

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
 Data File : VD078542.D
 Acq On : 23 Feb 2024 18:07
 Operator : RP/MD
 Sample : P1584-01
 Misc : 5.02G/5ml/MSVOA_D/SOIL/A
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 AU-06-02232024

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_D\Method\82D021324S.M
 Title : SW846 8260

Signal : TIC: VD078542.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.740	20	26	34	rBV	111155	188539	19.99%	1.625%
2	4.793	537	545	547	rBV3	5326	10450	1.11%	0.090%
3	7.805	1047	1057	1063	rBV2	87568	210733	22.34%	1.817%
4	7.875	1063	1069	1080	rVB	140280	315802	33.48%	2.722%
5	8.234	1121	1130	1140	rVB2	57055	131185	13.91%	1.131%
6	8.775	1214	1222	1230	rBV	226767	448614	47.55%	3.867%
7	10.269	1466	1476	1484	rBV	392600	676500	71.71%	5.832%
8	10.663	1538	1543	1548	rBV2	7455	11824	1.25%	0.102%
9	11.028	1598	1605	1611	rBV	7480	12136	1.29%	0.105%
10	11.581	1691	1699	1709	rVB	427613	716776	75.98%	6.179%
11	11.793	1727	1735	1741	rBV2	24473	49885	5.29%	0.430%
12	12.122	1786	1791	1796	rVB2	15810	26646	2.82%	0.230%
13	12.575	1861	1868	1879	rVB	352012	567660	60.17%	4.894%
14	12.763	1892	1900	1904	rBV2	21369	36919	3.91%	0.318%
15	12.828	1904	1911	1914	rBV2	77570	129252	13.70%	1.114%
16	12.857	1914	1916	1919	rVV2	34968	51373	5.45%	0.443%
17	12.898	1919	1923	1929	rVV2	94214	153227	16.24%	1.321%
18	12.945	1929	1931	1938	rVB4	7763	11958	1.27%	0.103%
19	13.063	1942	1951	1957	rBV	45433	87000	9.22%	0.750%
20	13.128	1958	1962	1969	rVB5	19404	36933	3.91%	0.318%
21	13.210	1969	1976	1982	rBV	197144	318198	33.73%	2.743%
22	13.345	1992	1999	2003	rVB3	19507	34776	3.69%	0.300%
23	13.404	2003	2009	2014	rBV5	39881	75358	7.99%	0.650%
24	13.457	2014	2018	2022	rVV3	18944	31529	3.34%	0.272%
25	13.516	2022	2028	2038	rVV2	512801	943391	100.00%	8.133%
26	13.592	2038	2041	2046	rVB5	9784	12560	1.33%	0.108%
27	13.716	2057	2062	2066	rVV2	169964	295633	31.34%	2.549%
28	13.757	2066	2069	2078	rVV4	116636	247943	26.28%	2.137%
29	13.839	2078	2083	2088	rVV3	124976	192047	20.36%	1.656%
30	13.916	2088	2096	2101	rVV	57927	105061	11.14%	0.906%
31	13.975	2101	2106	2108	rVV	63590	97671	10.35%	0.842%
32	14.004	2108	2111	2116	rVV	86497	131276	13.92%	1.132%
33	14.063	2116	2121	2126	rVV	115847	170237	18.05%	1.468%
34	14.122	2126	2131	2134	rVV5	23713	40411	4.28%	0.348%
35	14.169	2134	2139	2145	rVB	133449	219599	23.28%	1.893%
36	14.245	2145	2152	2157	rBV6	21774	54518	5.78%	0.470%

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Integrator: RTE

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Sampling : 1

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Max Peaks: 100

Peak Location: TOP

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Peak separation: 5

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37	14.304	2157	2162	2166	rVV3	34910	72501	7.69%	0.625%
38	14.351	2166	2170	2173	rVV5	43305	88118	9.34%	0.760%
39	14.387	2173	2176	2179	rVV	64508	104732	11.10%	0.903%
40	14.422	2179	2182	2187	rVV	91471	135946	14.41%	1.172%
41	14.469	2187	2190	2194	rVV	73820	92977	9.86%	0.802%
42	14.510	2194	2197	2204	rVB5	47175	69502	7.37%	0.599%
43	14.575	2204	2208	2209	rBV4	21335	26742	2.83%	0.231%
44	14.610	2209	2214	2217	rVV3	70971	109965	11.66%	0.948%
45	14.657	2217	2222	2226	rVV2	145221	290156	30.76%	2.501%
46	14.698	2226	2229	2234	rVB2	138921	204219	21.65%	1.761%
47	14.775	2234	2242	2247	rBV5	269870	589270	62.46%	5.080%
48	14.822	2247	2250	2256	rVB3	53077	98890	10.48%	0.852%
49	14.928	2262	2268	2275	rVB	187744	310259	32.89%	2.675%
50	14.998	2275	2280	2281	rBV3	37350	47249	5.01%	0.407%
51	15.034	2281	2286	2290	rVV4	103857	212231	22.50%	1.830%
52	15.075	2290	2293	2298	rVV2	60946	101782	10.79%	0.877%
53	15.151	2298	2306	2315	rVB3	107218	256085	27.15%	2.208%
54	15.216	2315	2317	2320	rBV4	8383	11446	1.21%	0.099%
55	15.304	2326	2332	2334	rBV5	37393	65718	6.97%	0.567%
56	15.392	2341	2347	2351	rBV	94618	160255	16.99%	1.382%
57	15.445	2351	2356	2363	rVV5	77906	229839	24.36%	1.981%
58	15.528	2367	2370	2379	rVB3	62244	116683	12.37%	1.006%
59	15.622	2379	2386	2395	rBV4	72592	167579	17.76%	1.445%
60	15.698	2395	2399	2405	rVB6	15809	22539	2.39%	0.194%
61	15.786	2407	2414	2416	rBV7	43149	81445	8.63%	0.702%
62	15.828	2416	2421	2430	rVB2	181184	360527	38.22%	3.108%
63	15.916	2432	2436	2442	rBV6	19638	34481	3.66%	0.297%
64	15.998	2442	2450	2462	rVB5	46858	151748	16.09%	1.308%
65	16.145	2466	2475	2481	rBV3	79786	185536	19.67%	1.599%
66	16.339	2498	2508	2514	rVV6	72041	163596	17.34%	1.410%
67	16.404	2514	2519	2527	rVV6	40693	83188	8.82%	0.717%
68	16.610	2546	2554	2559	rBV8	22989	68671	7.28%	0.592%
69	16.739	2568	2576	2584	rBV7	16728	42553	4.51%	0.367%

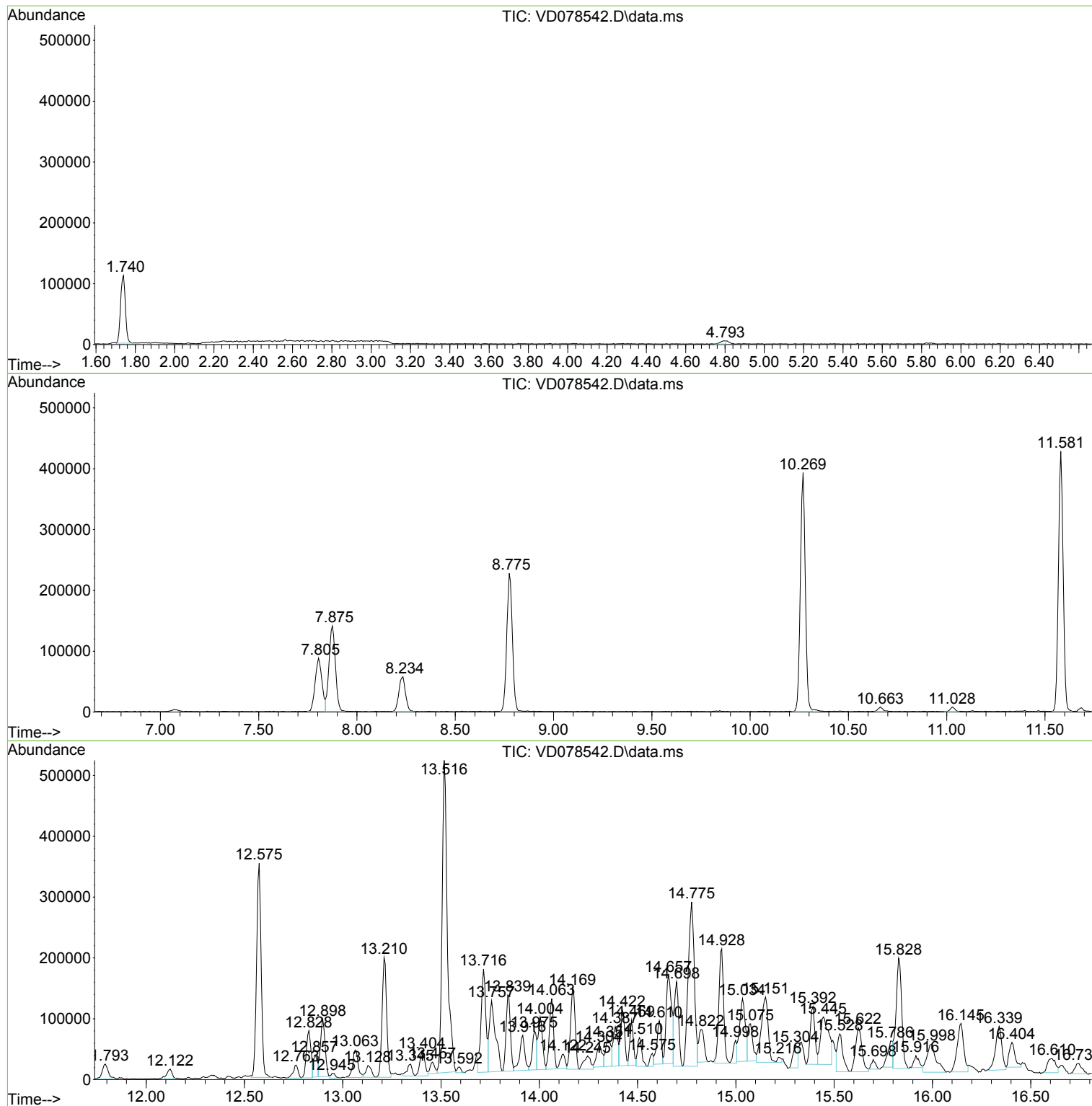
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ClientSampleId :
AU-06-02232024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.M
Quant Title : SW846 8260

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



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ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

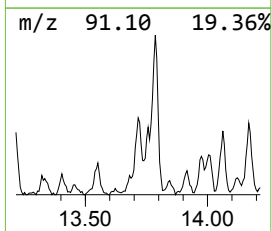
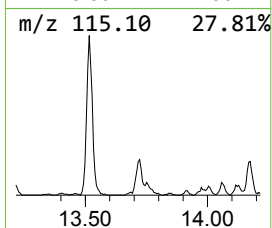
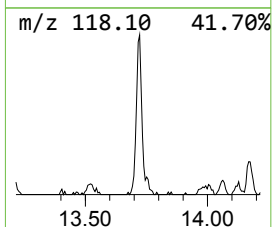
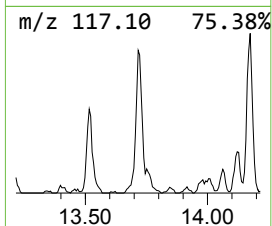
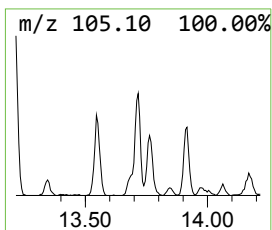
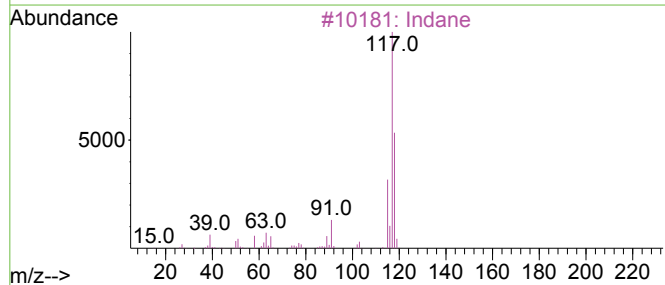
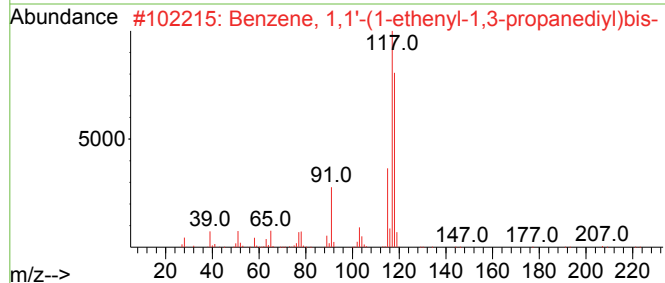
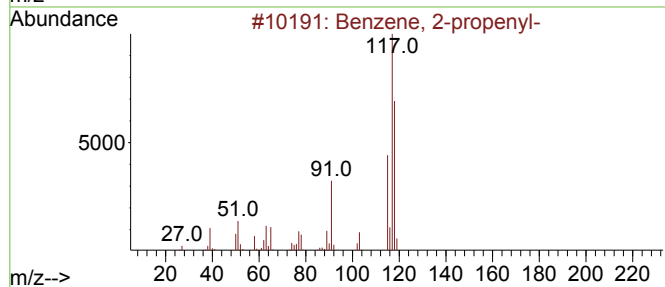
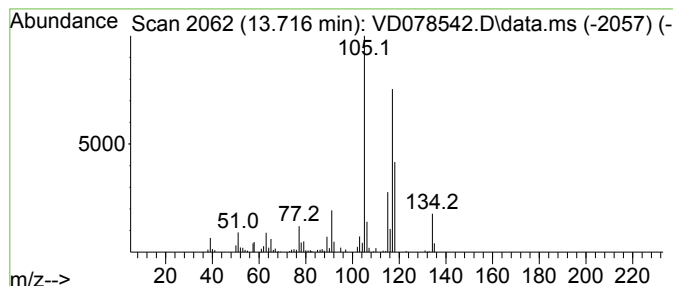
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 2-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.716	15.67 ug/l	295633	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
2			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	64
3			Indane	118	C9H10	000496-11-7	64
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	58
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	53



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ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

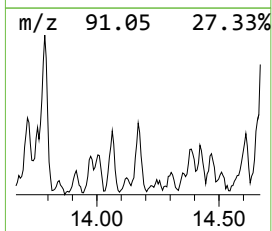
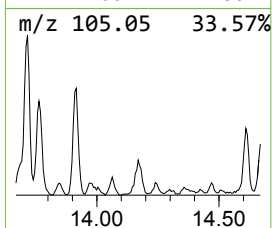
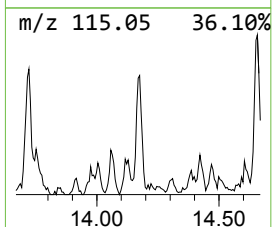
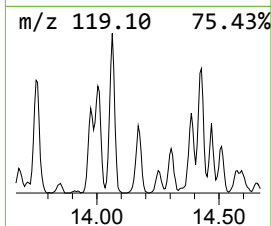
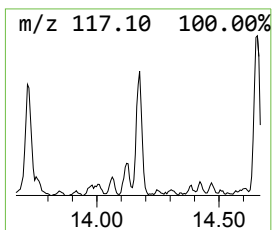
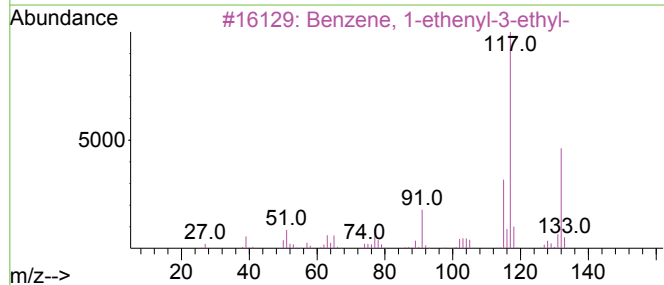
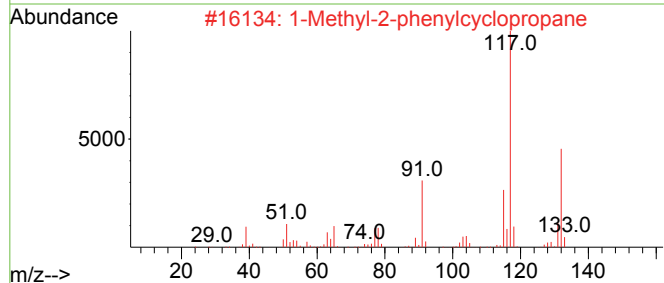
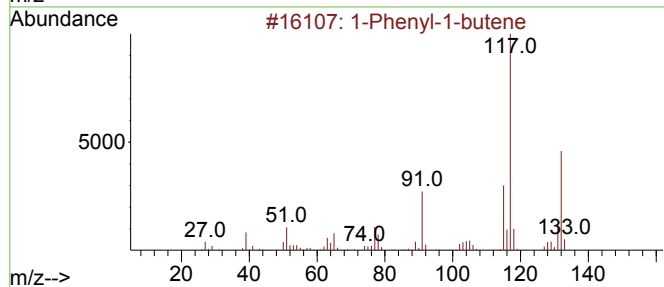
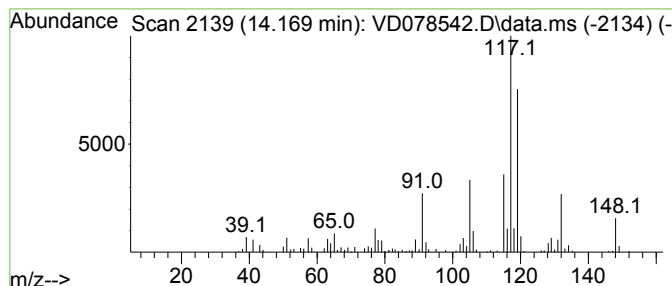
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.M
Quant Title : SW846 8260

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

Peak Number 3 1-Phenyl-1-butene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.169	11.64 ug/l	219599	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	55
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	55
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	49
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	46
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	46



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Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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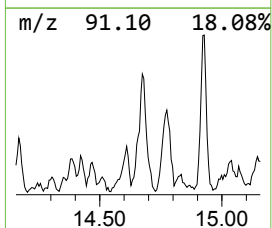
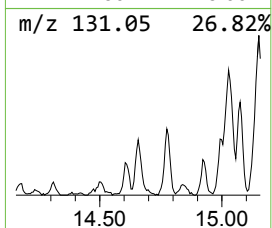
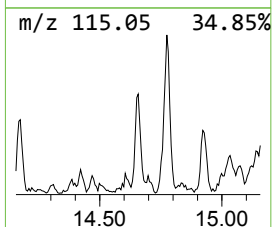
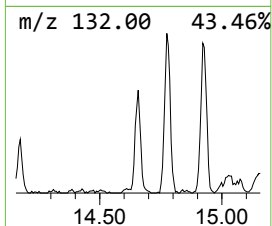
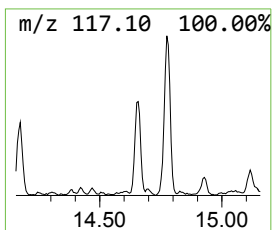
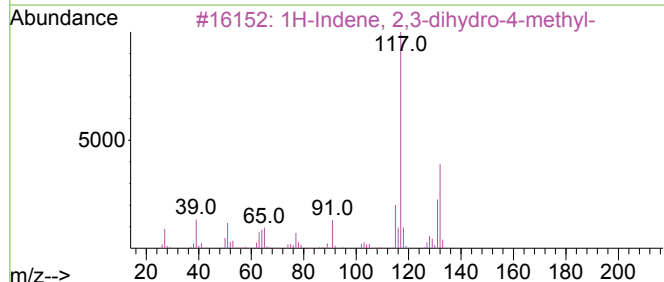
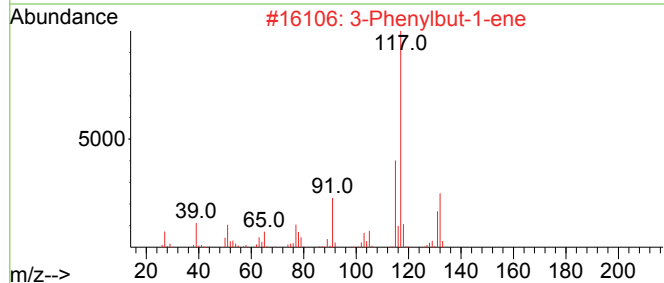
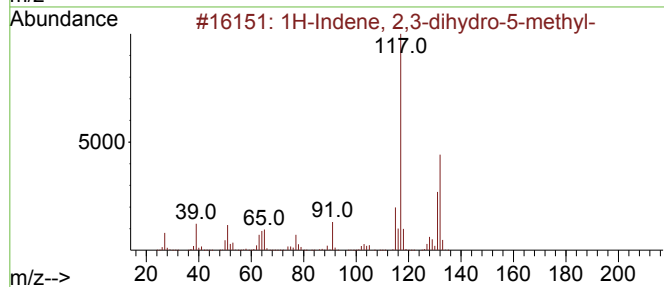
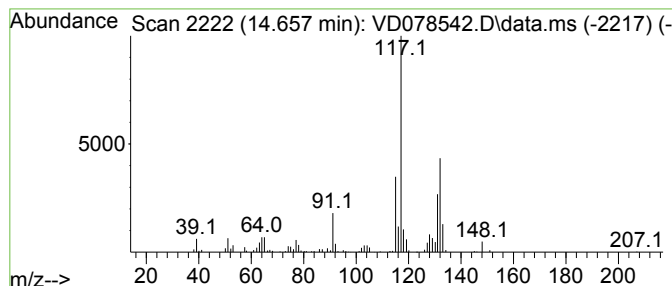
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.M
Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.657	15.38 ug/l	290156	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	86
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



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Instrument :
MSVOA_D
ClientSampleId :
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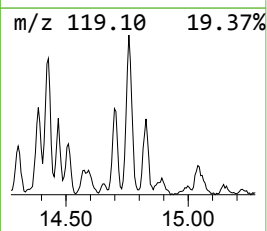
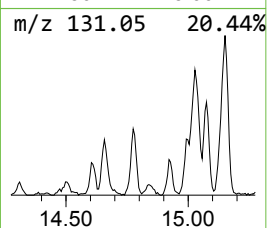
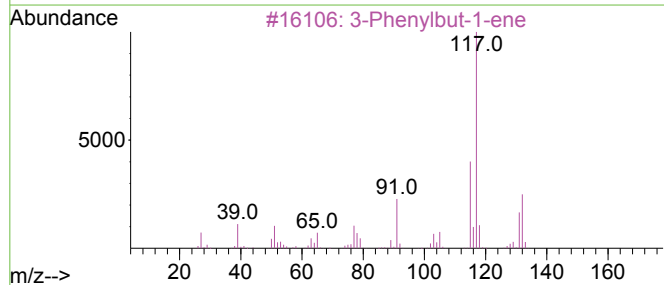
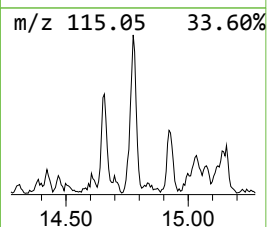
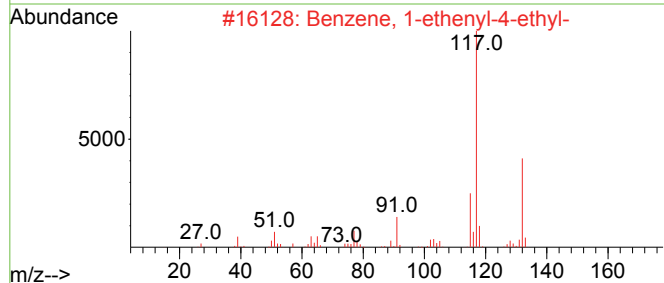
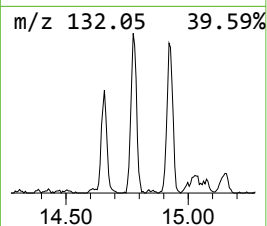
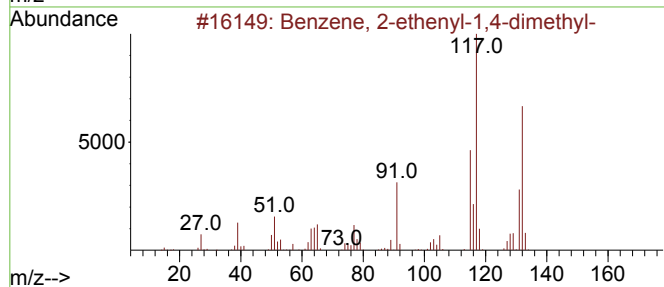
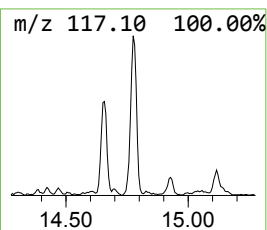
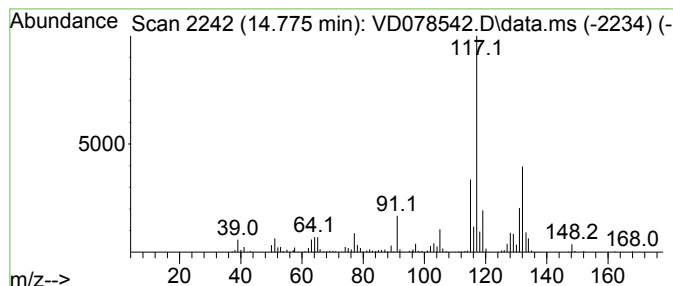
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.775	31.23 ug/l	589270	1,4-Dichlorobenzene-d4	13.516

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-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	81
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
Data File : VD078542.D
Acq On : 23 Feb 2024 18:07
Operator : RP/MD
Sample : P1584-01
Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

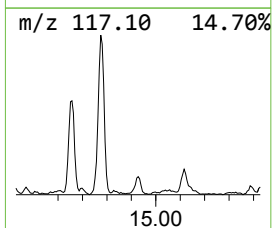
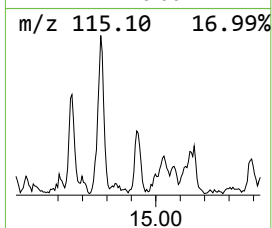
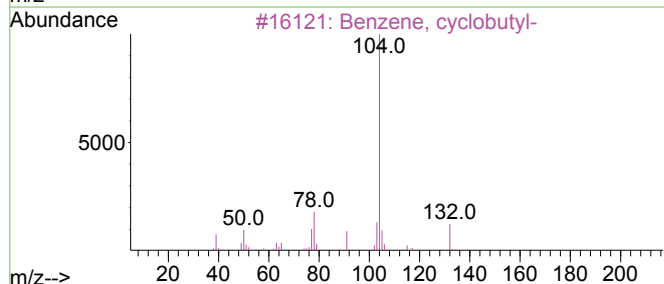
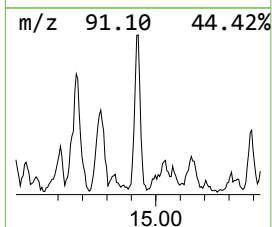
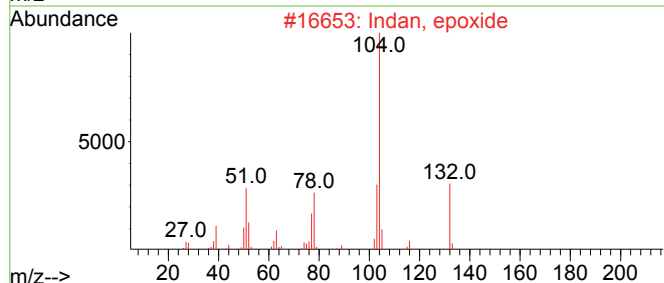
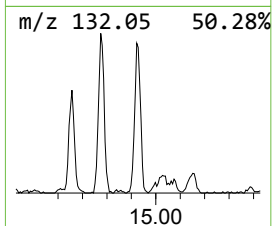
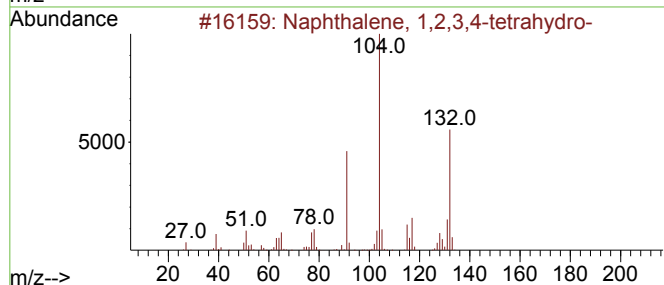
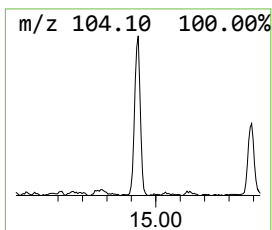
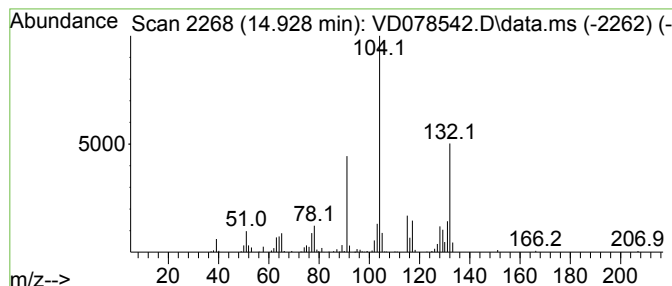
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.928	16.44 ug/l	310259	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Indan, epoxide	132	C9H8O	000768-22-9	64
3			Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
4			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	53
5			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
Data File : VD078542.D
Acq On : 23 Feb 2024 18:07
Operator : RP/MD
Sample : P1584-01
Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

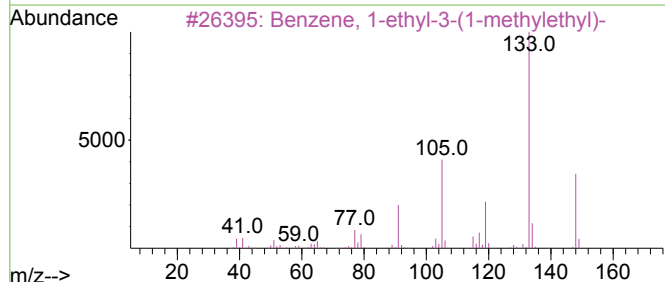
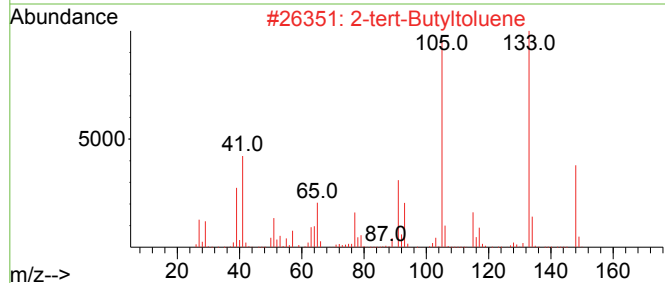
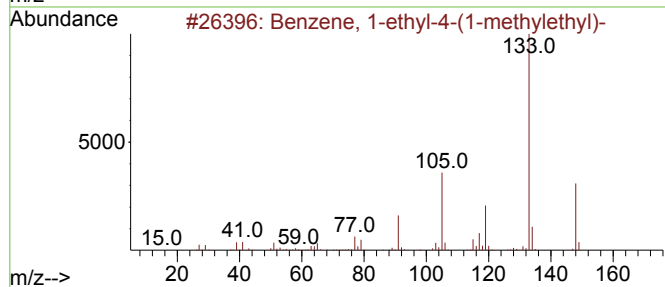
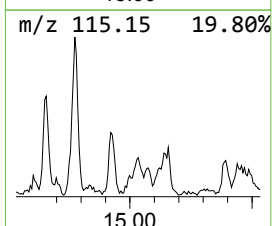
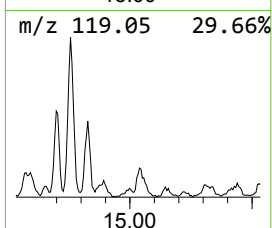
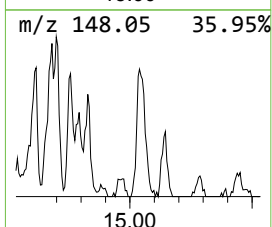
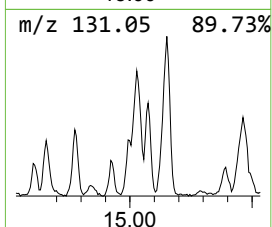
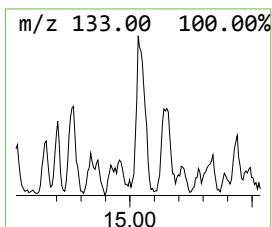
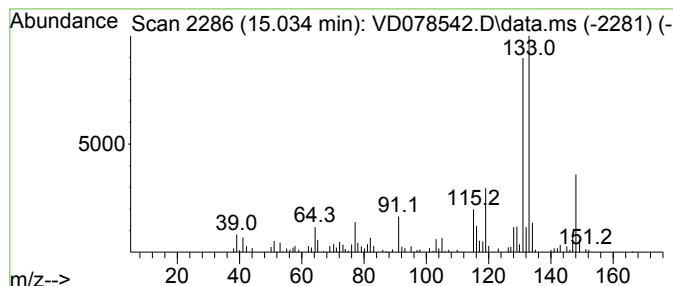
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Quant Title : SW846 8260

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

Peak Number 7 unknown15.034 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.034	11.25 ug/l	212231	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	45
2			2-tert-Butyltoluene	148	C11H16	001074-92-6	43
3			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	43
4			3,4-Dimethylcumene	148	C11H16	1000370-34-1	38
5			Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
Data File : VD078542.D
Acq On : 23 Feb 2024 18:07
Operator : RP/MD
Sample : P1584-01
Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

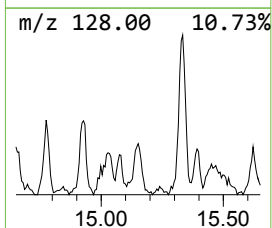
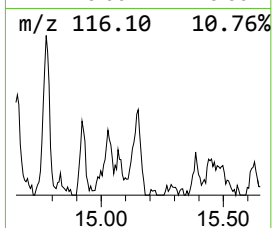
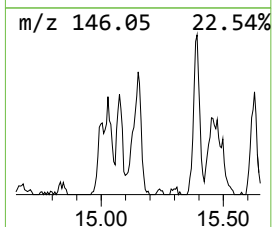
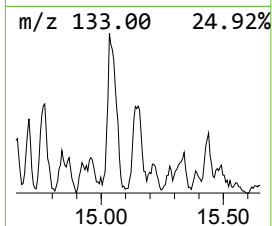
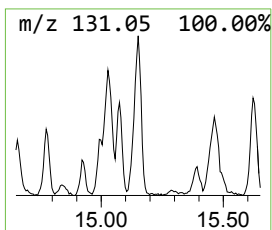
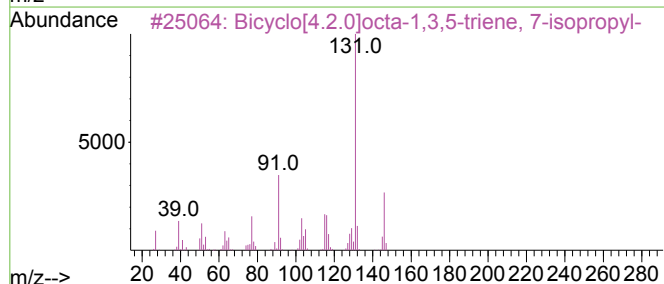
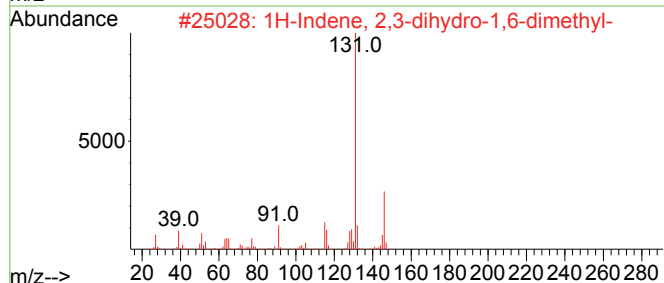
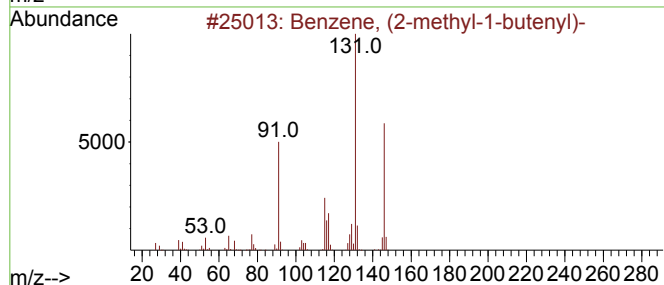
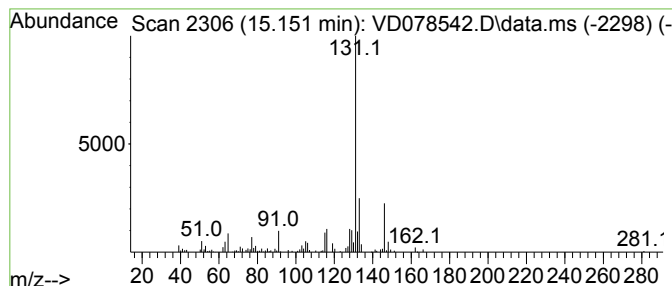
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.151	13.57 ug/l	256085	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	81
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
3			Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	81
4			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
Data File : VD078542.D
Acq On : 23 Feb 2024 18:07
Operator : RP/MD
Sample : P1584-01
Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

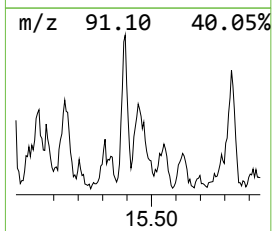
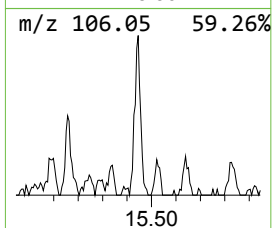
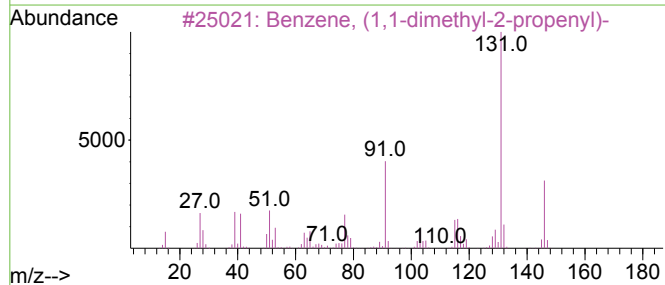
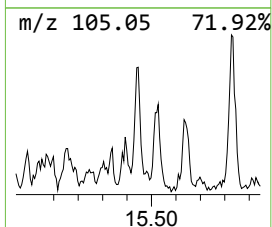
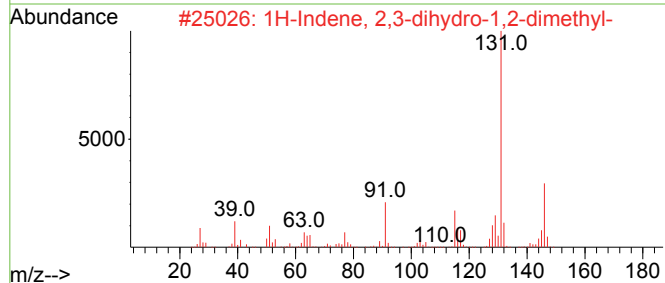
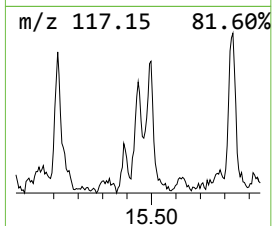
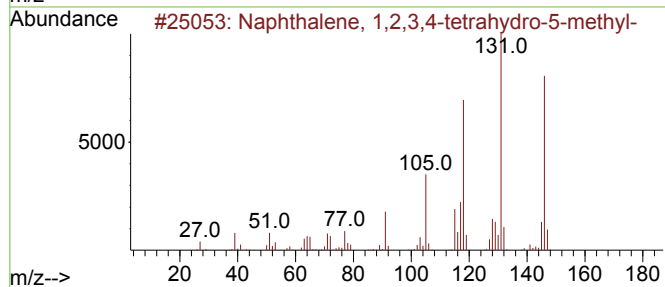
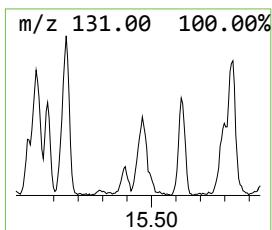
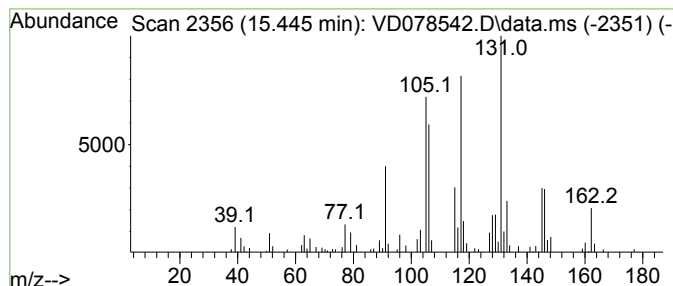
Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.445	12.18 ug/l	229839	1,4-Dichlorobenzene-d4		13.516	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	002809-64-5	50
2	1H-Indene, 2,3-dihydro-1,2-dimet...		146	C11H14	017057-82-8	46
3	Benzene, (1,1-dimethyl-2-propenyl)-		146	C11H14	018321-36-3	42
4	Benzene, (1-methyl-1-butenyl)-		146	C11H14	053172-84-2	41
5	2,2-Dimethylindene, 2,3-dihydro-		146	C11H14	020836-11-7	41



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
Data File : VD078542.D
Acq On : 23 Feb 2024 18:07
Operator : RP/MD
Sample : P1584-01
Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

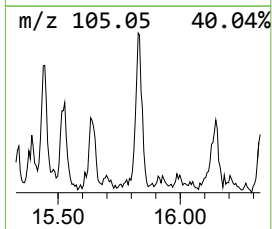
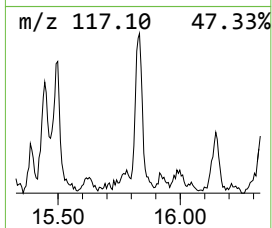
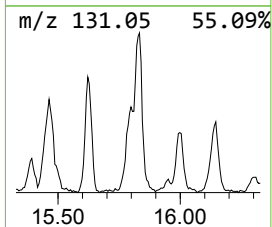
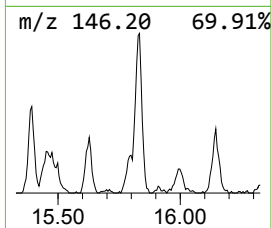
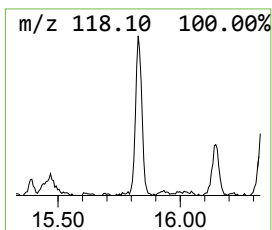
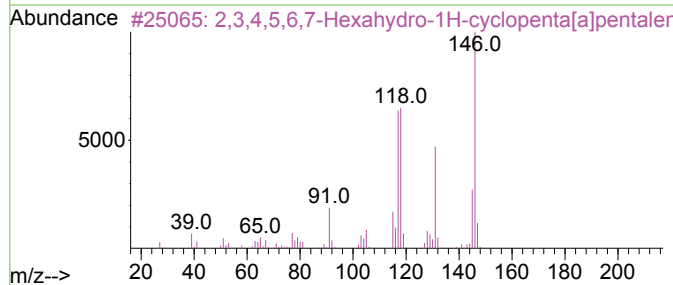
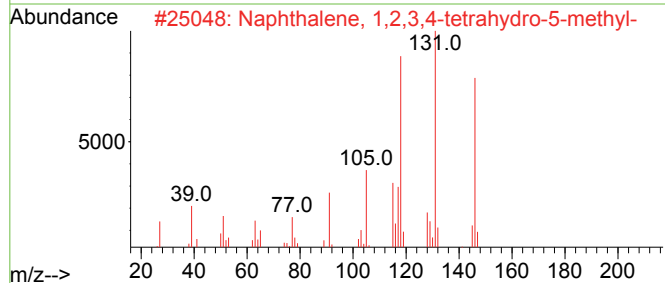
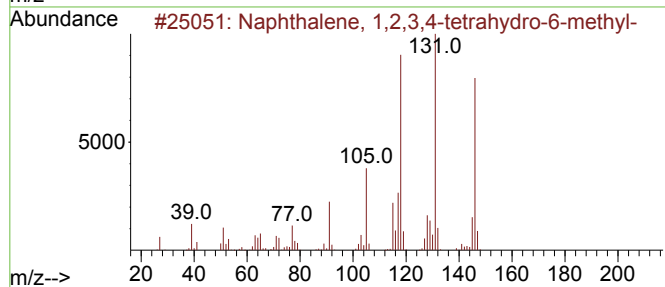
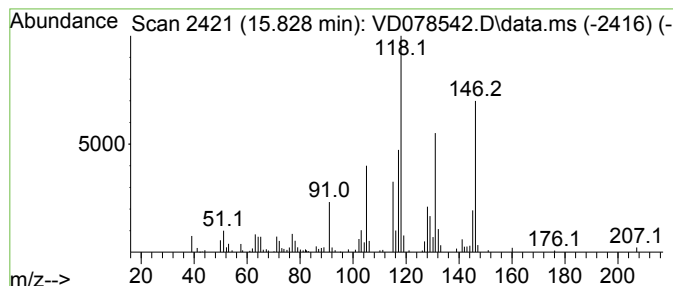
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.828	19.11 ug/l	360527	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	86
3			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	74
4			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	55
5			1H-Inden-1-one, 2,3-dihydro-3-me...	146	C10H10O	006072-57-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022324\
Data File : VD078542.D
Acq On : 23 Feb 2024 18:07
Operator : RP/MD
Sample : P1584-01
Misc : 5.02G/5ml/MSVOA_D/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
AU-06-02232024

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021324S.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
Benzene, 2-prop...	13.716	15.7	ug/l	295633	4	13.516	943391	50.0
1-Phenyl-1-butene	14.169	11.6	ug/l	219599	4	13.516	943391	50.0
1H-Indene, 2,3-...	14.657	15.4	ug/l	290156	4	13.516	943391	50.0
Benzene, 2-ethe...	14.775	31.2	ug/l	589270	4	13.516	943391	50.0
Naphthalene, 1,...	14.928	16.4	ug/l	310259	4	13.516	943391	50.0
unknown15.034	15.034	11.3	ug/l	212231	4	13.516	943391	50.0
Benzene, (2-met...	15.151	13.6	ug/l	256085	4	13.516	943391	50.0
Naphthalene, 1,...	15.445	12.2	ug/l	229839	4	13.516	943391	50.0
Naphthalene, 1,...	15.828	19.1	ug/l	360527	4	13.516	943391	50.0