

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
 Data File : VN075803.D
 Acq On : 15 Dec 2022 06:27
 Operator : JC\MD
 Sample : N5990-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DUP120822

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Title : SW846 8260

Signal : TIC: VN075803.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.336	888	898	908	rBV	809916	2343996	2.47%	0.389%
2	8.236	1044	1051	1062	rVV2	690052	2529710	2.67%	0.420%
3	8.612	1103	1115	1126	rBV2	755747	2152095	2.27%	0.357%
4	8.741	1126	1137	1143	rBV4	338127	1151572	1.22%	0.191%
5	9.106	1190	1199	1206	rBV	964931	2197866	2.32%	0.365%
6	9.606	1268	1284	1300	rBV2	564638	1798111	1.90%	0.299%
7	10.106	1362	1369	1373	rBV	503249	960402	1.01%	0.159%
8	10.571	1440	1448	1454	rBV	1088309	2021393	2.13%	0.336%
9	11.865	1660	1668	1677	rBV	969475	1826485	1.93%	0.303%
10	11.971	1677	1686	1695	rBV2	55961423	94770751	100.00%	15.738%
11	12.076	1695	1704	1715	rVB2	46297579	86650421	91.43%	14.390%
12	12.400	1749	1759	1775	rBV	25364127	41216425	43.49%	6.845%
13	12.700	1799	1810	1818	rBV	3086520	5108205	5.39%	0.848%
14	12.853	1827	1836	1845	rVB	770918	1355882	1.43%	0.225%
15	13.041	1860	1868	1873	rBV	8879527	14197749	14.98%	2.358%
16	13.100	1873	1878	1881	rVV	13504020	21774883	22.98%	3.616%
17	13.135	1881	1884	1888	rVV	12852319	19666075	20.75%	3.266%
18	13.176	1888	1891	1899	rVB	10371023	15077042	15.91%	2.504%
19	13.341	1911	1919	1929	rBV	17799696	28610924	30.19%	4.751%
20	13.488	1934	1944	1957	rVV2	56259300	89211344	94.13%	14.815%
21	13.618	1957	1966	1971	rVV2	392865	981762	1.04%	0.163%
22	13.823	1990	2001	2011	rBV	20654217	34405442	36.30%	5.714%
23	13.953	2017	2023	2025	rBV	2209127	3232774	3.41%	0.537%
24	14.000	2025	2031	2047	rVV2	15773990	34118703	36.00%	5.666%
25	14.182	2056	2062	2068	rVV2	1823144	3105372	3.28%	0.516%
26	14.247	2068	2073	2076	rVV	2960422	4997855	5.27%	0.830%
27	14.276	2076	2078	2083	rVV	2127877	2836649	2.99%	0.471%
28	14.335	2083	2088	2094	rVV	5930573	8985652	9.48%	1.492%
29	14.400	2094	2099	2102	rVV	872427	1411487	1.49%	0.234%
30	14.447	2102	2107	2116	rVB2	3175934	5516965	5.82%	0.916%
31	14.576	2123	2129	2135	rBV	1446427	2211938	2.33%	0.367%
32	14.659	2135	2143	2146	rBV	3573331	5941940	6.27%	0.987%
33	14.694	2146	2149	2154	rVB	4166511	6035129	6.37%	1.002%
34	14.929	2184	2189	2195	rBV	3296046	5366667	5.66%	0.891%
35	15.053	2201	2210	2216	rBV2	6109585	13277972	14.01%	2.205%
36	15.117	2216	2221	2230	rVB2	683535	1403260	1.48%	0.233%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Title : SW846 8260

37	15.212	2230	2237	2245	rBV2	1239740	2318044	2.45%	0.385%
38	15.323	2245	2256	2261	rBV3	848637	2229430	2.35%	0.370%
39	15.441	2268	2276	2286	rVB3	689934	1770275	1.87%	0.294%
40	15.647	2288	2311	2319	rBV	14106359	27406219	28.92%	4.551%

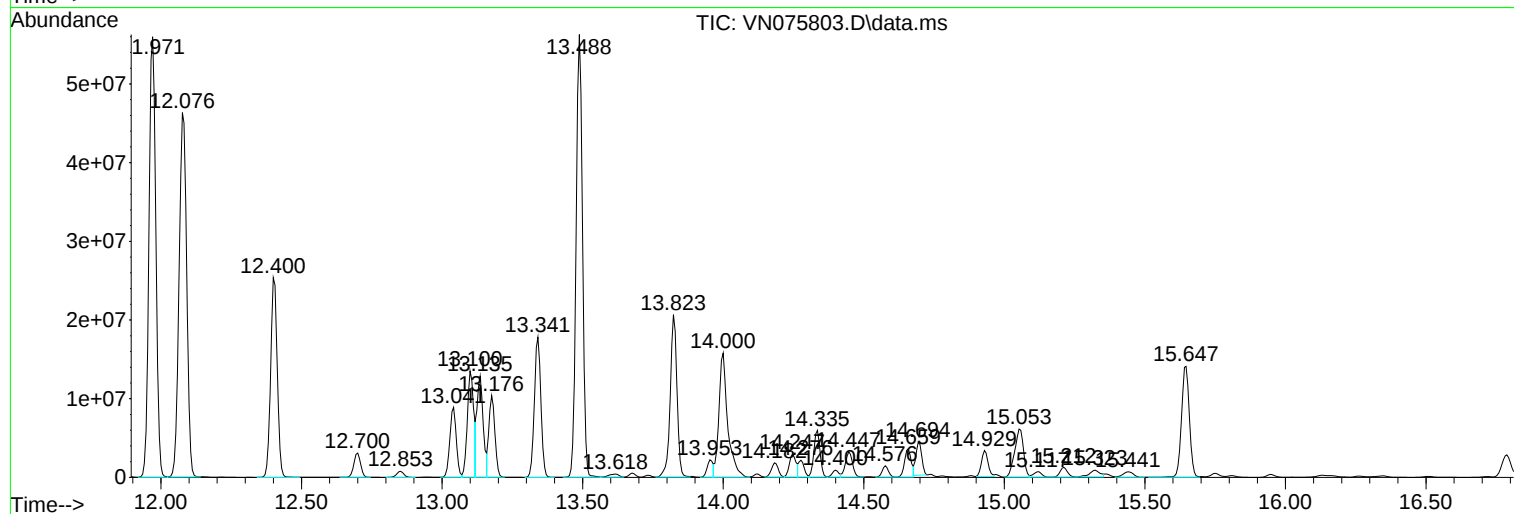
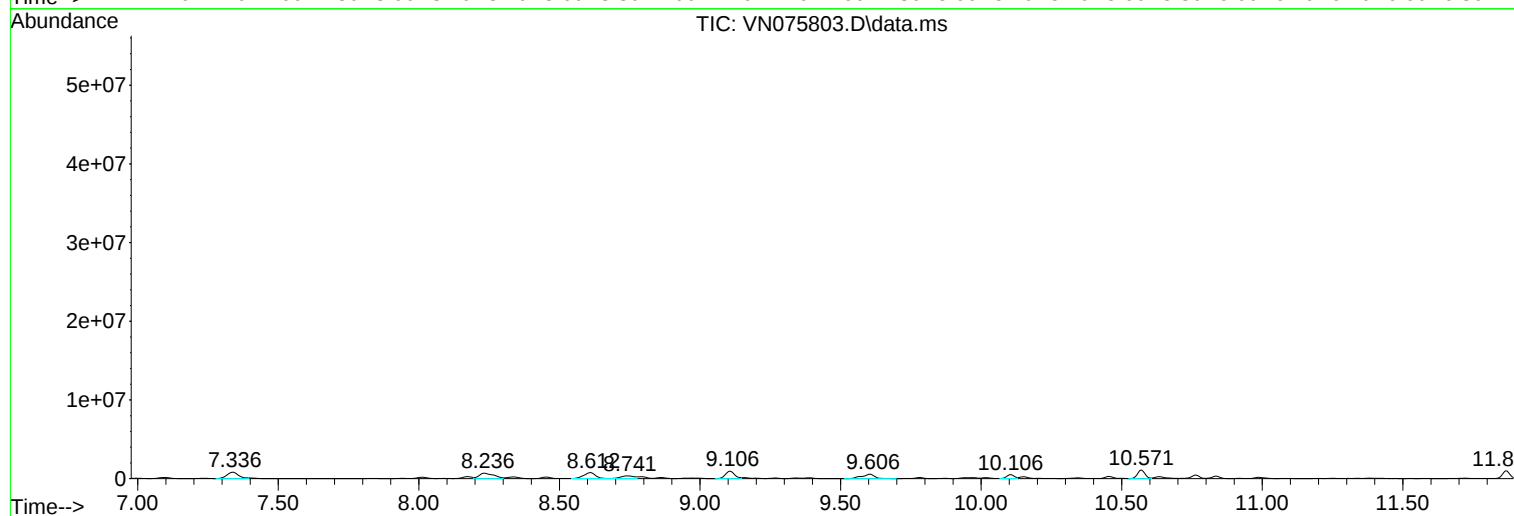
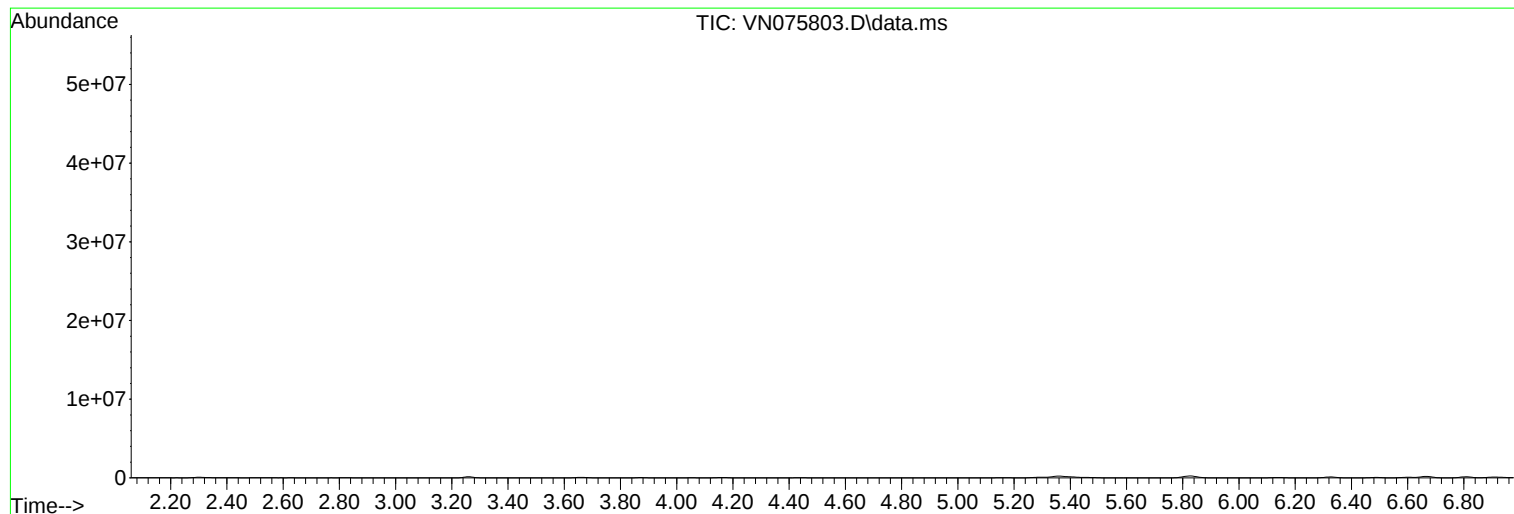
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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



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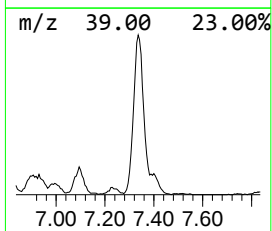
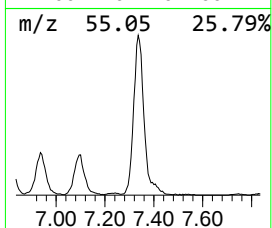
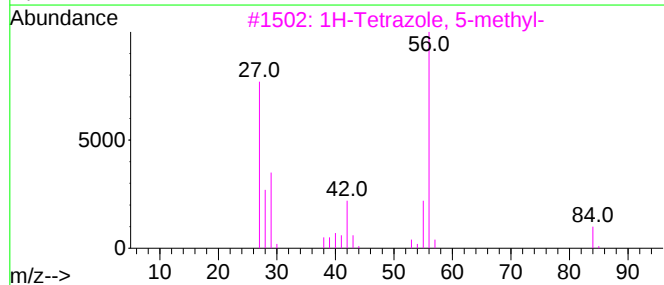
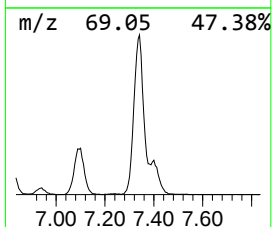
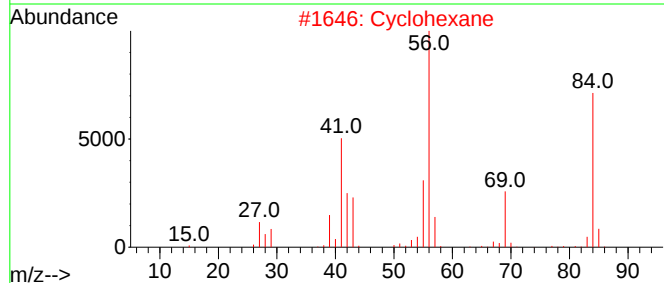
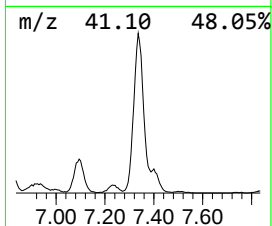
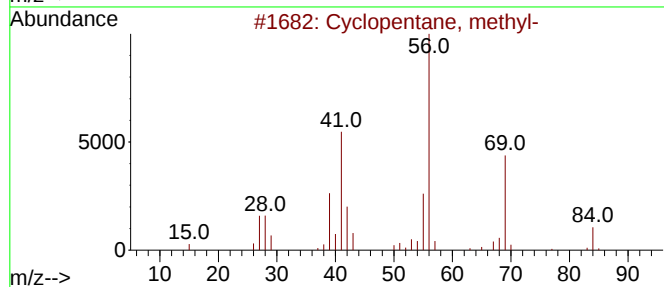
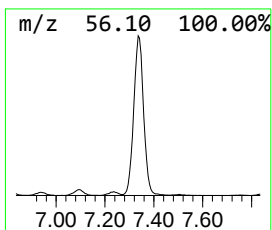
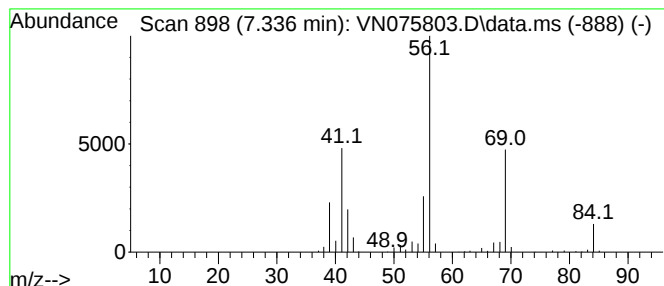
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.336	46.33 ug/l	2344000	Pentafluorobenzene	8.230

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclohexane	84	C6H12	000110-82-7	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5		Cyclobutane, butyl-	112	C8H16	013152-44-8	50



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Instrument :
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ClientSampleId :
DUP120822

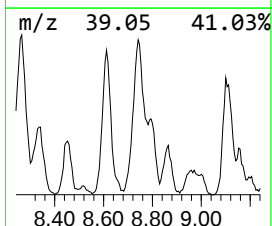
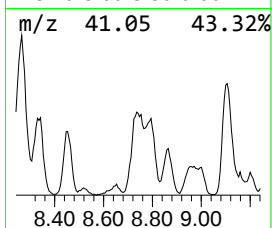
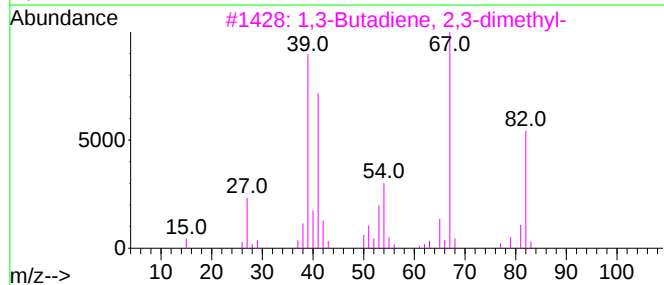
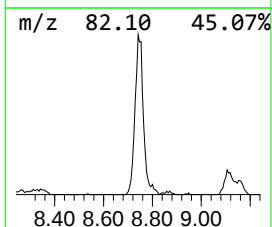
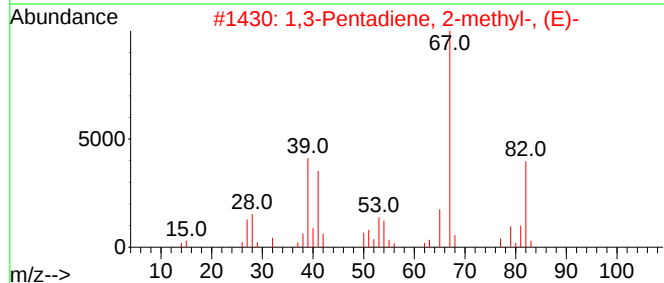
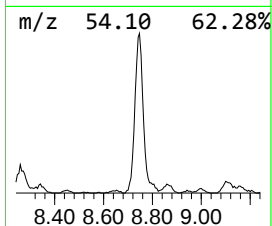
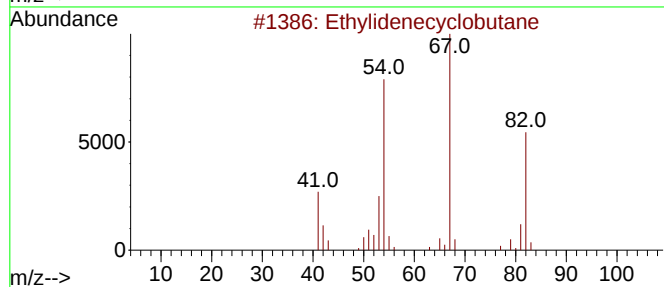
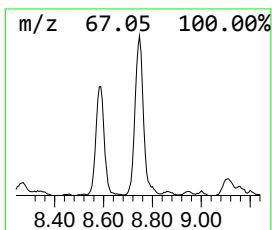
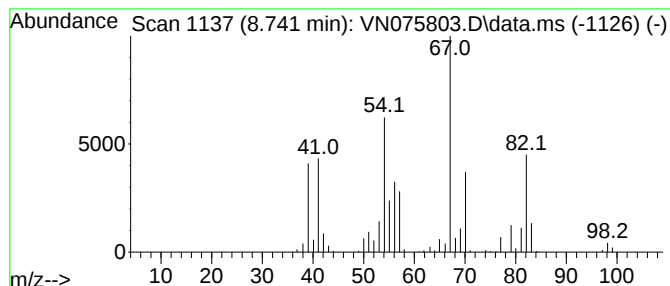
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 2 Ethylidenecyclobutane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.741	26.20 ug/l	1151570	1,4-Difluorobenzene	9.106

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylidenecyclobutane	82	C6H10	001528-21-8	76
2			1,3-Pentadiene, 2-methyl-, (E)-	82	C6H10	000926-54-5	64
3			1,3-Butadiene, 2,3-dimethyl-	82	C6H10	000513-81-5	62
4			Cyclohexene	82	C6H10	000110-83-8	60
5			Cyclobutane, ethenyl-	82	C6H10	002597-49-1	60



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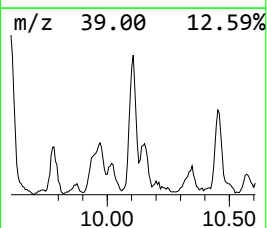
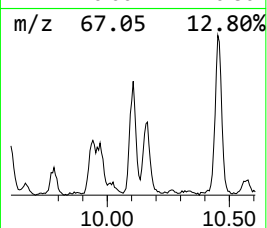
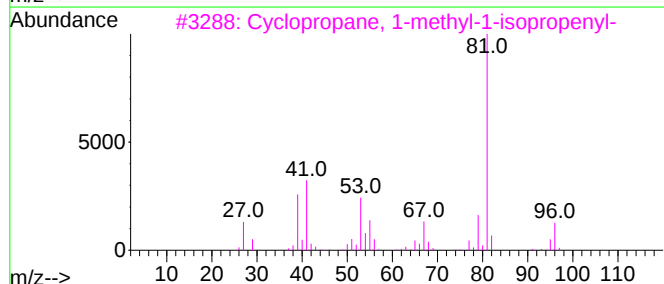
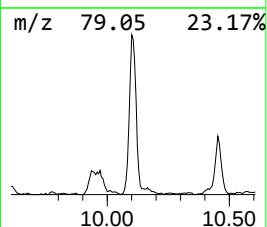
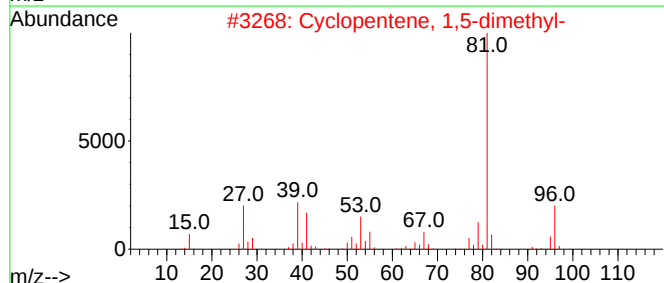
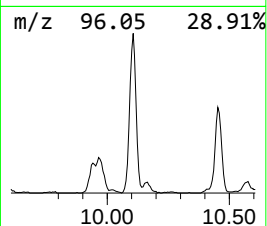
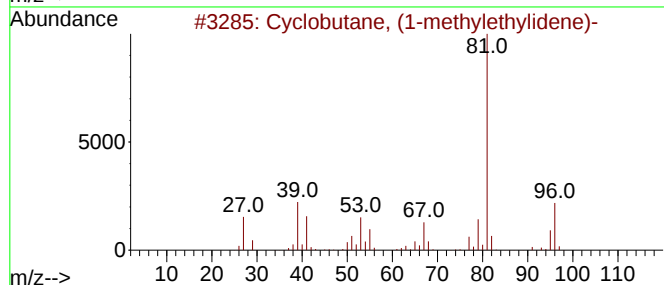
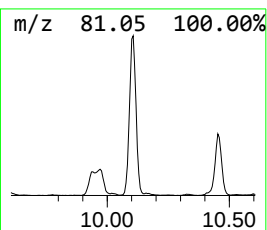
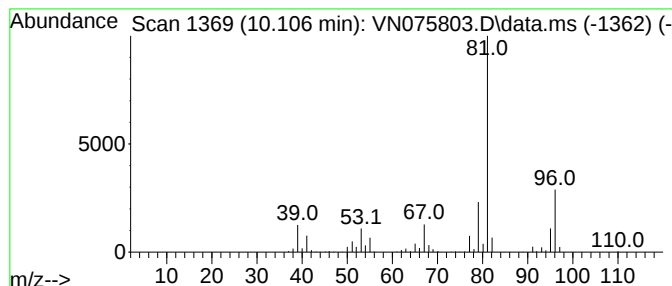
Instrument :
MSVOA_N
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclobutane, (1-methylethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.106	21.85 ug/l	960402	1,4-Difluorobenzene	9.106	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	91
2	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	91
3	Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	90
4	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	86
5	1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	78



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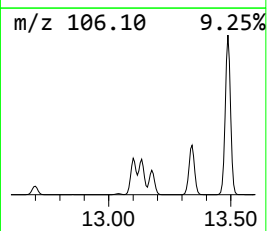
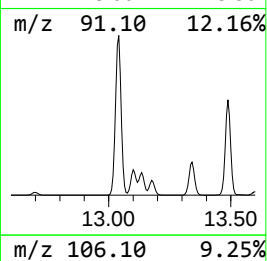
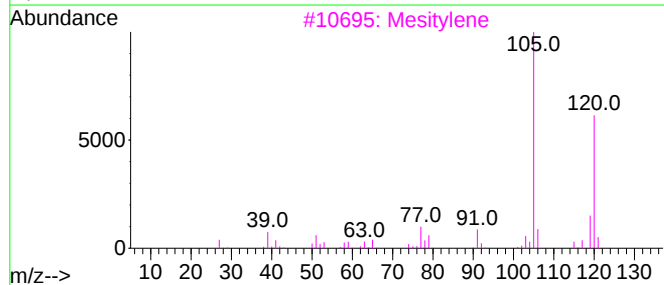
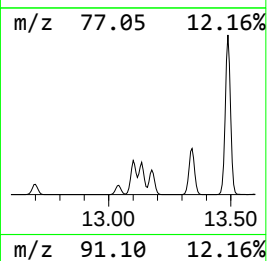
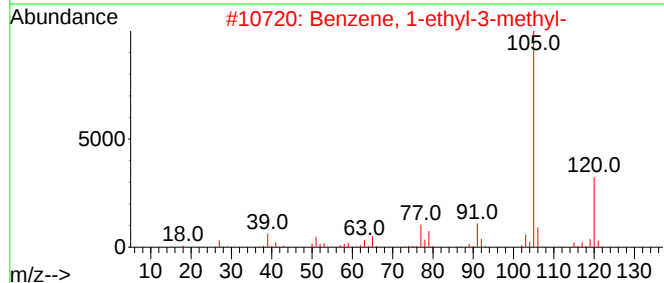
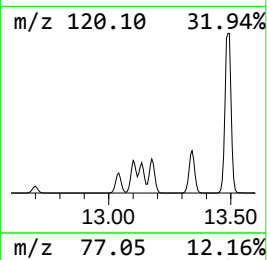
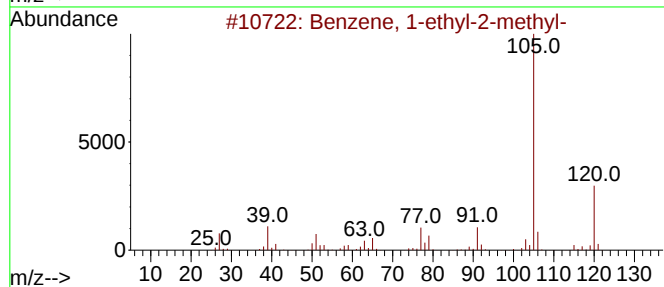
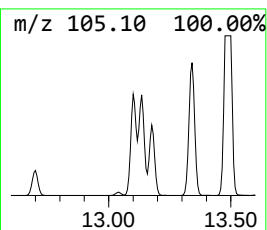
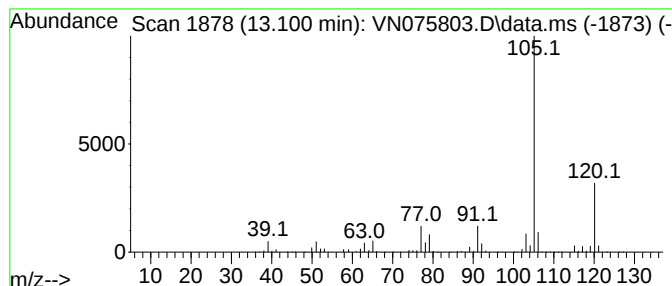
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	31.64 ug/l	21774900	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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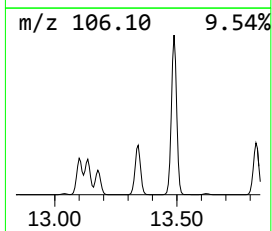
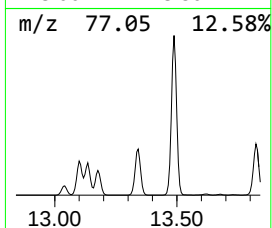
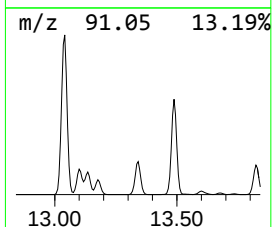
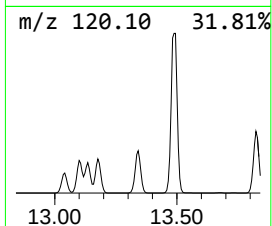
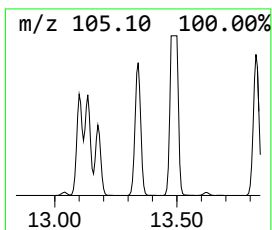
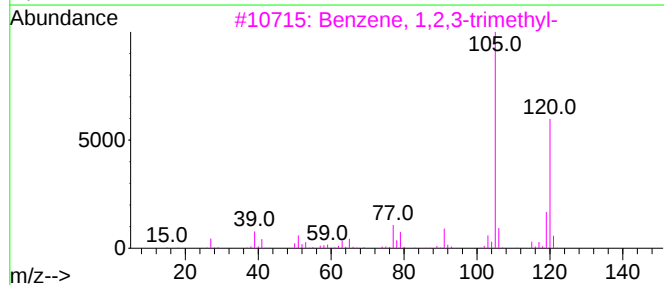
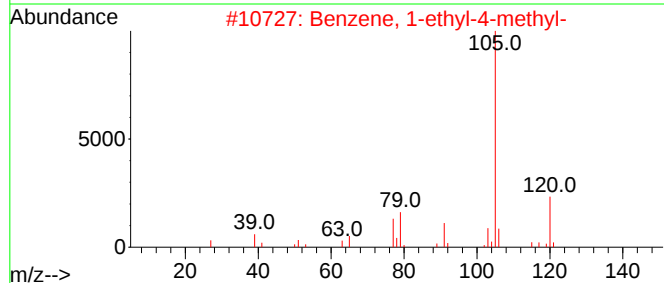
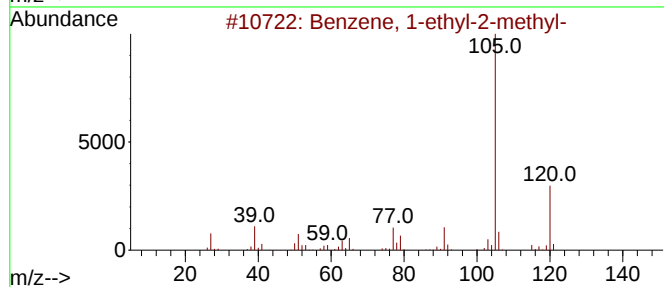
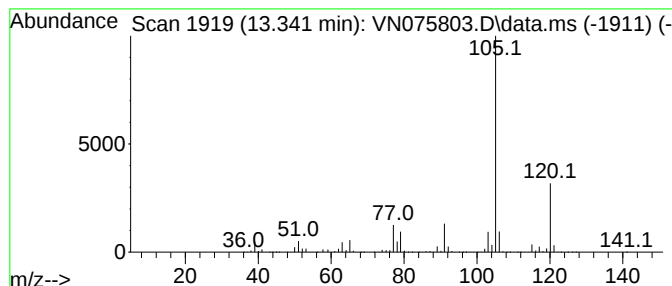
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Quant Title : SW846 8260

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.
13.341	41.58 ug/l	28610900	1,4-Dichlorobenzene-d4	13.794

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_N\Data\VN121422\
Data File : VN075803.D
Acq On : 15 Dec 2022 06:27
Operator : JC\MD
Sample : N5990-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP120822

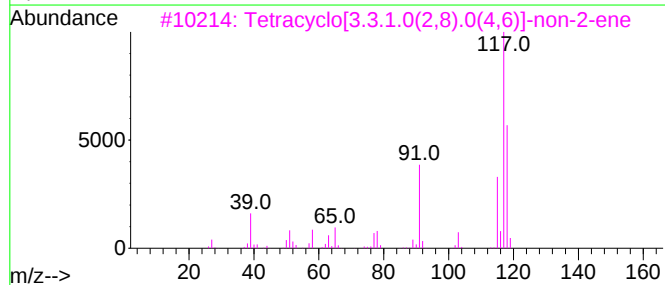
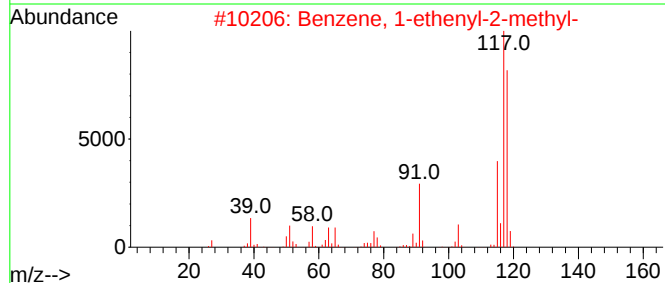
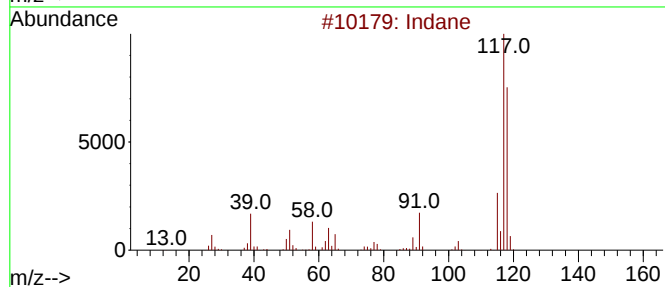
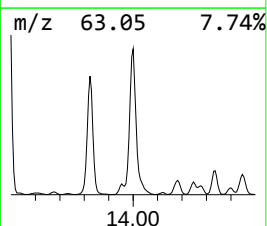
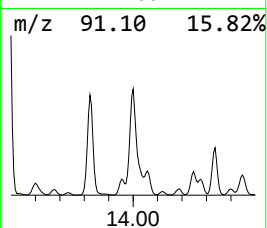
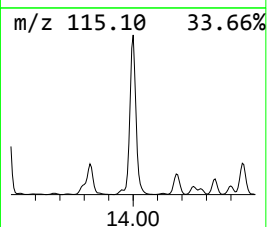
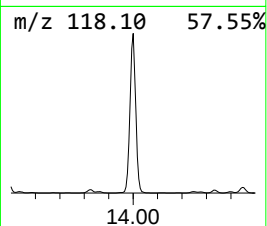
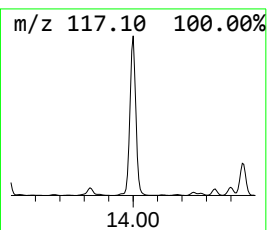
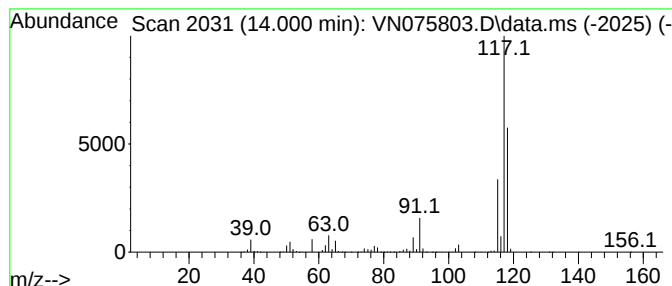
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Quant Title : SW846 8260

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

Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.000	49.58 ug/l	34118700	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
4			Deltacyclene	118	C9H10	007785-10-6	64
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075803.D
Acq On : 15 Dec 2022 06:27
Operator : JC\MD
Sample : N5990-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP120822

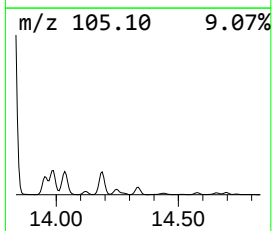
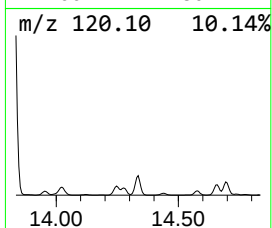
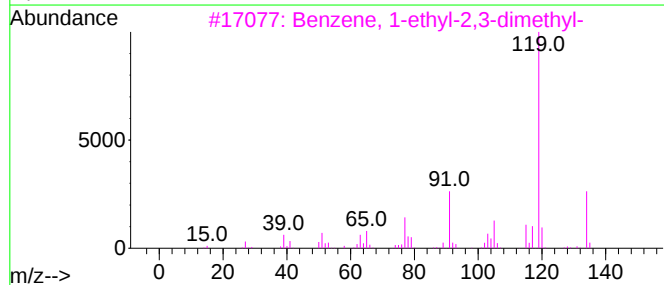
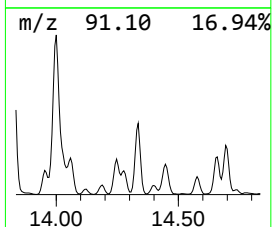
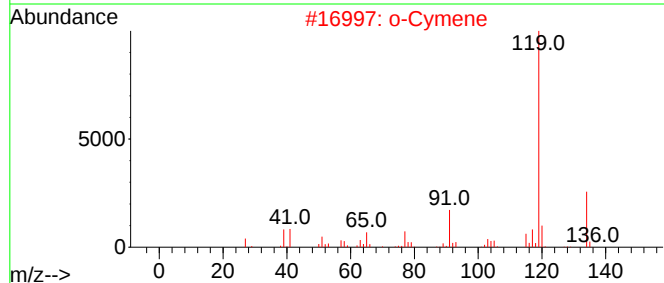
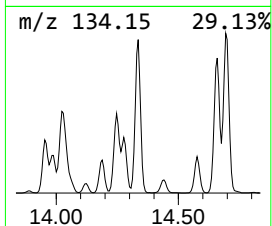
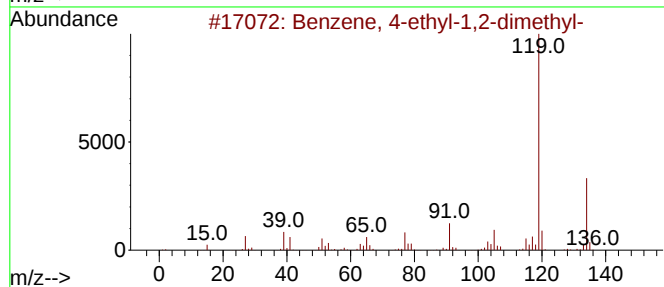
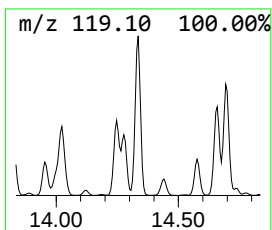
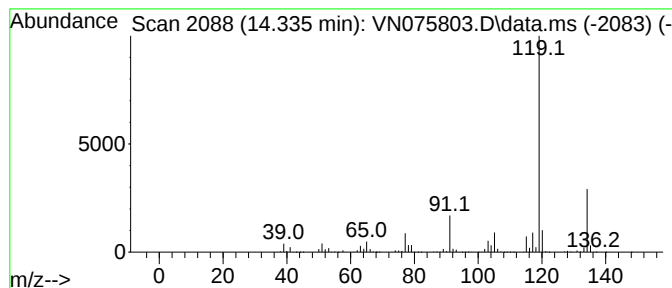
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.335	13.06 ug/l	8985650	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075803.D
Acq On : 15 Dec 2022 06:27
Operator : JC\MD
Sample : N5990-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP120822

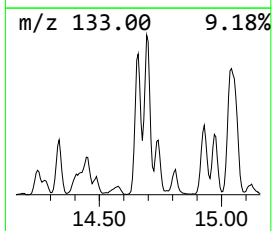
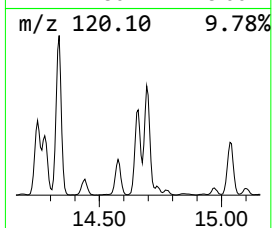
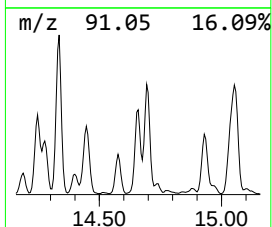
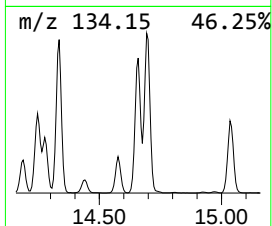
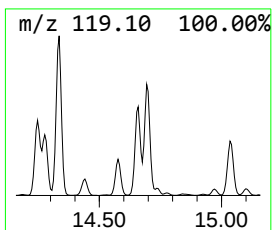
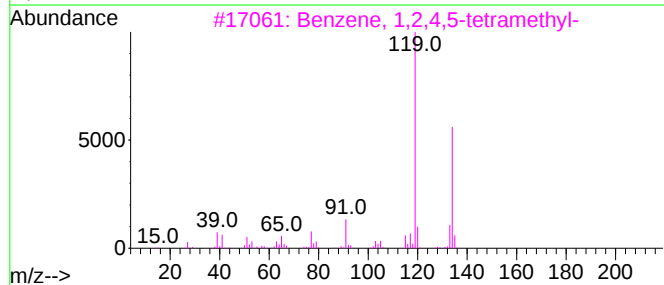
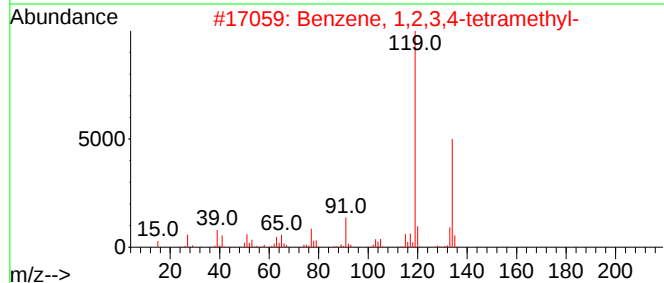
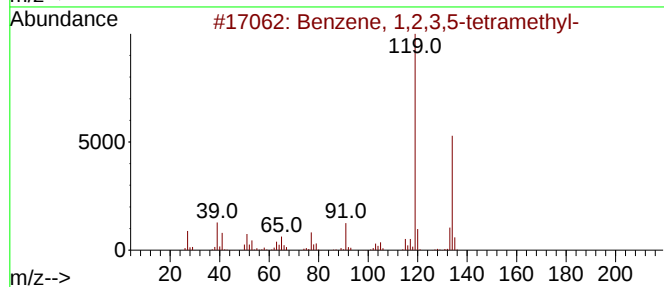
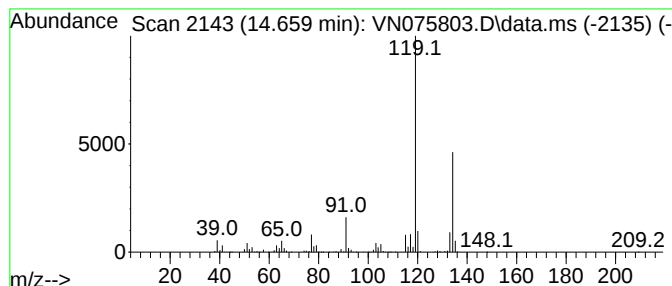
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.659	8.64 ug/l	5941940	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
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_N\Data\VN121422\
Data File : VN075803.D
Acq On : 15 Dec 2022 06:27
Operator : JC\MD
Sample : N5990-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP120822

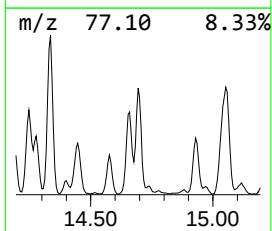
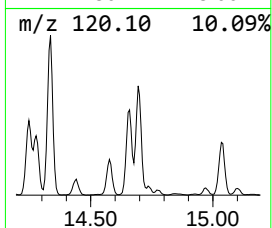
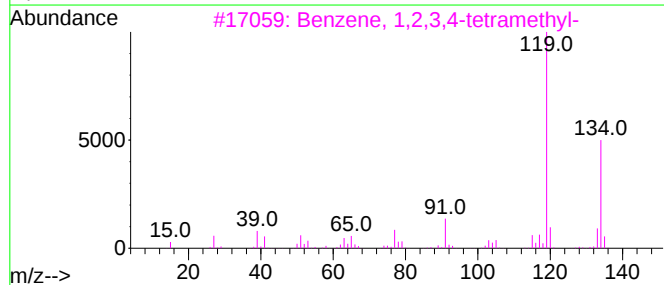
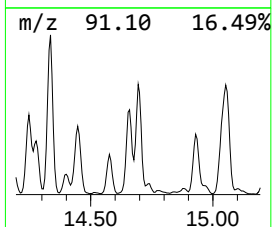
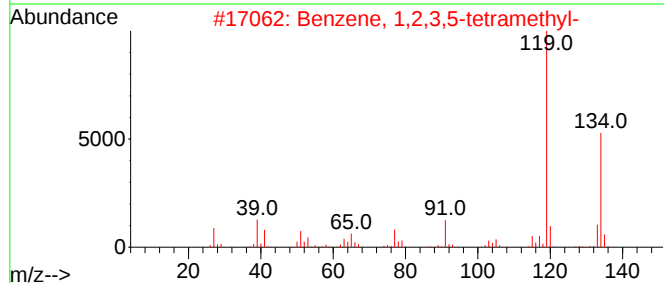
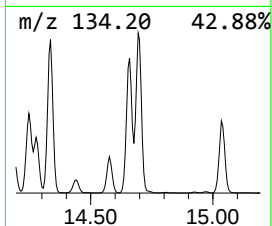
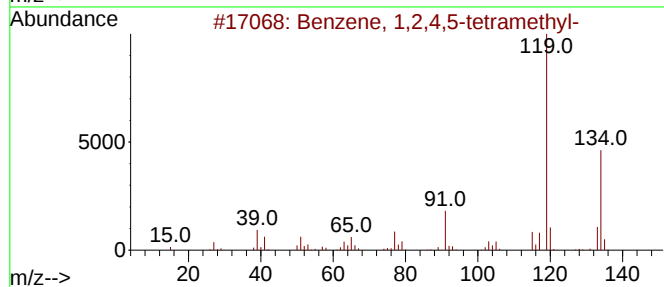
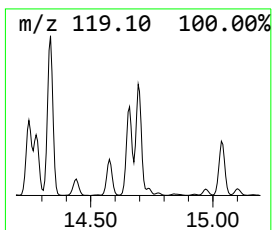
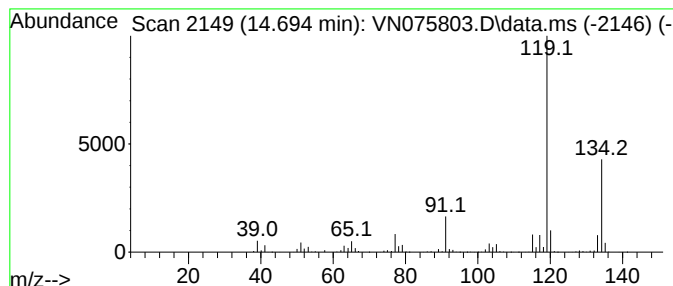
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.694	8.77 ug/l	6035130	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075803.D
Acq On : 15 Dec 2022 06:27
Operator : JC\MD
Sample : N5990-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP120822

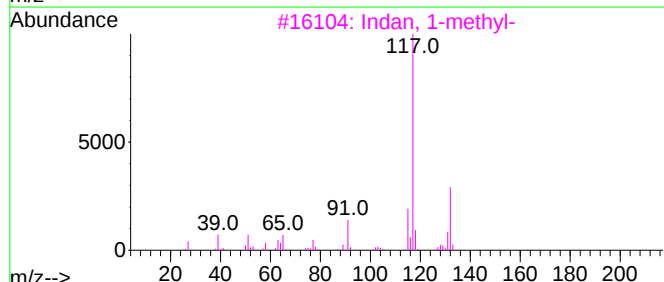
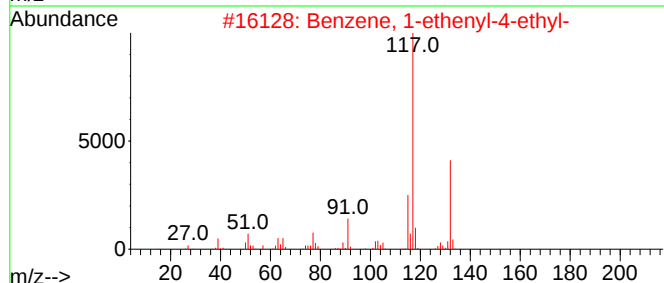
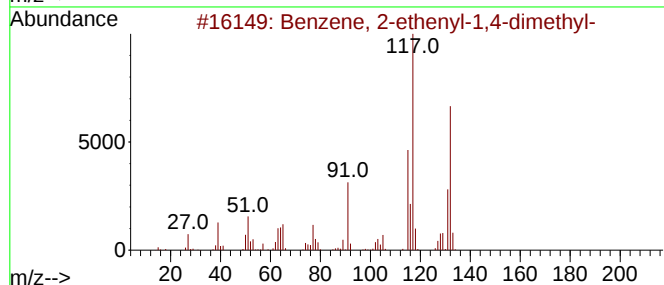
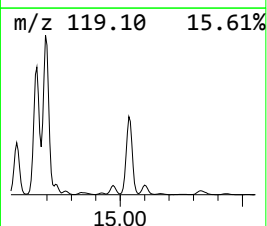
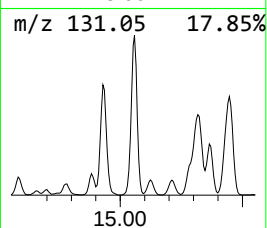
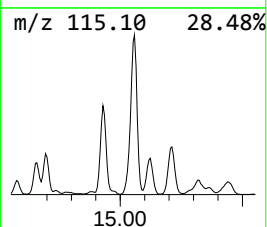
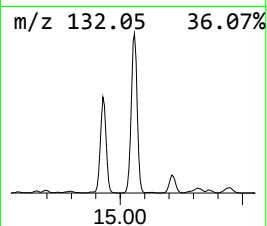
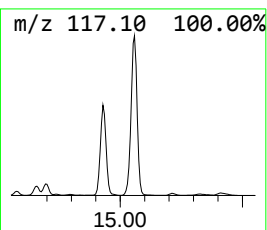
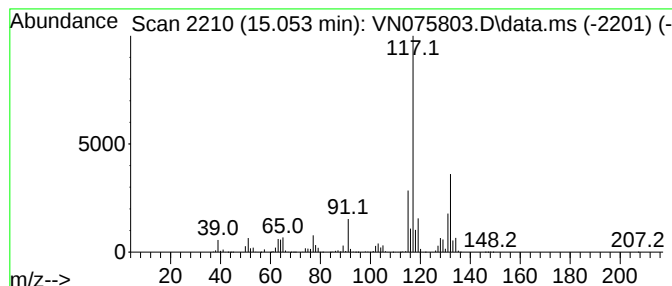
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	19.30 ug/l	13278000	1,4-Dichlorobenzene-d4	13.794

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	94
3		Indan, 1-methyl-	132	C10H12	000767-58-8	90
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121422\
Data File : VN075803.D
Acq On : 15 Dec 2022 06:27
Operator : JC\MD
Sample : N5990-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP120822

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Ethylidenecyclo...	8.741	26.2	ug/l	1151570	2	9.106	2197870	50.0
Cyclobutane, (1...	10.106	21.9	ug/l	960402	2	9.106	2197870	50.0
Benzene, 1-ethy...	13.100	31.6	ug/l	21774900	4	13.794	34405400	50.0
Benzene, 1-ethy...	13.341	41.6	ug/l	28610900	4	13.794	34405400	50.0
Indane	14.000	49.6	ug/l	34118700	4	13.794	34405400	50.0
Benzene, 4-ethy...	14.335	13.1	ug/l	8985650	4	13.794	34405400	50.0
Benzene, 1,2,3,...	14.659	8.6	ug/l	5941940	4	13.794	34405400	50.0
Benzene, 1,2,4,...	14.694	8.8	ug/l	6035130	4	13.794	34405400	50.0
Benzene, 2-ethe...	15.053	19.3	ug/l	13278000	4	13.794	34405400	50.0