

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
 Data File : VX028648.D
 Acq On : 12 May 2022 19:12
 Operator : JC/MD
 Sample : N2799-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2A

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_X\Method\82X051222W.M
 Title : SW846 8260

Signal : TIC: VX028648.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.746	103	108	121	rVB	2058547	2865418	2.94%	0.684%
2	1.953	137	142	154	rVB	1416401	1988051	2.04%	0.475%
3	2.825	280	285	289	rVV	1367464	2807521	2.88%	0.670%
4	2.867	289	292	316	rVB2	708333	1611023	1.65%	0.385%
5	3.099	319	330	348	rVB	1014419	2307684	2.36%	0.551%
6	3.447	377	387	401	rBV	698133	1546986	1.59%	0.369%
7	4.306	518	528	540	rVB	1359569	3848782	3.94%	0.919%
8	5.477	712	720	729	rBV3	544399	1897909	1.94%	0.453%
9	6.202	822	839	843	rVV3	271078	1190290	1.22%	0.284%
10	6.763	924	931	938	rBV	425601	1055177	1.08%	0.252%
11	7.379	1020	1032	1044	rBV4	757008	2020493	2.07%	0.482%
12	8.653	1235	1241	1246	rVB	905245	1362525	1.40%	0.325%
13	8.726	1246	1253	1260	rBV	14788598	23439836	24.02%	5.597%
14	10.055	1461	1471	1476	rBV	991685	1492723	1.53%	0.356%
15	10.202	1489	1495	1501	rVB	22966223	31447228	32.23%	7.510%
16	10.323	1506	1515	1520	rBV2	51891533	97581987	100.00%	23.303%
17	10.653	1562	1569	1574	rVB	30457028	41604989	42.64%	9.935%
18	10.964	1613	1620	1627	rVB	2067682	2572623	2.64%	0.614%
19	11.086	1635	1640	1645	rVB	872875	1095904	1.12%	0.262%
20	11.305	1668	1676	1683	rBV	5024631	6367442	6.53%	1.521%
21	11.384	1683	1689	1697	rVV	20864024	39361288	40.34%	9.400%
22	11.457	1697	1701	1706	rVB	12258876	14259648	14.61%	3.405%
23	11.616	1721	1727	1737	rBV	11378471	13875099	14.22%	3.313%
24	11.762	1744	1751	1756	rBV	34615204	45685073	46.82%	10.910%
25	12.024	1789	1794	1799	rVB2	1255001	1700904	1.74%	0.406%
26	12.091	1799	1805	1810	rBV	11024634	13028690	13.35%	3.111%
27	12.244	1825	1830	1838	rBV2	9697049	14218575	14.57%	3.395%
28	12.317	1838	1842	1850	rVB2	3273688	4712589	4.83%	1.125%
29	12.415	1852	1858	1862	rBV2	839394	1144920	1.17%	0.273%
30	12.463	1862	1866	1872	rVB	834949	997880	1.02%	0.238%
31	12.530	1872	1877	1879	rBV	1921934	2342256	2.40%	0.559%
32	12.555	1879	1881	1886	rVB	1617590	1885274	1.93%	0.450%
33	12.616	1886	1891	1894	rBV	3669917	4250449	4.36%	1.015%
34	12.701	1900	1905	1913	rBV2	2032044	2725785	2.79%	0.651%
35	12.921	1934	1941	1945	rBV	1645183	2105764	2.16%	0.503%
36	12.963	1945	1948	1954	rVB	2413988	2710455	2.78%	0.647%

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Integration Parameters: RTEINT.P

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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_X\Method\82X051222W.M

Title : SW846 8260

37	13.171	1978	1982	1986	rVB	2014050	2283296	2.34%	0.545%
38	13.292	1998	2002	2007	rVB2	3799066	4900825	5.02%	1.170%
39	13.427	2019	2024	2032	rVB	917882	1162474	1.19%	0.278%
40	13.780	2076	2082	2088	rBV	7792159	10002877	10.25%	2.389%
41	14.640	2218	2223	2233	rVB	2815769	3486414	3.57%	0.833%
42	14.780	2241	2246	2256	rBV	1457981	1812736	1.86%	0.433%

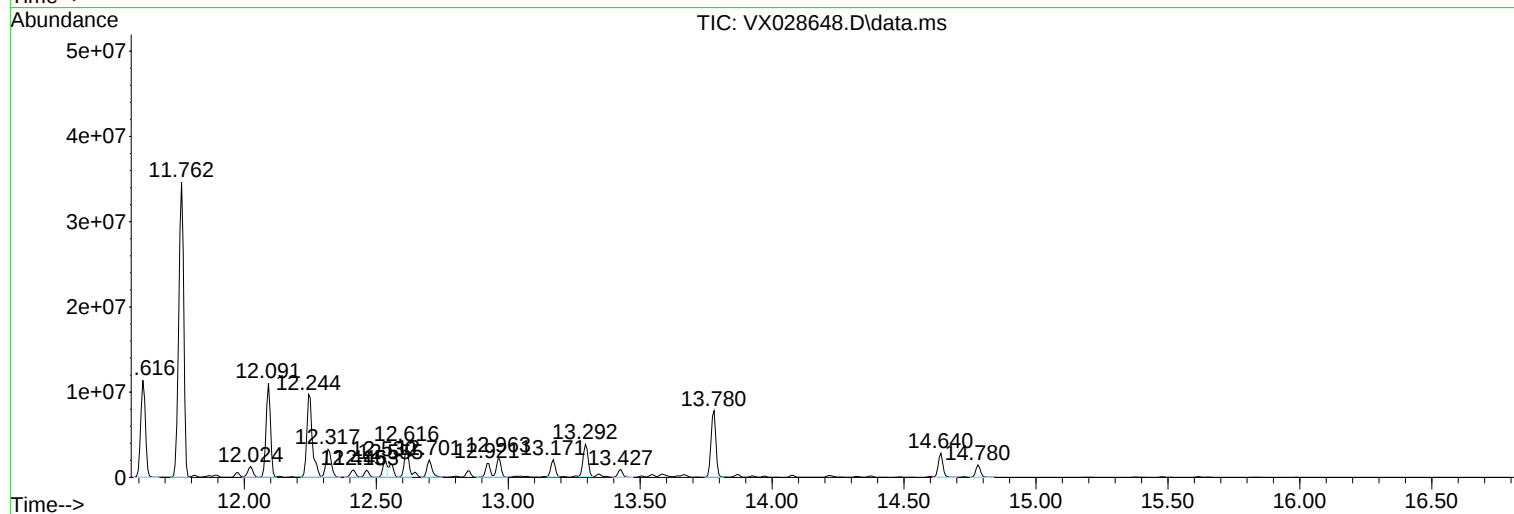
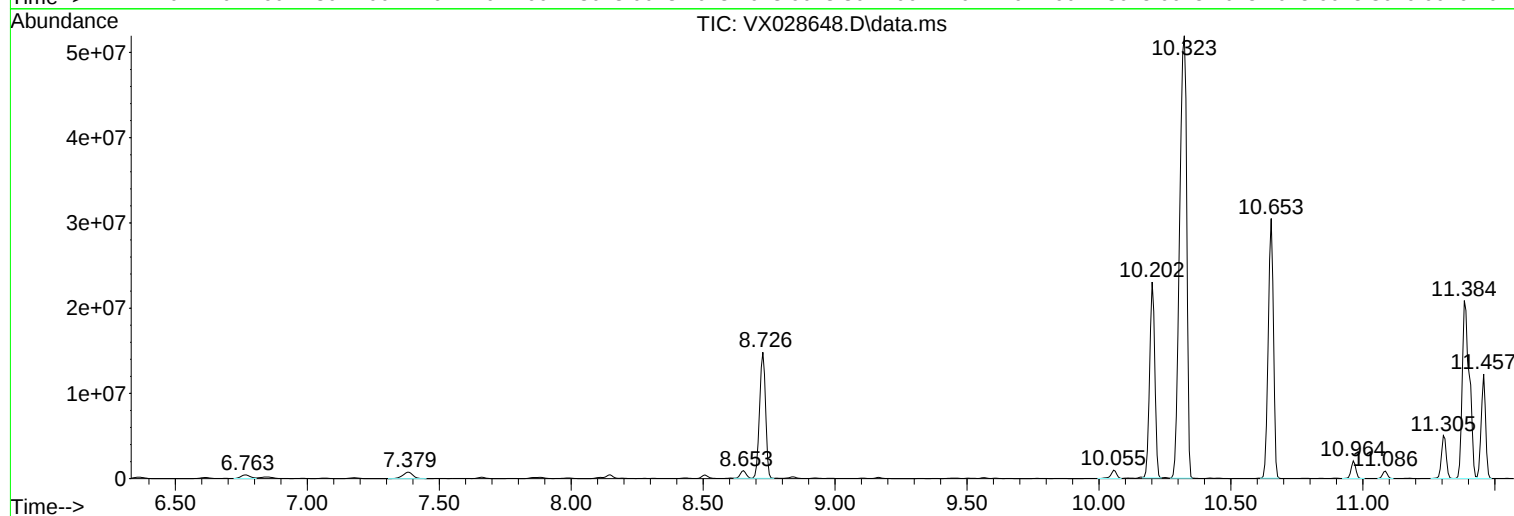
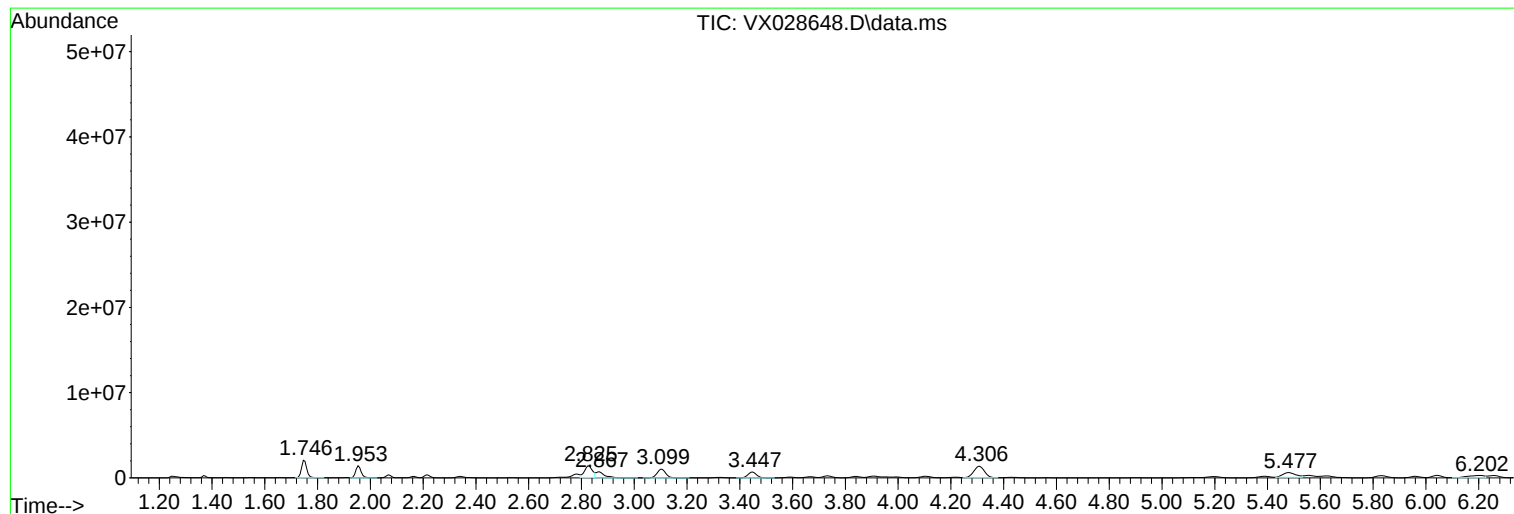
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

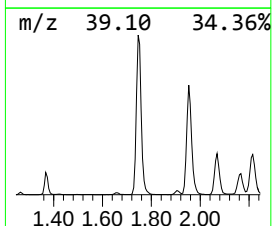
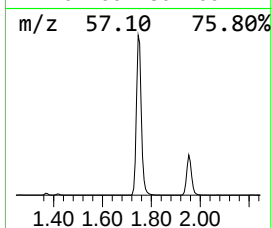
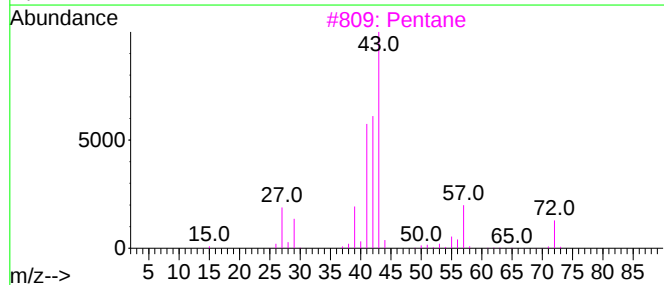
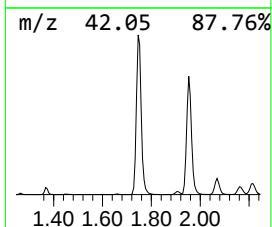
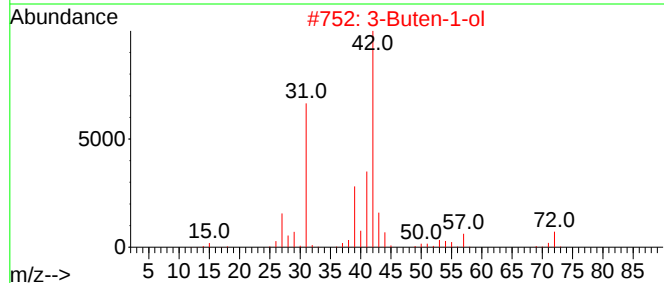
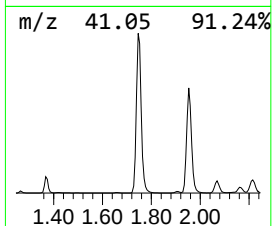
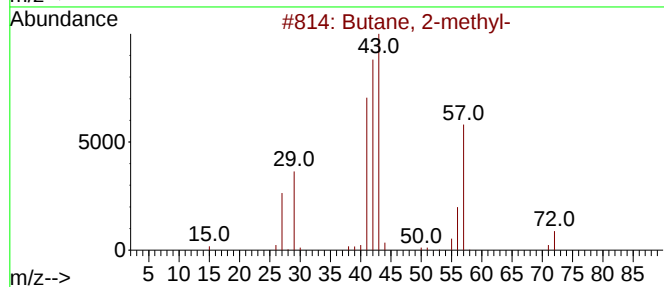
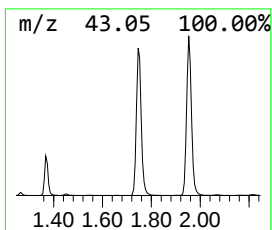
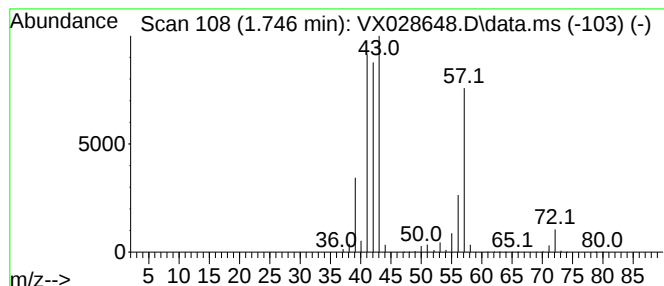
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.746	75.49 ug/l	2865420	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	87
2			3-Buten-1-ol	72	C4H8O	000627-27-0	43
3			Pentane	72	C5H12	000109-66-0	28
4			1-Butene	56	C4H8	000106-98-9	10
5			Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	10



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

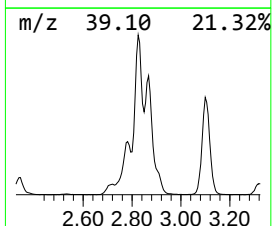
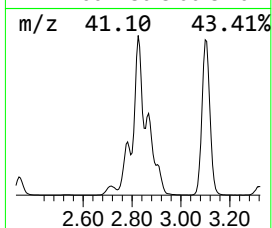
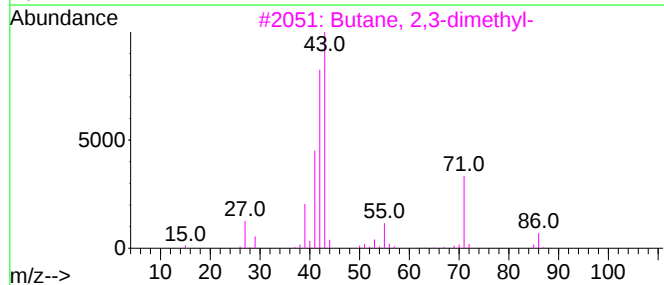
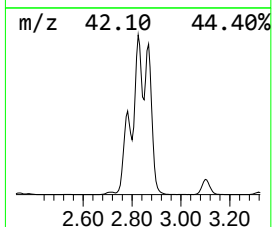
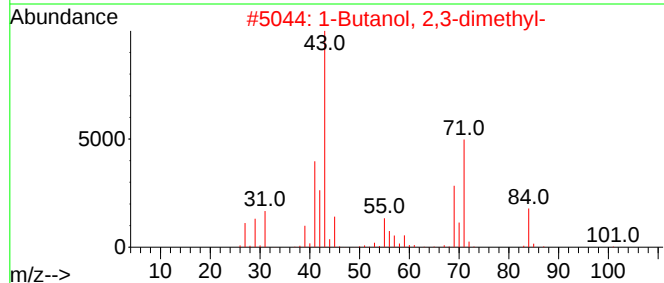
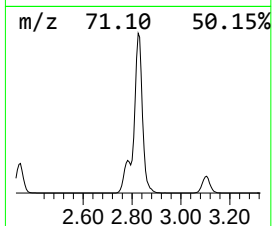
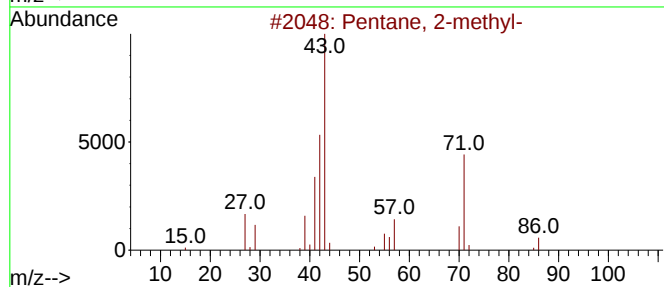
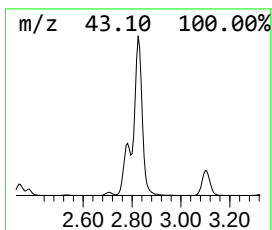
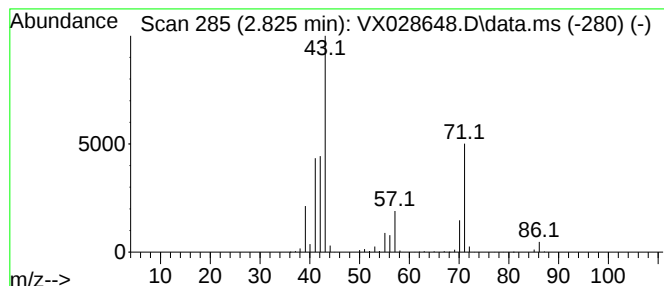
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.825	73.96 ug/l	2807520	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2			1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	43
4			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	40
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

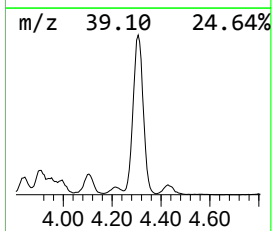
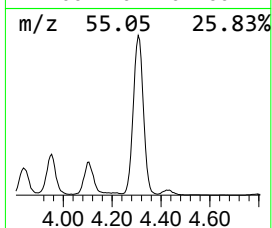
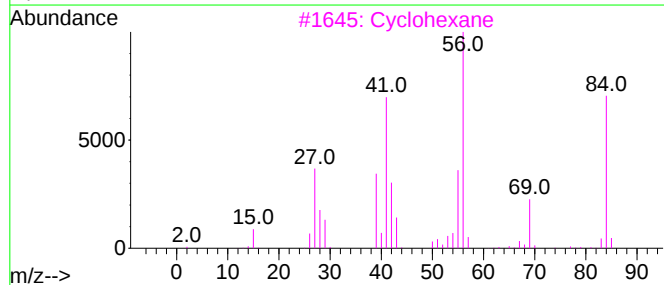
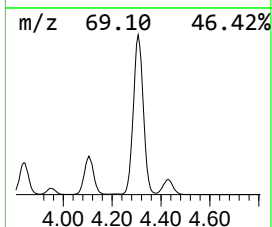
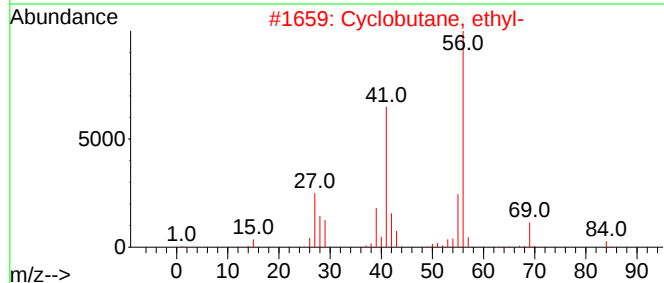
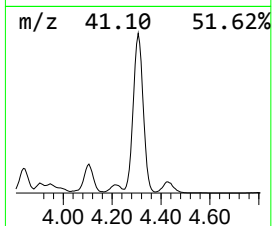
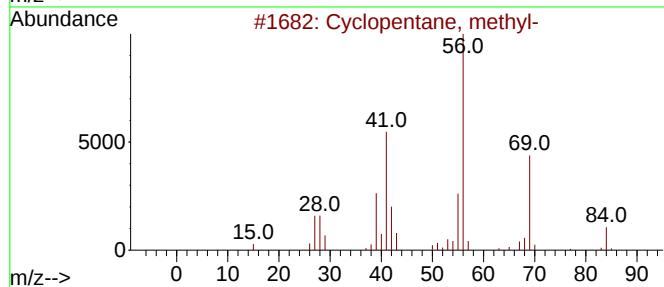
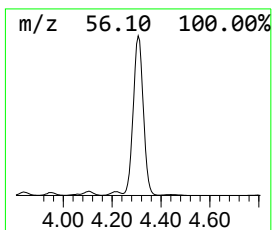
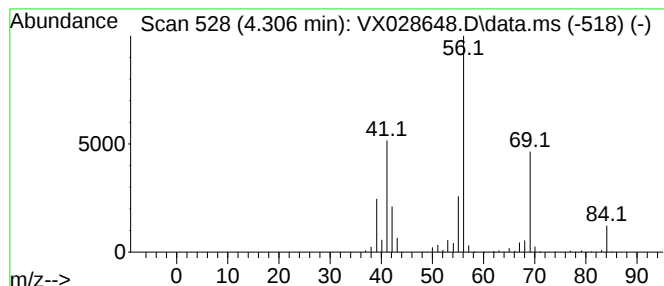
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	101.40 ug/l	3848780	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3			Cyclohexane	84	C6H12	000110-82-7	72
4			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43



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Instrument :
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ClientSampleId :
MW-2A

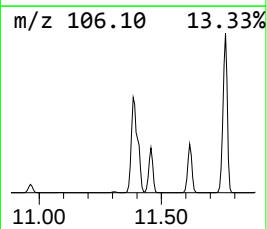
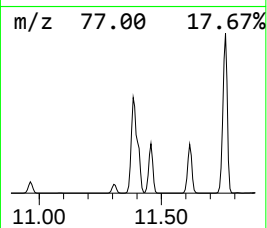
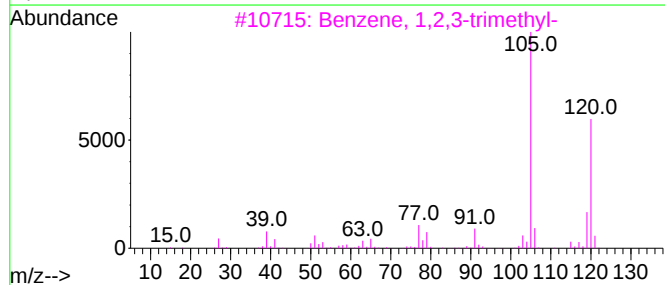
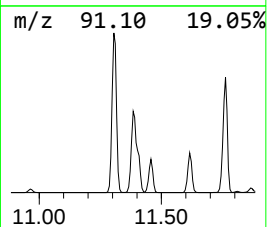
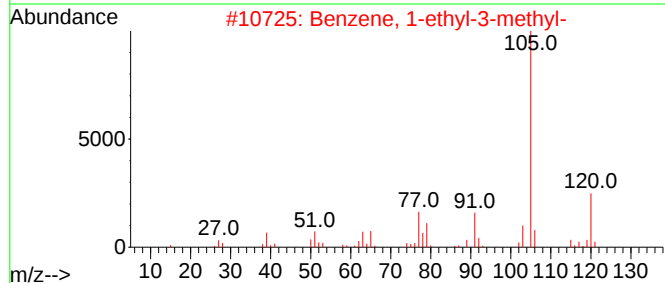
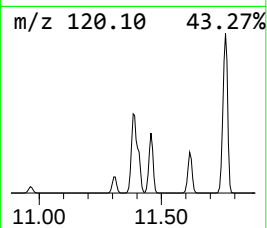
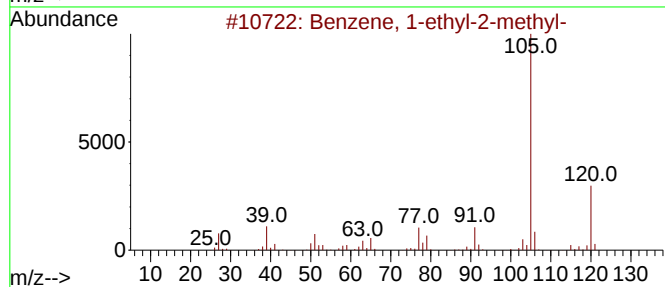
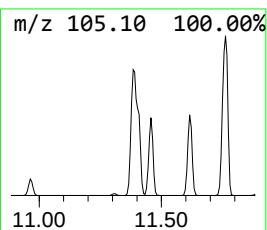
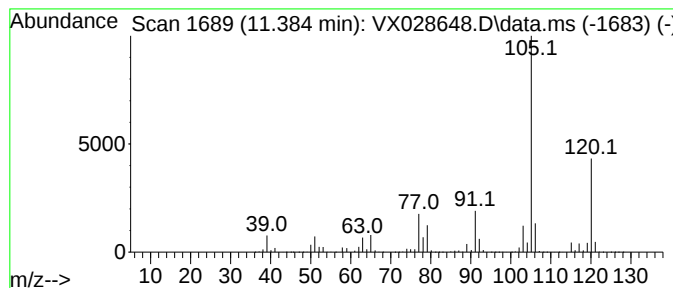
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	1157.07 ug/l	39361300	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87
5			Mesitylene	120	C9H12	000108-67-8	87



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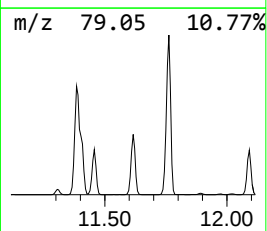
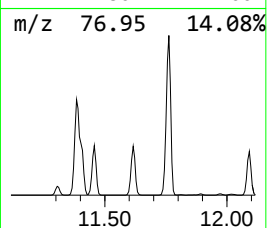
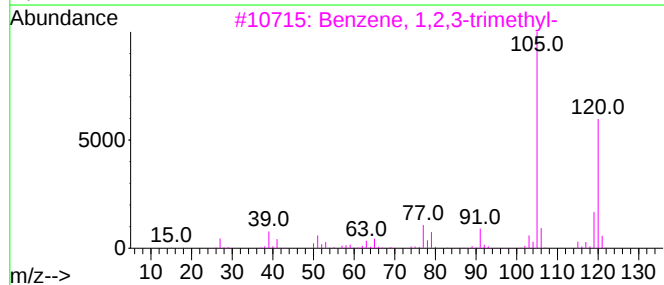
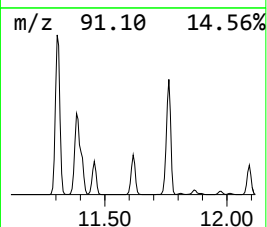
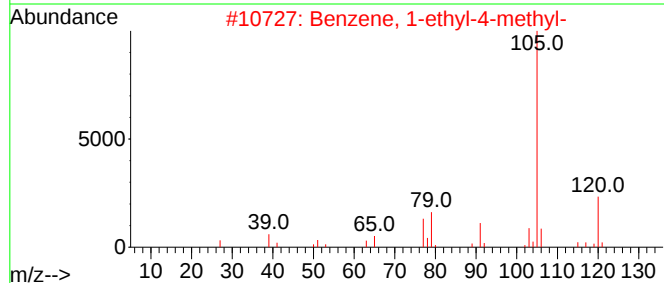
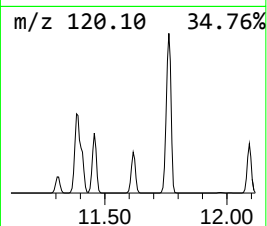
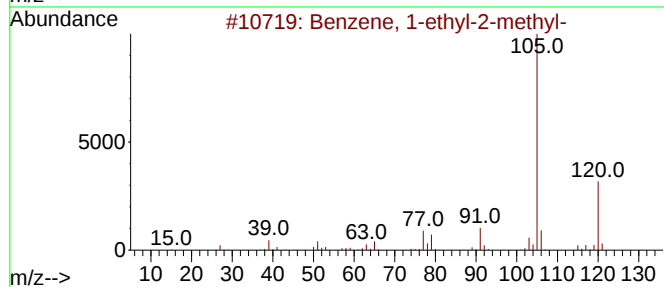
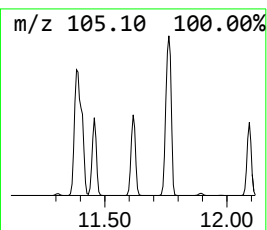
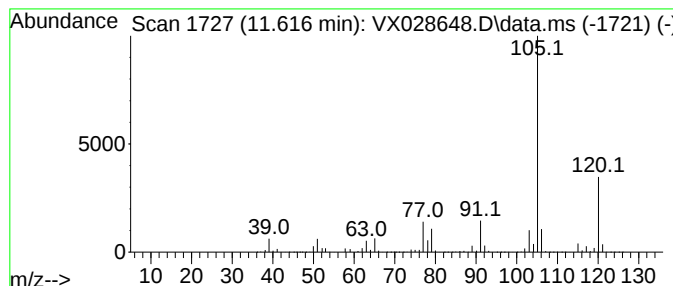
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	407.87 ug/l	13875100	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

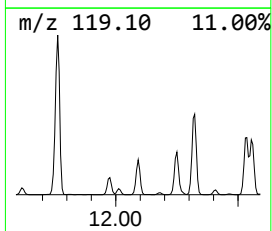
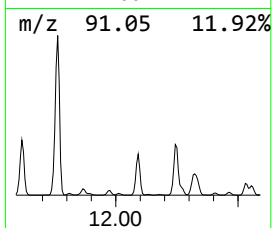
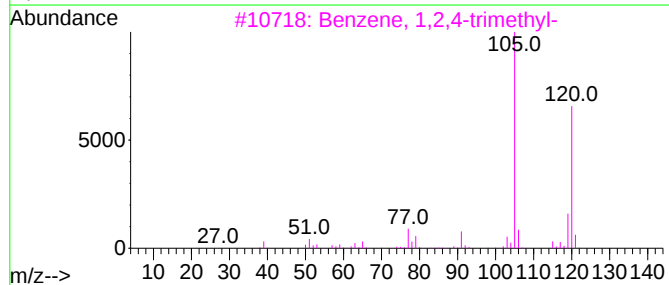
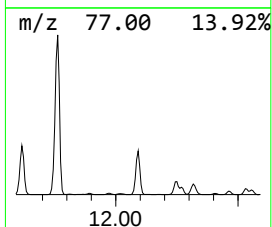
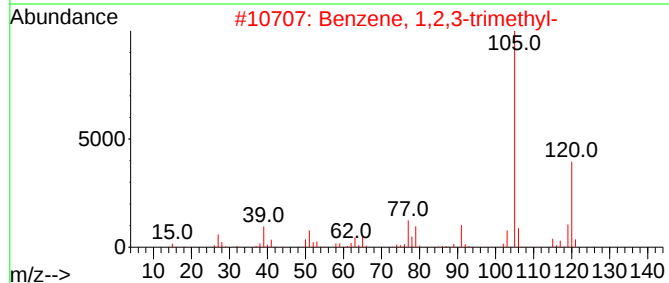
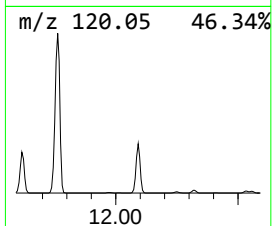
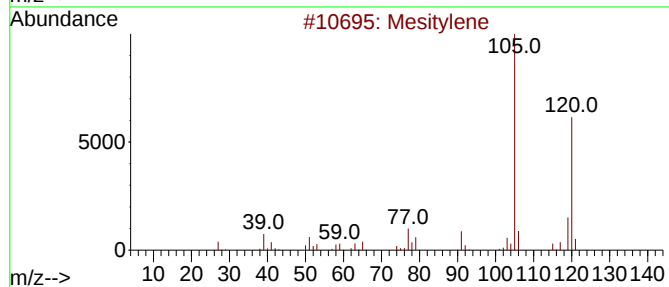
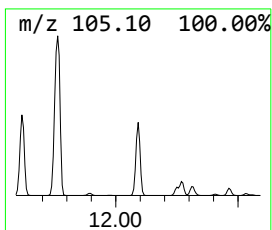
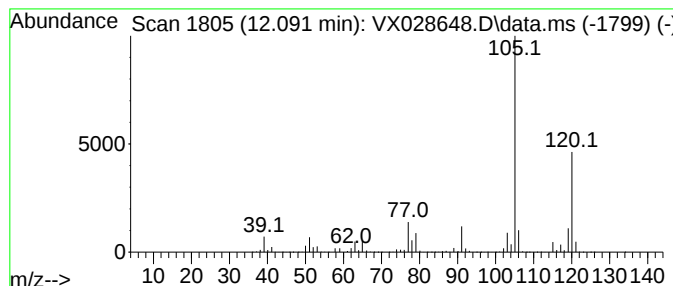
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	382.99 ug/l	13028700	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

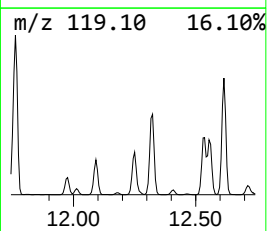
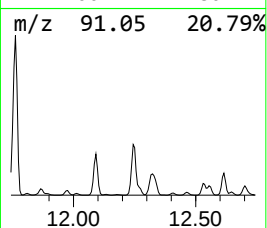
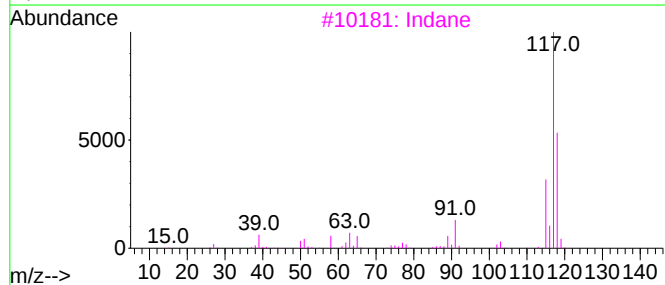
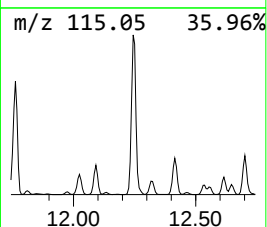
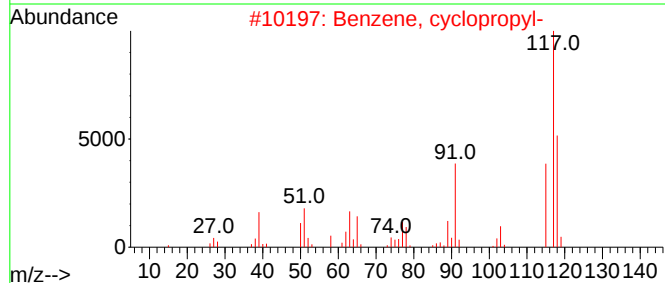
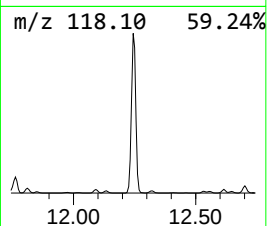
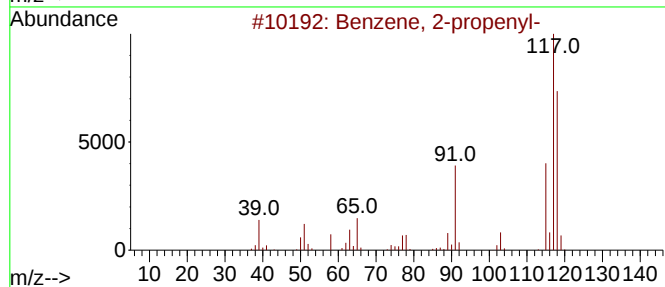
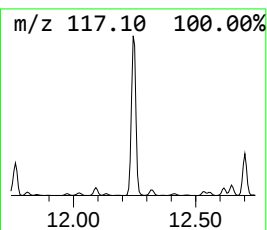
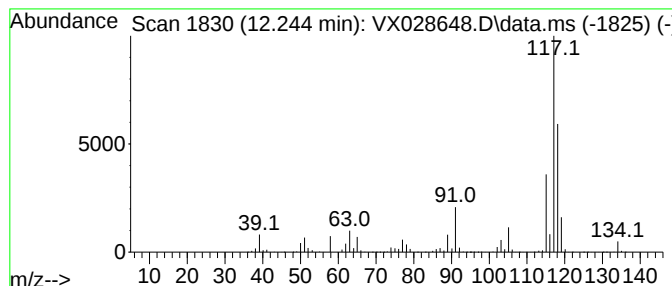
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-propenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	417.97 ug/l	14218600	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3			Indane	118	C9H10	000496-11-7	76
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	70
5			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

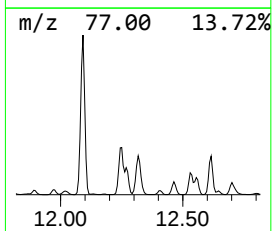
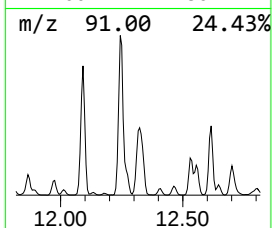
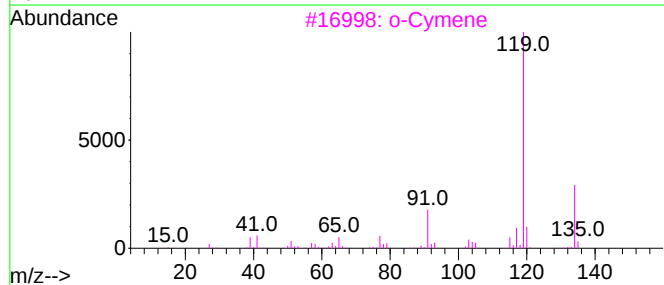
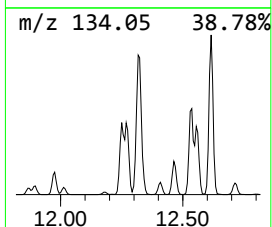
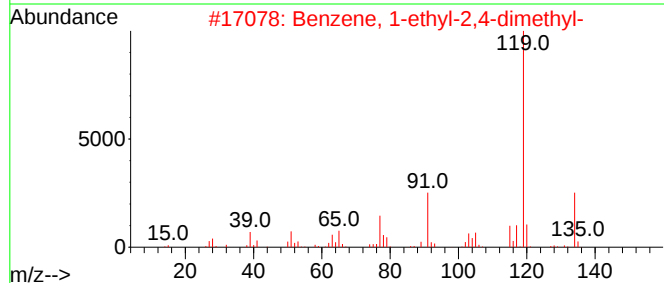
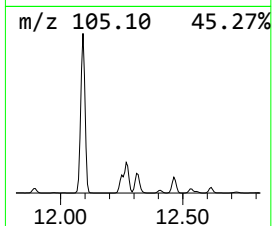
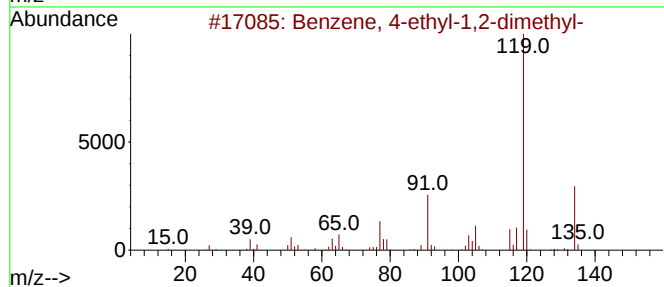
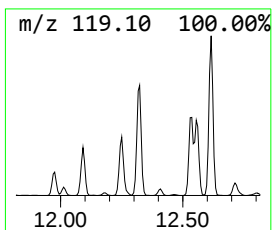
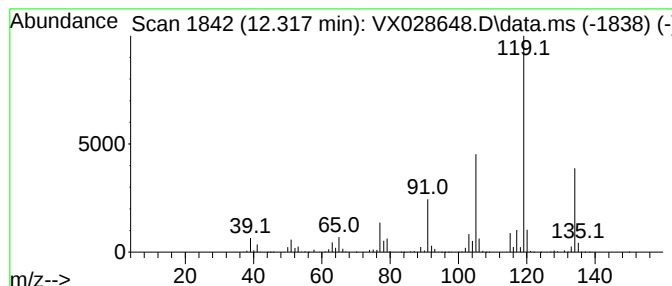
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	138.53 ug/l	4712590	1,4-Dichlorobenzene-d4	12.024

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-2,4-dimethyl-	134	C10H14	000874-41-9	94
3			o-Cymene	134	C10H14	000527-84-4	94
4			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
5			p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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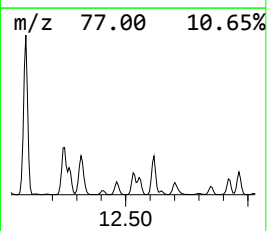
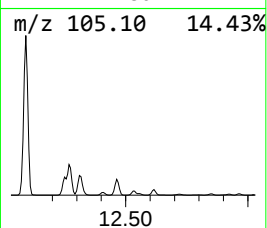
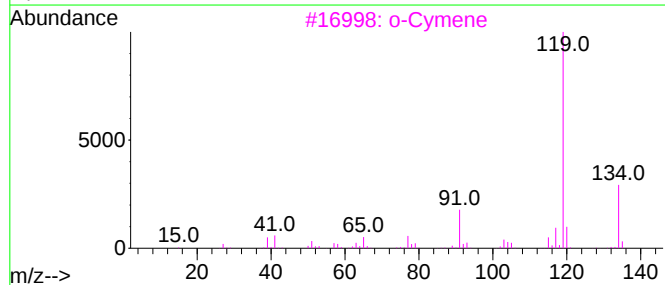
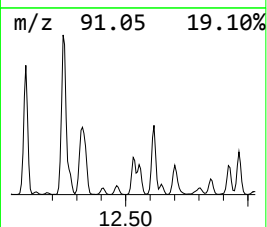
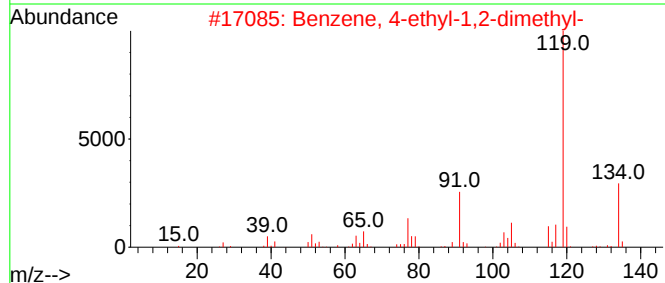
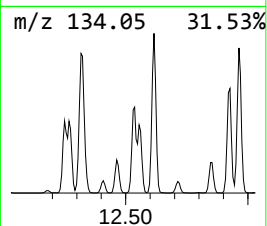
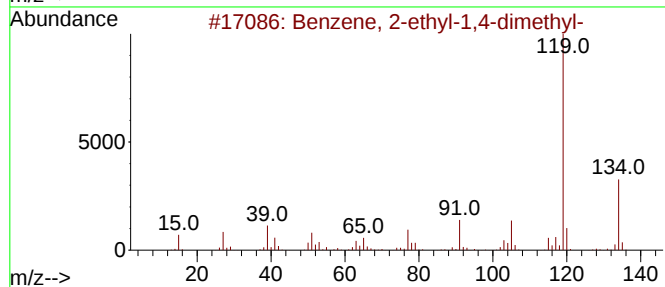
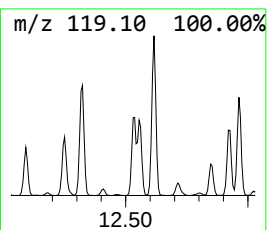
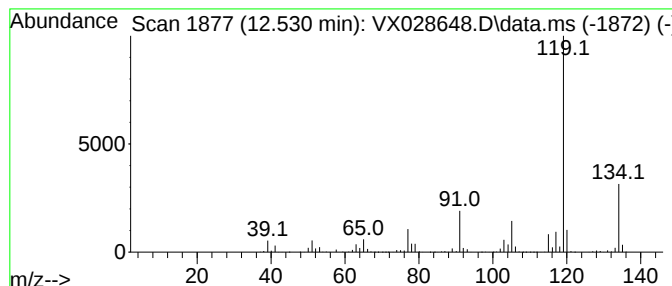
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.530	68.85 ug/l	2342260	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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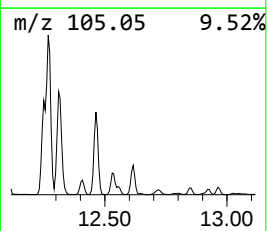
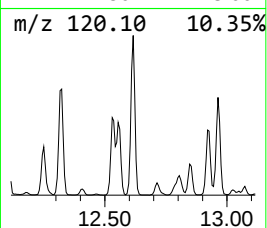
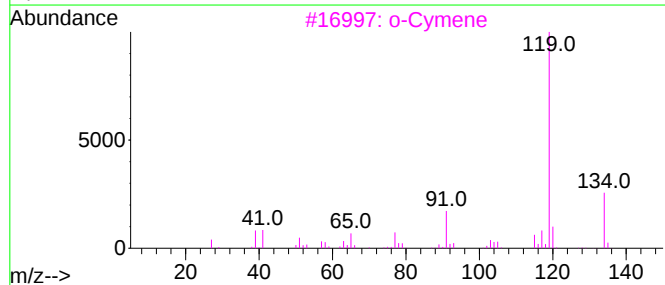
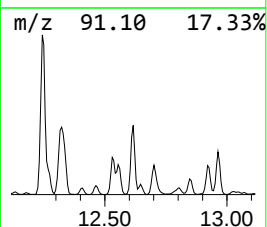
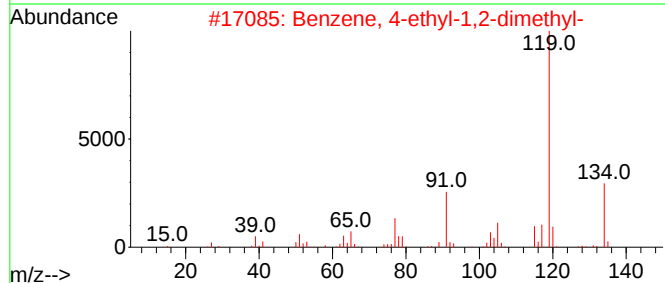
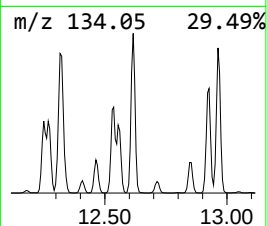
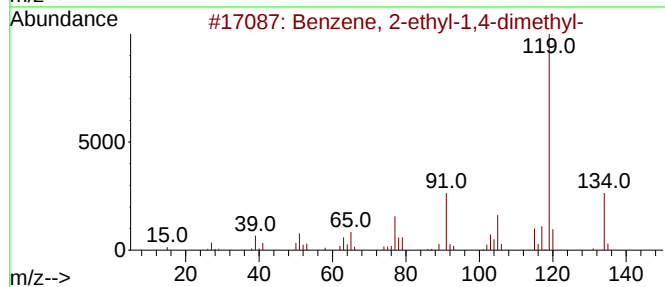
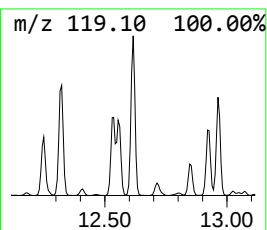
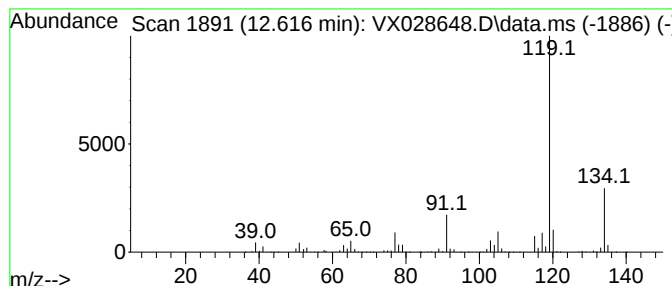
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Quant Title : SW846 8260

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

Peak Number 10 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	124.95 ug/l	4250450	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
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ALS Vial : 31 Sample Multiplier: 1

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ClientSampleId :
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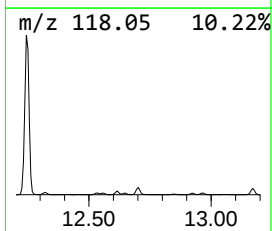
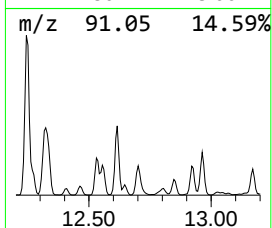
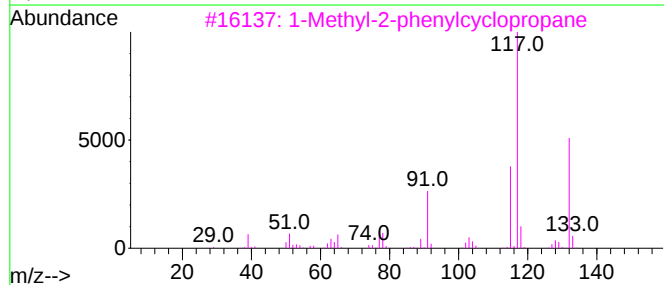
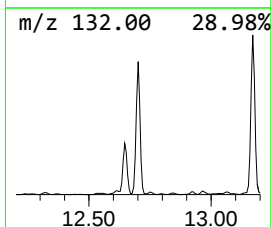
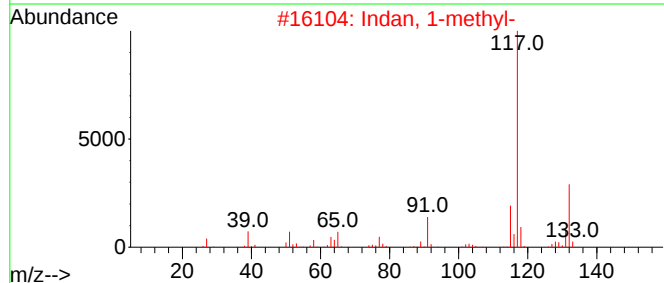
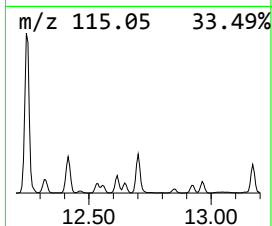
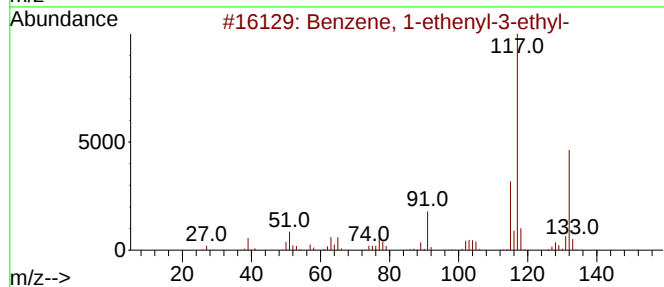
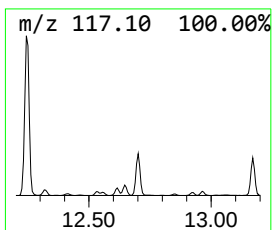
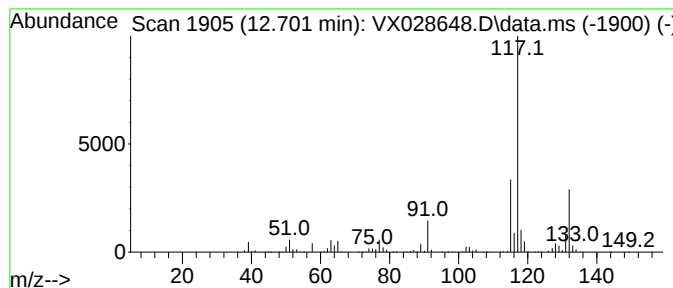
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Quant Title : SW846 8260

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

Peak Number 11 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	80.13 ug/l	2725790	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2			Indan, 1-methyl-	132	C10H12	000767-58-8	87
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

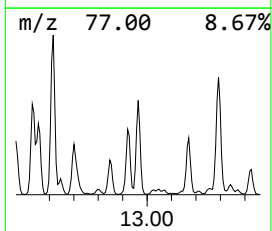
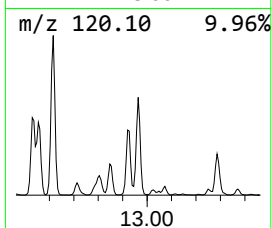
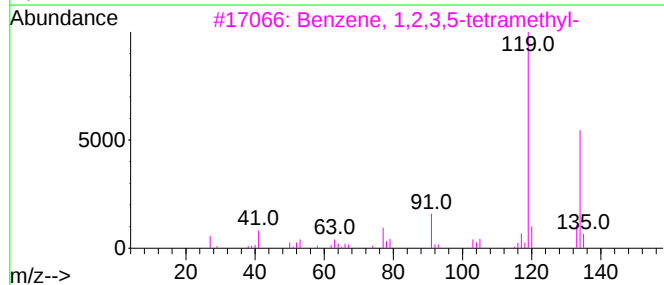
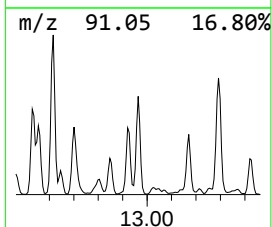
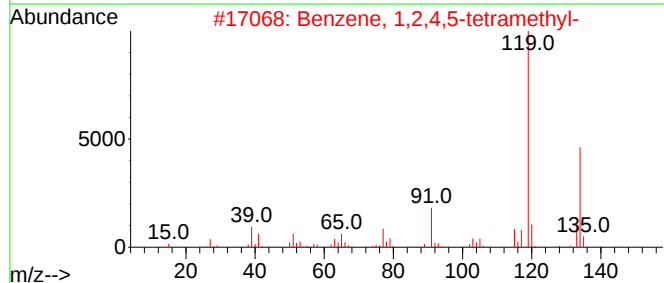
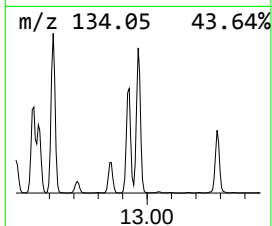
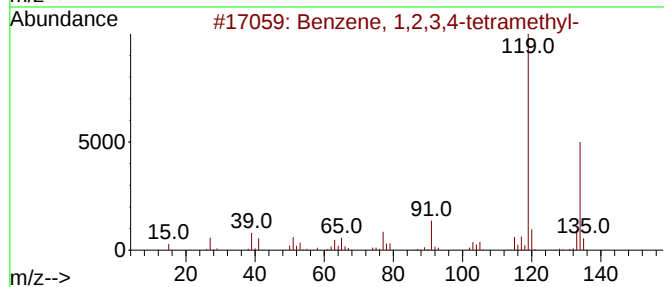
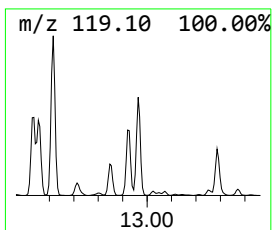
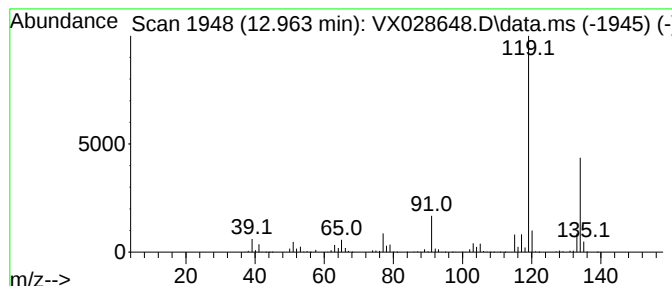
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Quant Title : SW846 8260

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

Peak Number 12 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	79.68 ug/l	2710460	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

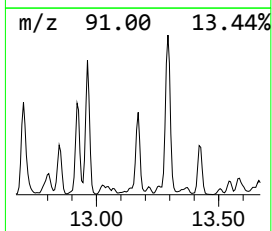
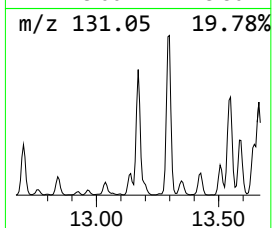
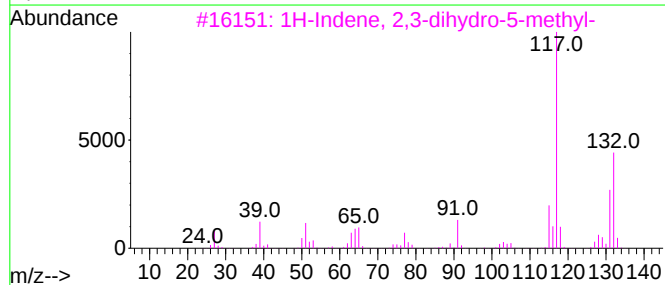
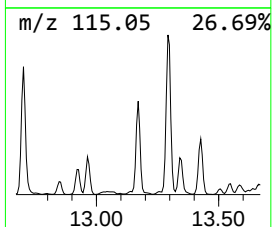
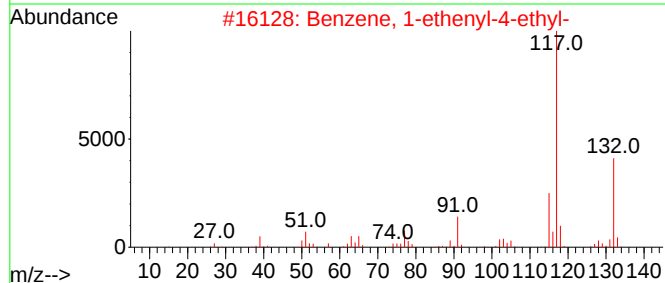
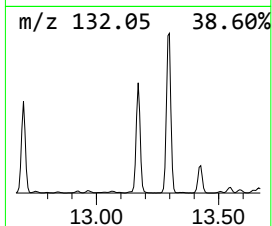
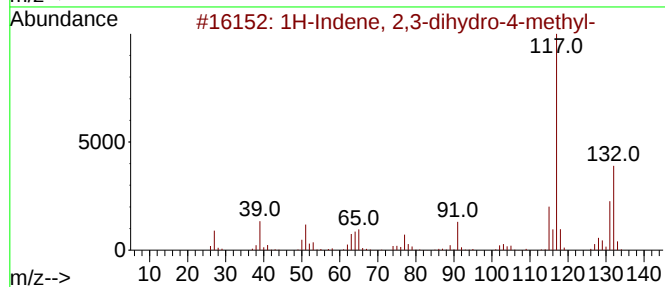
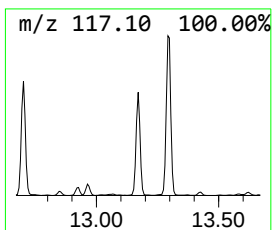
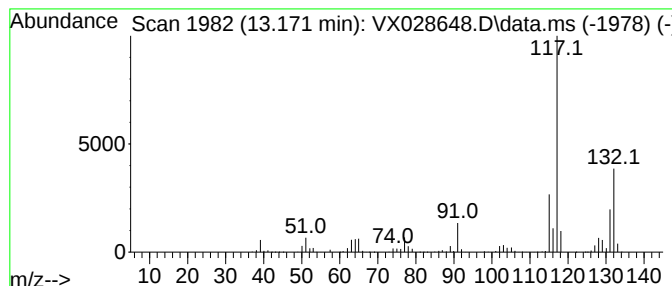
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Quant Title : SW846 8260

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

Peak Number 13 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.171	67.12 ug/l	2283300	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5		Indan, 1-methyl-	132	C10H12	000767-58-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

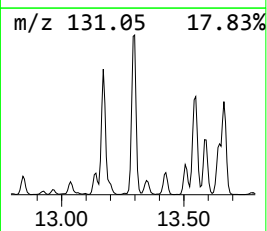
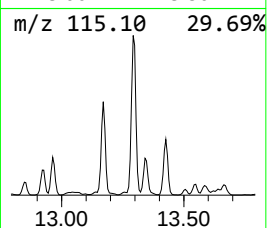
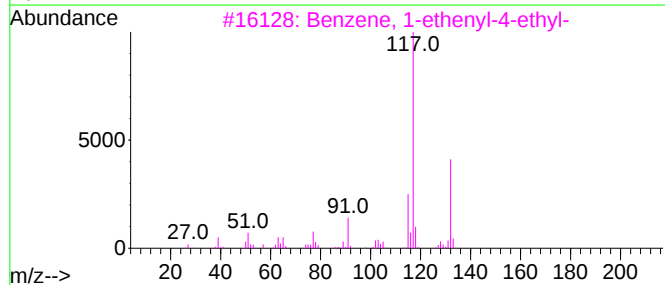
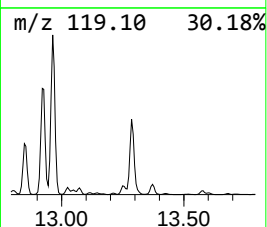
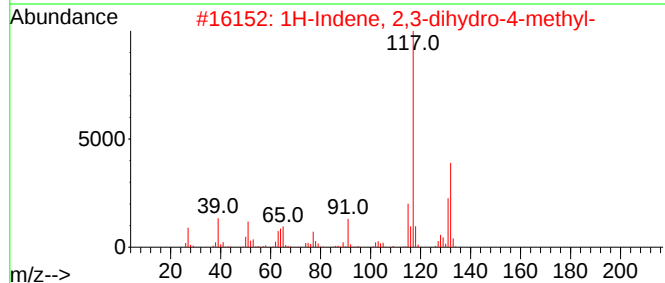
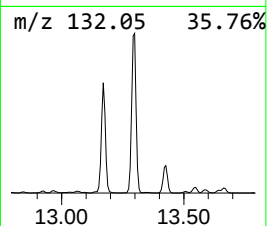
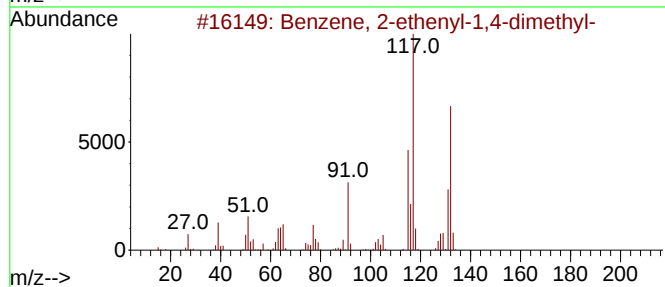
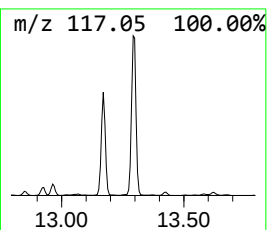
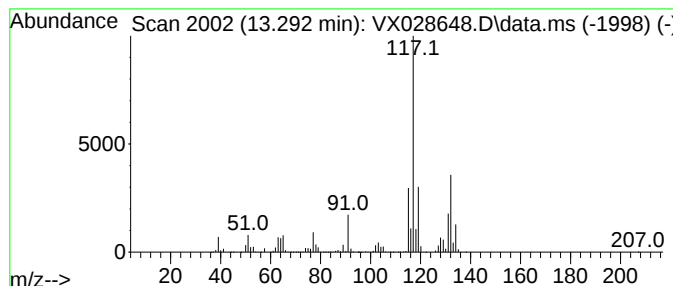
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Quant Title : SW846 8260

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

Peak Number 14 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	144.07 ug/l	4900830	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
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	87
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

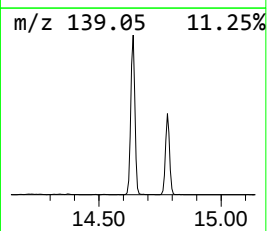
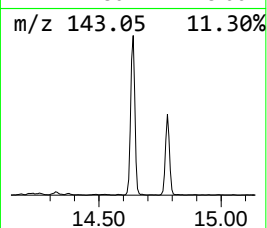
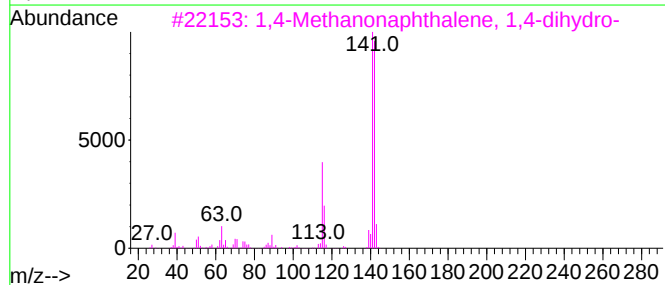
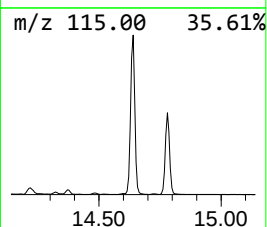
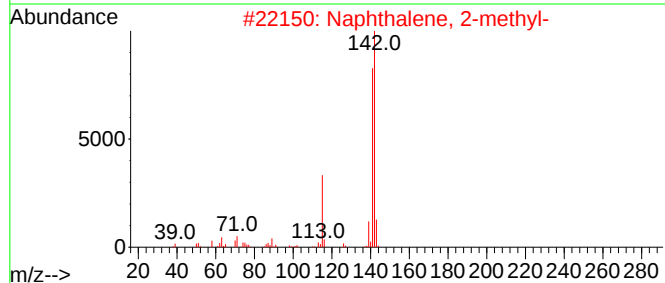
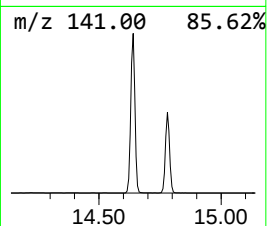
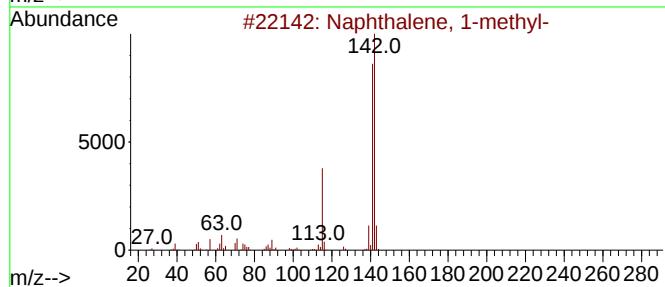
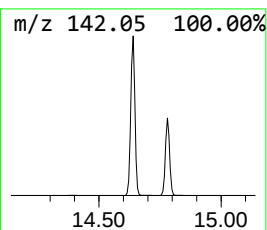
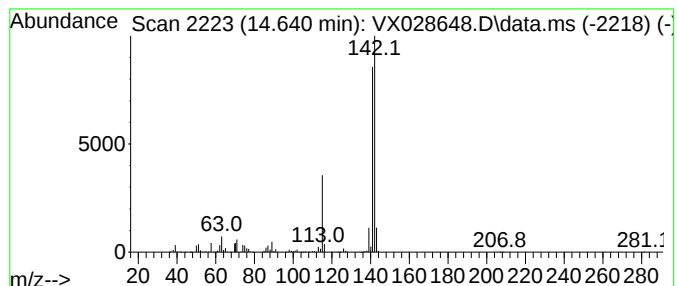
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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

Peak Number 15 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	102.49 ug/l	3486410	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051222\
Data File : VX028648.D
Acq On : 12 May 2022 19:12
Operator : JC/MD
Sample : N2799-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 31 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2A

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.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
Butane, 2-methyl-	1.746	75.5	ug/l	2865420	1	5.556	1897910	50.0
Pentane, 2-methyl-	2.825	74.0	ug/l	2807520	1	5.556	1897910	50.0
Cyclopentane, m...	4.306	101.4	ug/l	3848780	1	5.556	1897910	50.0
Benzene, 1-ethy...	11.384	1157.1	ug/l	39361300	4	12.024	1700900	50.0
Benzene, 1-ethy...	11.616	407.9	ug/l	13875100	4	12.024	1700900	50.0
Benzene, 1,2,3-...	12.091	383.0	ug/l	13028700	4	12.024	1700900	50.0
Benzene, 2-prop...	12.244	418.0	ug/l	14218600	4	12.024	1700900	50.0
Benzene, 4-ethy...	12.317	138.5	ug/l	4712590	4	12.024	1700900	50.0
Benzene, 2-ethy...	12.530	68.8	ug/l	2342260	4	12.024	1700900	50.0
o-Cymene	12.616	125.0	ug/l	4250450	4	12.024	1700900	50.0
Benzene, 1-ethe...	12.701	80.1	ug/l	2725790	4	12.024	1700900	50.0
Benzene, 1,2,3,...	12.963	79.7	ug/l	2710460	4	12.024	1700900	50.0
1H-Indene, 2,3-...	13.171	67.1	ug/l	2283300	4	12.024	1700900	50.0
Benzene, 2-ethe...	13.292	144.1	ug/l	4900830	4	12.024	1700900	50.0
Naphthalene, 1-...	14.640	102.5	ug/l	3486410	4	12.024	1700900	50.0