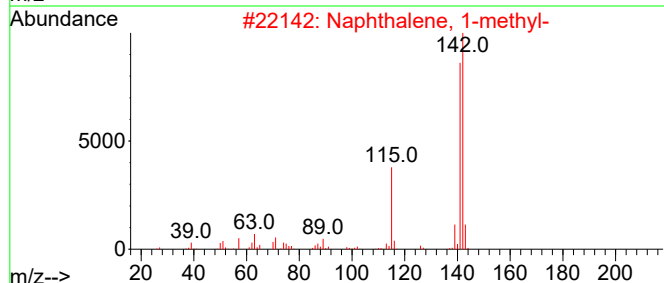
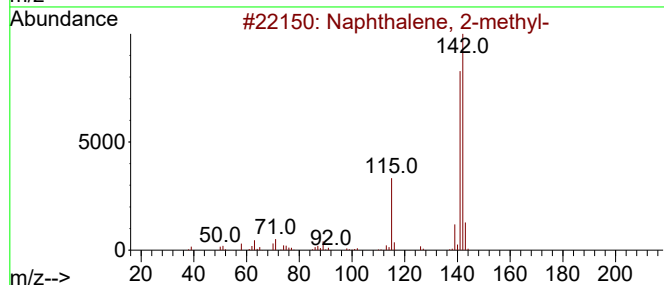
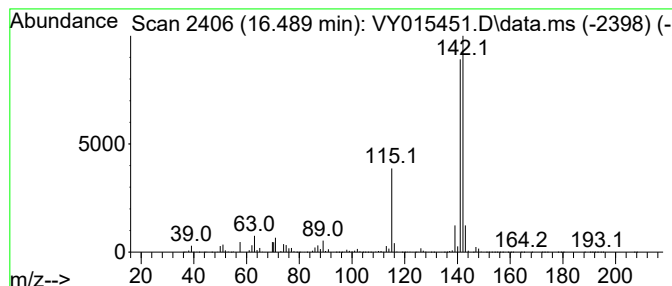
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.M
Quant Title : SW846 8260

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

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.489	295.39 ug/l	28478800	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5			Benzocycloheptatriene	142	C11H10	000264-09-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY090623\
Data File : VY015451.D
Acq On : 06 Sep 2023 15:44
Operator : SY/MD
Sample : 04299-05
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-03-9.5-10.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y082923S.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
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Benzene, 1-ethy...	12.971	30.5	ug/l	2940290	4	13.422	4820540	50.0
Benzene, 1-prop...	13.495	30.3	ug/l	2917720	4	13.422	4820540	50.0
Indane	13.629	757.1	ug/l	72995500	4	13.422	4820540	50.0
Benzene, 2-ethe...	14.562	69.8	ug/l	6731340	4	13.422	4820540	50.0
Benzene, (1-met...	14.745	88.7	ug/l	8550400	4	13.422	4820540	50.0
Benzene,1-methy...	14.830	190.2	ug/l	18339500	4	13.422	4820540	50.0
4-Phenylbut-3-e...	15.245	928.3	ug/l	89497100	4	13.422	4820540	50.0
Naphthalene, 1-...	16.294	450.2	ug/l	43407300	4	13.422	4820540	50.0
1,4-Methanonaph...	16.489	295.4	ug/l	28478800	4	13.422	4820540	50.0