

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
 Data File : VN083971.D
 Acq On : 18 Sep 2024 08:44
 Operator : JC\MD
 Sample : P3964-10
 Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CONVEYANA-2

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\82N091024W.M
 Title : SW846 8260

Signal : TIC: VN083971.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.159	1046	1055	1059	rBV	141129	323229	2.59%	0.374%
2	8.218	1059	1065	1075	rVV2	352827	795658	6.39%	0.921%
3	8.436	1093	1102	1113	rBV2	68698	158271	1.27%	0.183%
4	8.577	1117	1126	1138	rBV	163153	371675	2.98%	0.430%
5	8.753	1138	1156	1168	rBV	469950	1169848	9.39%	1.355%
6	8.959	1182	1191	1203	rBV2	59829	132204	1.06%	0.153%
7	9.094	1206	1214	1222	rBV	414584	849716	6.82%	0.984%
8	9.583	1285	1297	1302	rBV	173523	366060	2.94%	0.424%
9	9.641	1302	1307	1314	rVB	121137	234866	1.89%	0.272%
10	10.012	1362	1370	1381	rBV	352306	706786	5.67%	0.818%
11	10.141	1381	1392	1398	rBV	464513	945172	7.59%	1.094%
12	10.200	1398	1402	1406	rBV2	81764	127696	1.03%	0.148%
13	10.335	1416	1425	1438	rBV3	115869	286853	2.30%	0.332%
14	10.488	1444	1451	1458	rBV	411178	754281	6.06%	0.873%
15	10.565	1458	1464	1474	rVB	600258	1116485	8.96%	1.293%
16	10.741	1487	1494	1501	rVB	128179	234845	1.89%	0.272%
17	11.012	1531	1540	1547	rVB2	83969	167144	1.34%	0.194%
18	11.271	1574	1584	1593	rVV	79648	175831	1.41%	0.204%
19	11.641	1639	1647	1659	rVB2	183003	400020	3.21%	0.463%
20	11.741	1659	1664	1674	rVB	142929	245569	1.97%	0.284%
21	11.865	1674	1685	1691	rBV2	584565	1249301	10.03%	1.447%
22	11.959	1698	1701	1707	rVB	126300	206042	1.65%	0.239%
23	12.065	1711	1719	1725	rBV2	805506	1565107	12.57%	1.812%
24	12.394	1766	1775	1782	rVV	364089	738658	5.93%	0.855%
25	12.476	1784	1789	1795	rVB	94907	173353	1.39%	0.201%
26	12.694	1820	1826	1831	rBV	409498	723682	5.81%	0.838%
27	12.824	1840	1848	1851	rBV	522750	1034336	8.30%	1.198%
28	12.929	1861	1866	1871	rVB3	87261	154879	1.24%	0.179%
29	12.988	1871	1876	1879	rBV	193718	364317	2.92%	0.422%
30	13.035	1879	1884	1889	rVV	991869	1606495	12.90%	1.860%
31	13.094	1889	1894	1898	rVV	3591422	6149513	49.37%	7.120%
32	13.129	1898	1900	1903	rVV	1542427	2090196	16.78%	2.420%
33	13.171	1903	1907	1919	rVB	2341815	4098130	32.90%	4.745%
34	13.335	1926	1935	1942	rBV	1574317	2778115	22.30%	3.217%
35	13.423	1942	1950	1954	rBV	681488	1205465	9.68%	1.396%
36	13.482	1954	1960	1966	rBV	7785776	12456019	100.00%	14.422%

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Integration Parameters: RTEINT.P

Integrator: RTE

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Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Title : SW846 8260

37	13.529	1966	1968	1973	rVB	308269	416790	3.35%	0.483%
38	13.641	1974	1987	1996	rBV4	836566	2677099	21.49%	3.100%
39	13.712	1996	1999	2002	rBV	124220	173706	1.39%	0.201%
40	13.753	2002	2006	2008	rBV2	126682	157711	1.27%	0.183%
41	13.823	2008	2018	2030	rVB3	2474977	7186275	57.69%	8.321%
42	13.953	2034	2040	2041	rBV2	545542	828450	6.65%	0.959%
43	13.988	2041	2046	2049	rBV3	1710989	3447635	27.68%	3.992%
44	14.106	2062	2066	2074	rBV2	246906	423797	3.40%	0.491%
45	14.182	2074	2079	2084	rVB	596084	1037836	8.33%	1.202%
46	14.241	2084	2089	2092	rBV	965892	1660954	13.33%	1.923%
47	14.270	2092	2094	2099	rVV	893297	1185160	9.51%	1.372%
48	14.329	2099	2104	2112	rVB	1748895	2878643	23.11%	3.333%
49	14.441	2117	2123	2128	rVB2	443973	750236	6.02%	0.869%
50	14.570	2141	2145	2151	rVB	375332	603437	4.84%	0.699%
51	14.653	2151	2159	2162	rBV	1172222	2049793	16.46%	2.373%
52	14.694	2162	2166	2169	rVV	1405638	2088792	16.77%	2.419%
53	14.735	2169	2173	2177	rVB2	594874	868948	6.98%	1.006%
54	14.806	2182	2185	2192	rVB	291055	424933	3.41%	0.492%
55	14.876	2192	2197	2200	rBV	239207	329118	2.64%	0.381%
56	14.923	2200	2205	2209	rBV2	632265	1185293	9.52%	1.372%
57	15.029	2216	2223	2224	rBV4	930338	1707772	13.71%	1.977%
58	15.212	2250	2254	2260	rVB6	133618	259749	2.09%	0.301%
59	15.317	2261	2272	2284	rVV2	445092	1291161	10.37%	1.495%
60	15.429	2286	2291	2300	rVB4	279103	689790	5.54%	0.799%
61	15.635	2314	2326	2333	rVB	971774	2186685	17.56%	2.532%
62	15.747	2339	2345	2349	rBV4	148271	317378	2.55%	0.367%
63	15.823	2350	2358	2370	rVB2	296962	990399	7.95%	1.147%
64	15.941	2370	2378	2385	rBV	232730	582912	4.68%	0.675%
65	16.253	2426	2431	2437	rBV4	113601	247385	1.99%	0.286%
66	16.335	2439	2445	2452	rVB3	185575	443294	3.56%	0.513%
67	16.782	2513	2521	2522	rBV	645176	1118263	8.98%	1.295%

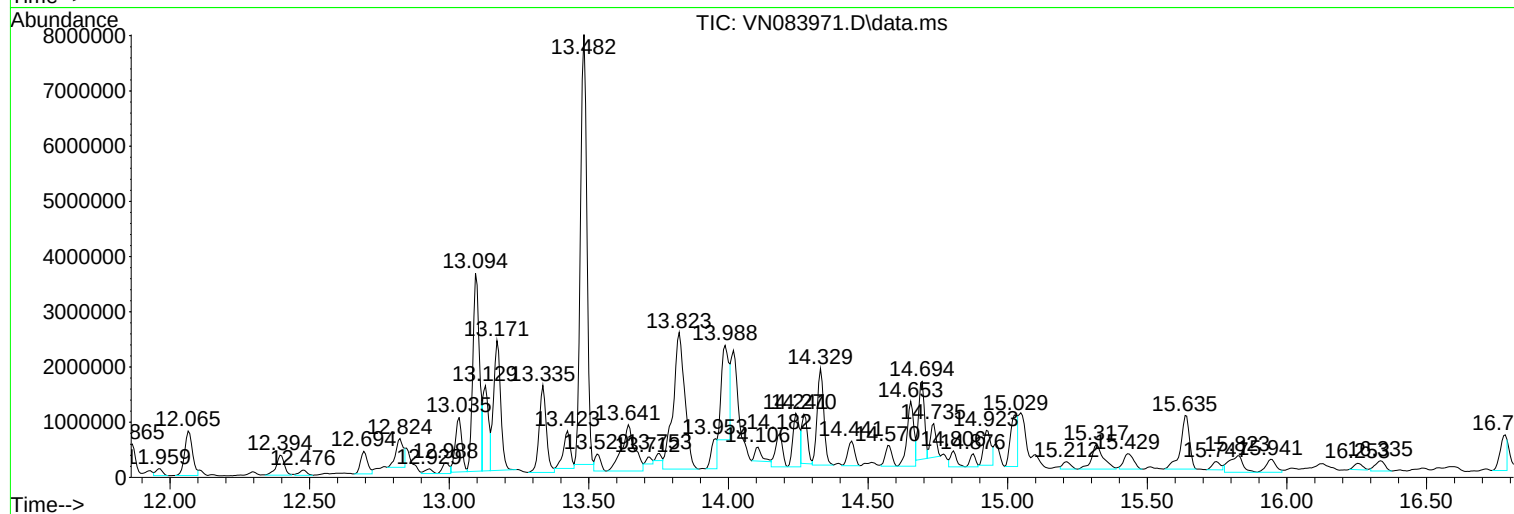
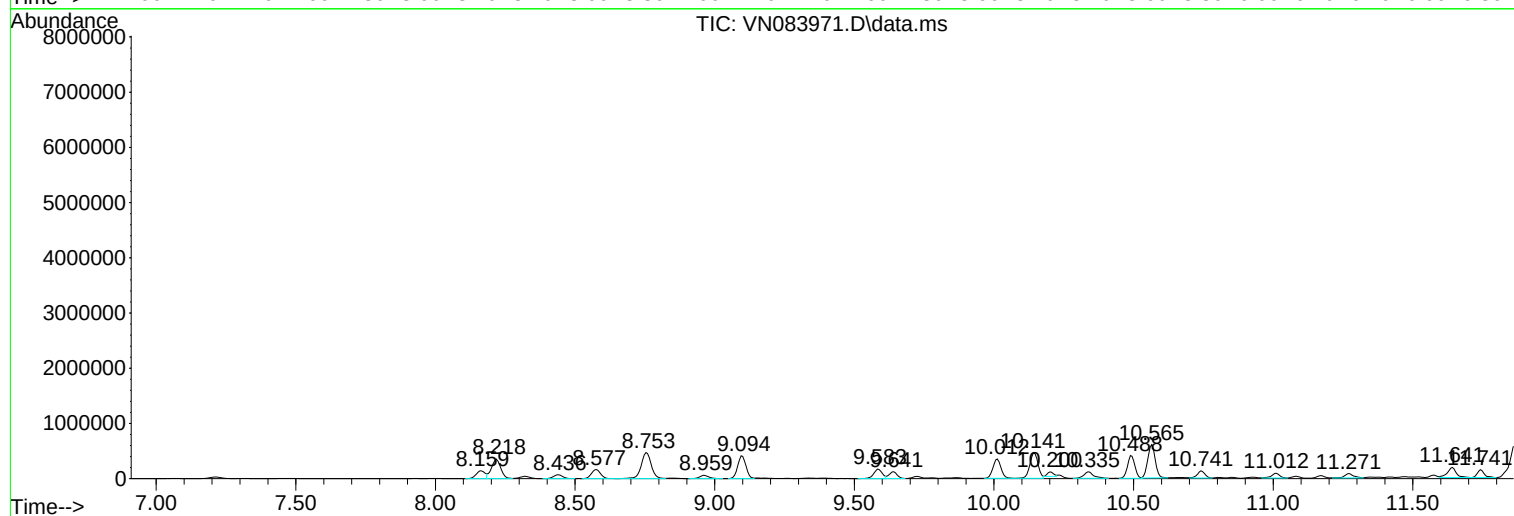
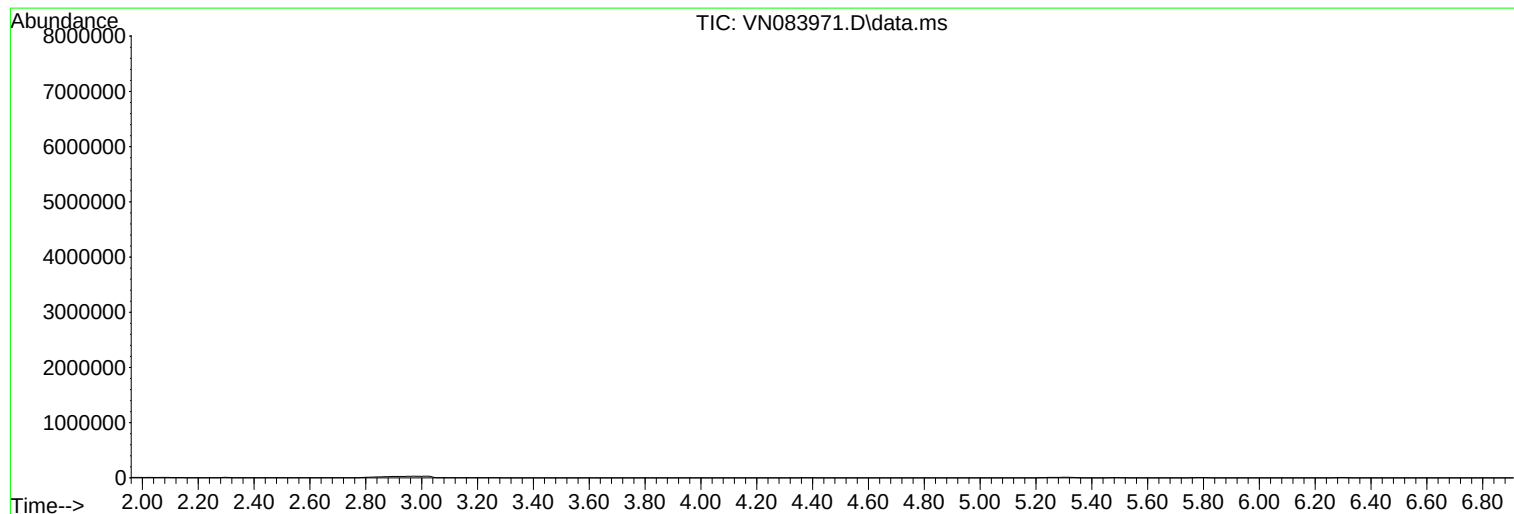
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Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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



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Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

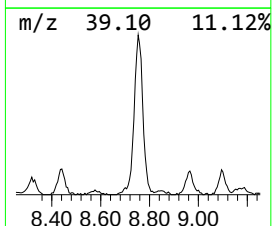
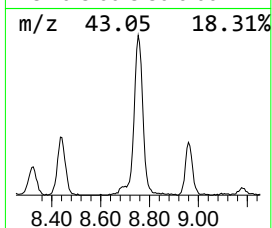
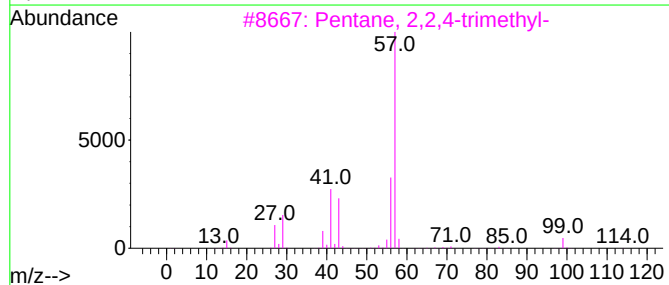
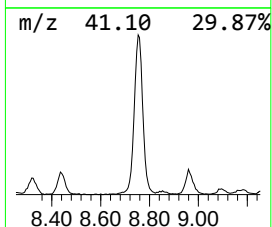
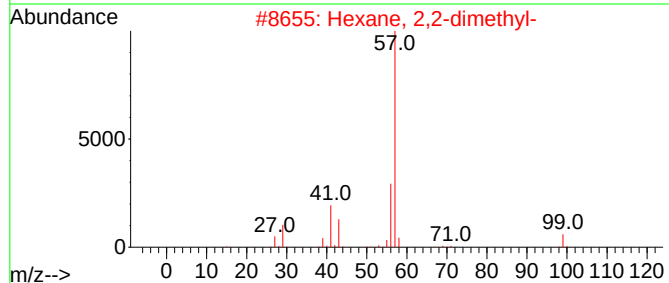
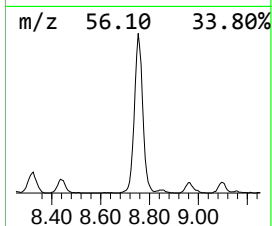
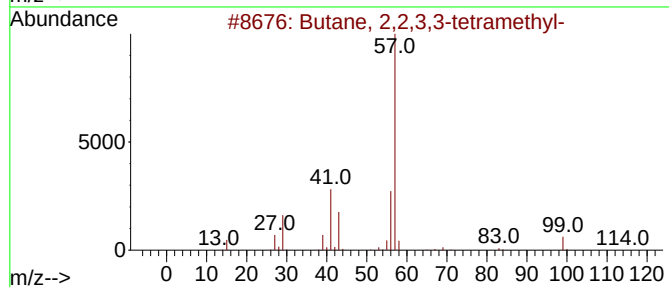
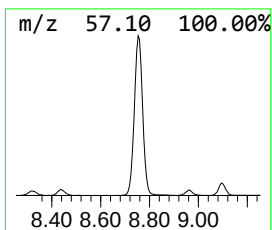
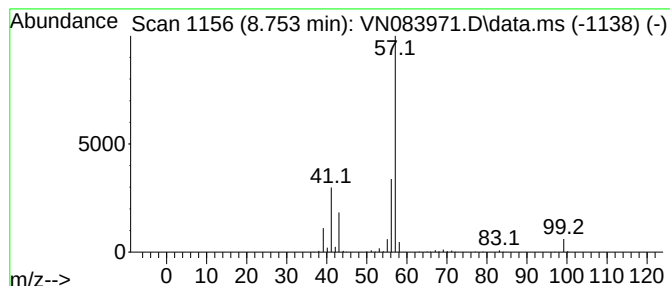
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.753	68.84 ug/l	1169850	1,4-Difluorobenzene	9.100

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
3			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	72
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	64



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ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
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CONVEYANA-2

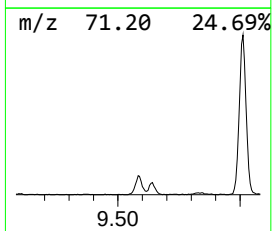
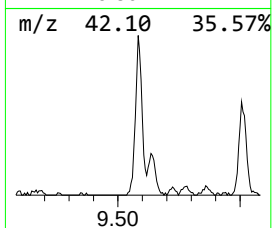
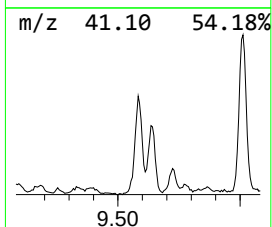
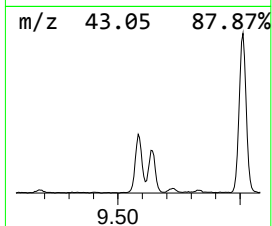
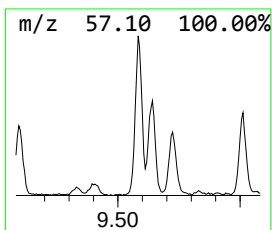
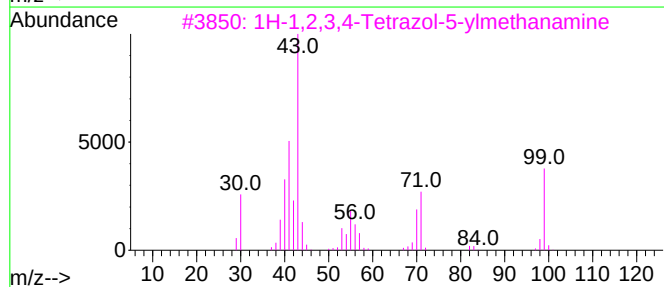
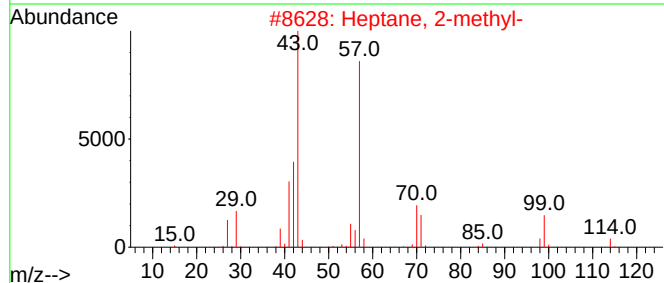
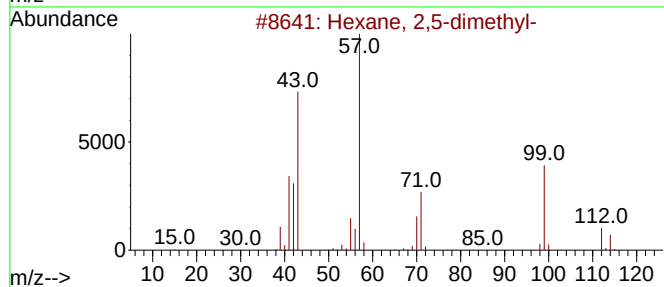
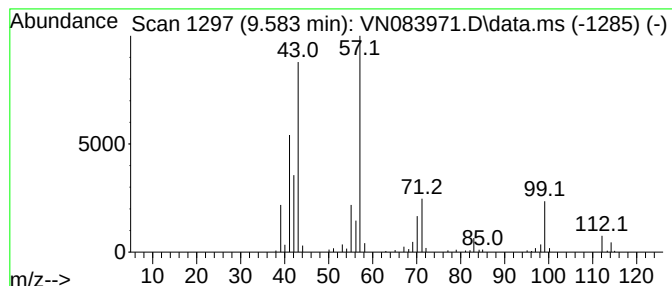
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 2 Hexane, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.583	21.54 ug/l	366060	1,4-Difluorobenzene	9.100

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	94
2		Heptane, 2-methyl-	114	C8H18	000592-27-8	72
3		1H-1,2,3,4-Tetrazol-5-ylmethanamine	99	C2H5N5	031602-63-8	64
4		Octane, 2,7-dimethyl-	142	C10H22	001072-16-8	50
5		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	47



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Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

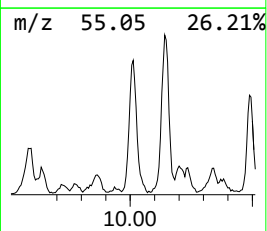
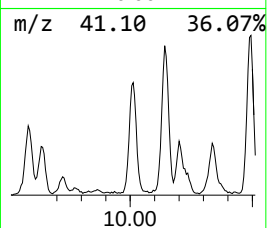
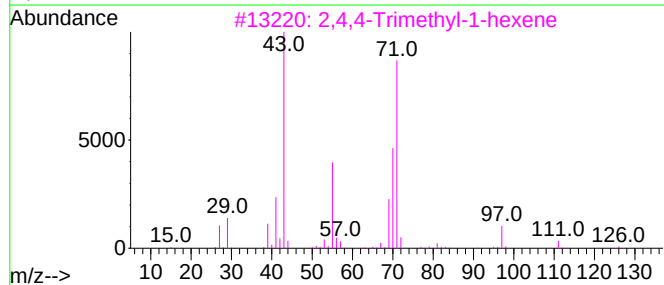
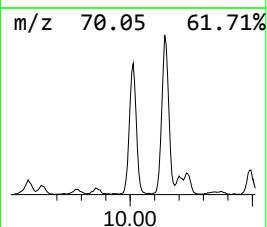
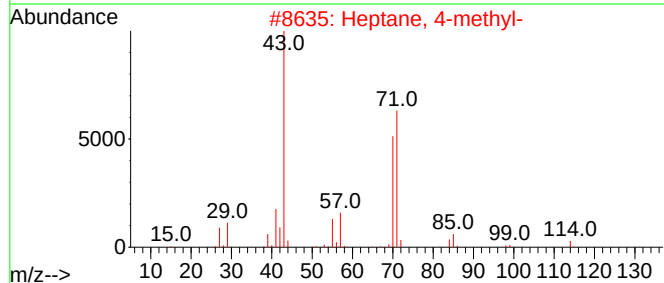
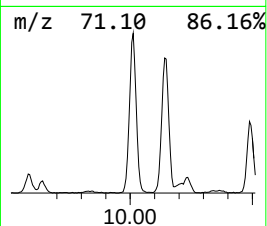
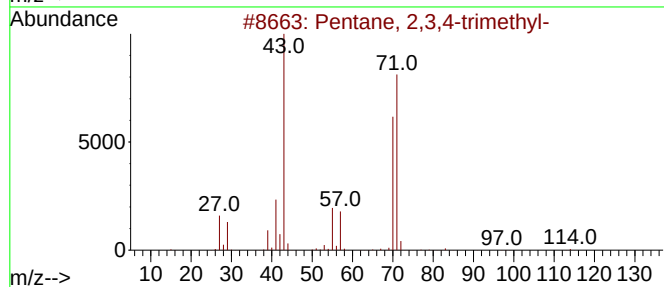
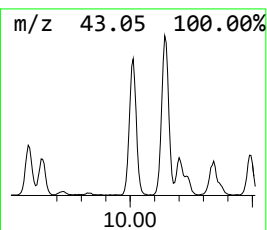
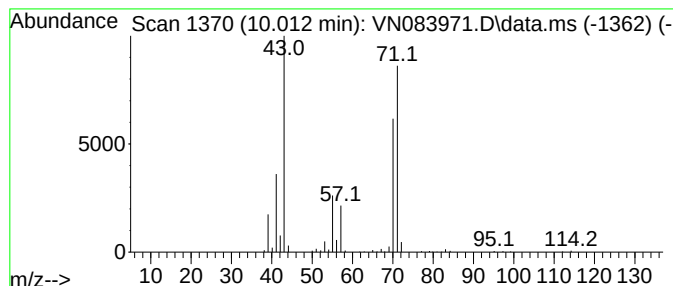
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.012	41.59 ug/l	706786	1,4-Difluorobenzene	9.100

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
4			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64



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ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

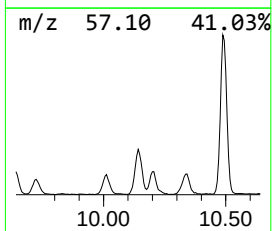
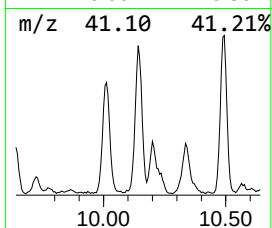
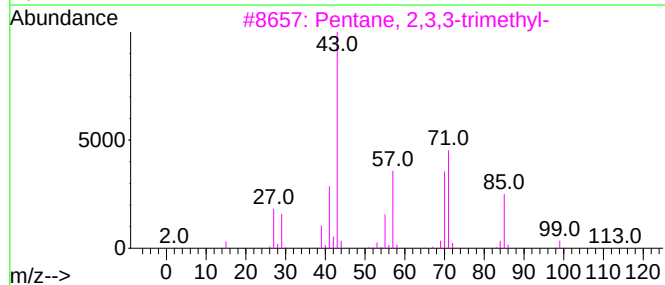
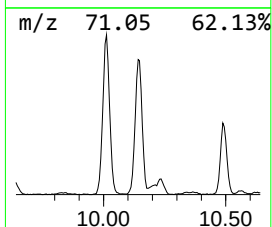
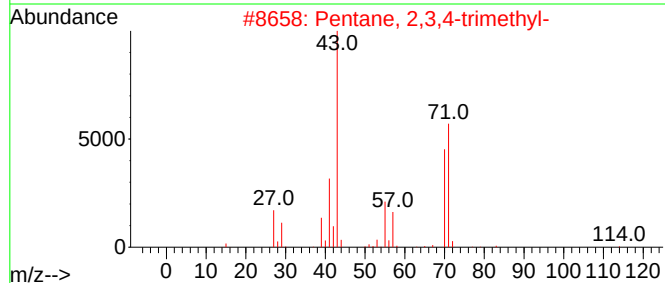
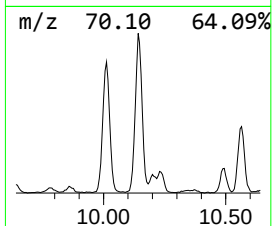
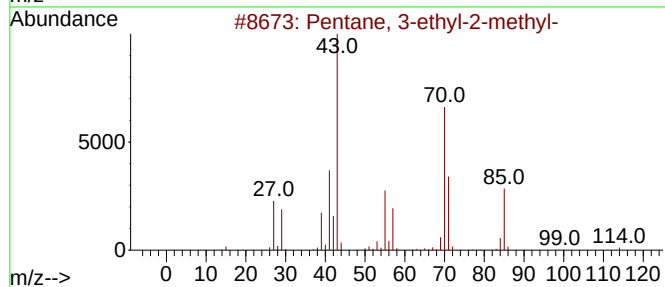
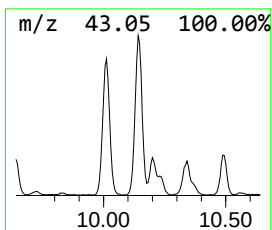
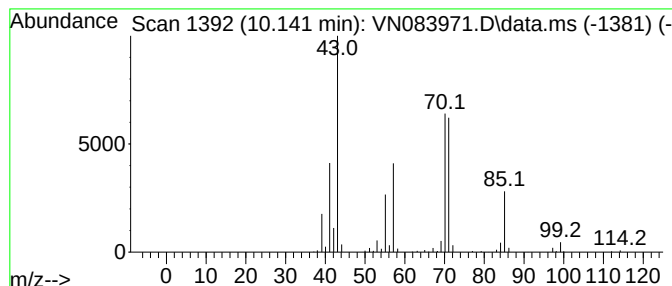
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.141	55.62 ug/l	945172	1,4-Difluorobenzene	9.100

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	81
2			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	81
3			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	64
4			Heptane, 3-ethyl-4-methyl-	142	C10H22	052896-91-0	59
5			Diisooamyl ether	158	C10H22O	000544-01-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

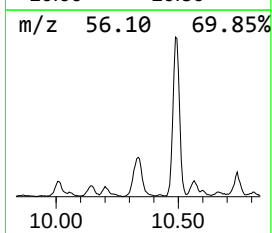
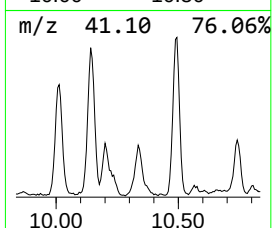
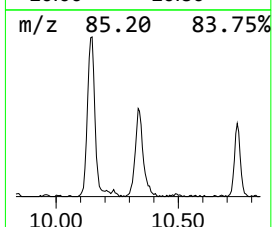
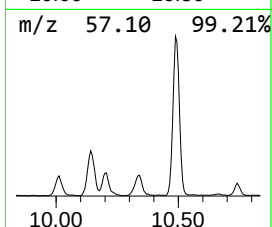
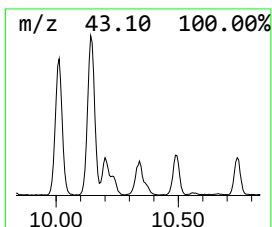
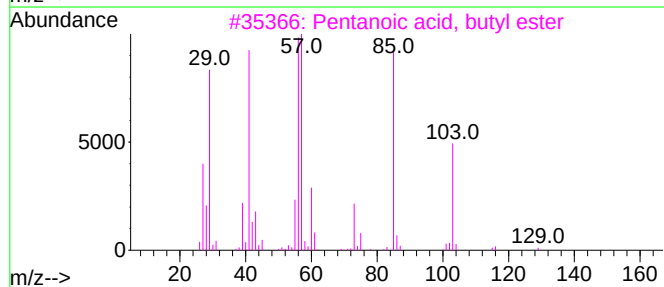
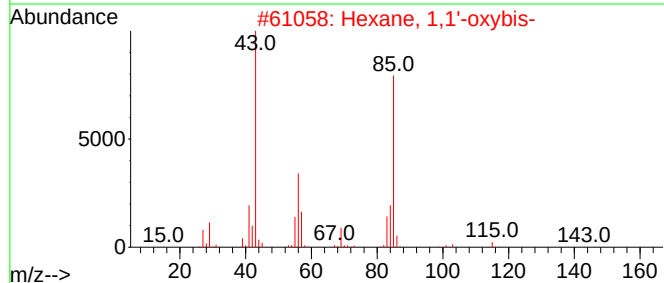
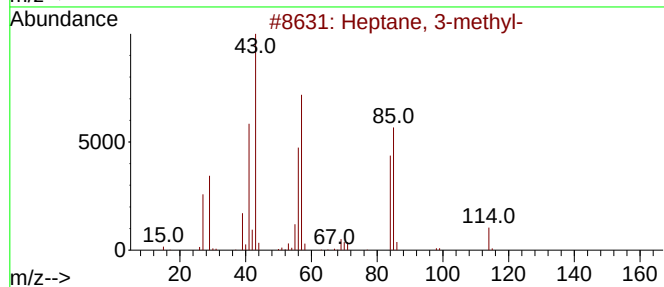
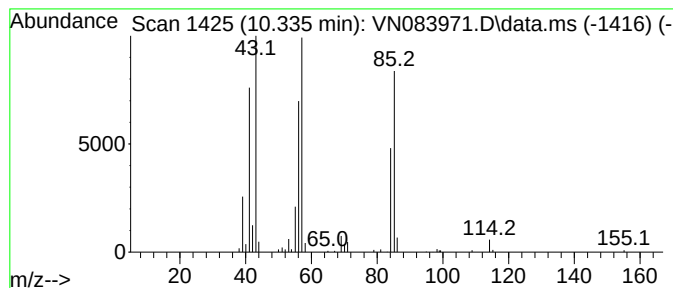
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Quant Title : SW846 8260

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

Peak Number 5 Heptane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.335	16.88 ug/l	286853	1,4-Difluorobenzene	9.100

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	90
2			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	59
3			Pentanoic acid, butyl ester	158	C9H18O2	000591-68-4	53
4			Butanoic acid, 3-methyl-, butyl ...	158	C9H18O2	000109-19-3	53
5			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

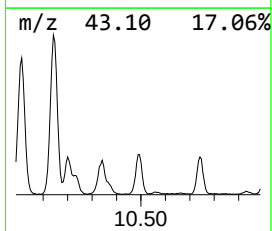
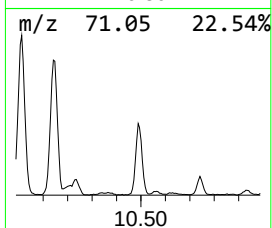
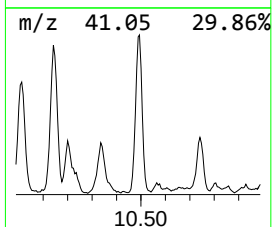
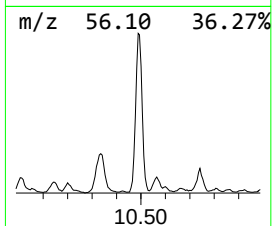
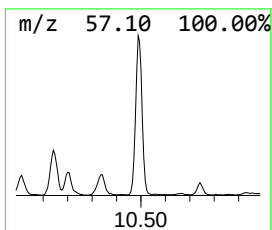
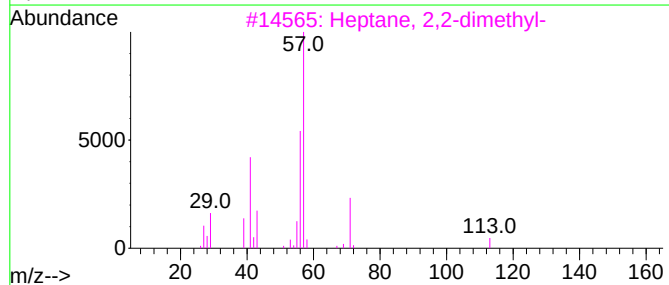
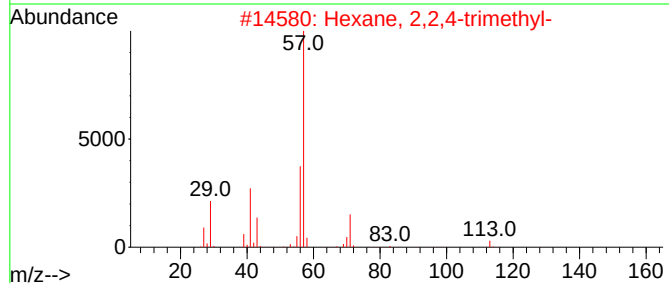
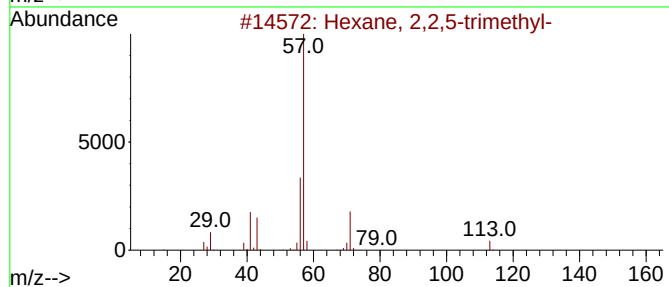
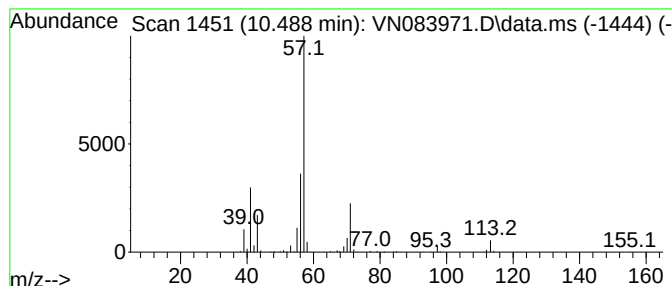
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 2,2,5-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.488	30.19 ug/l	754281	Chlorobenzene-d5	11.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	78
2			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	78
3			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64
4			Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	64
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

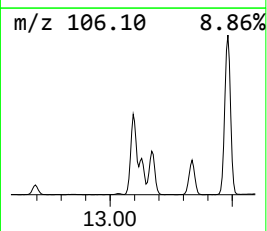
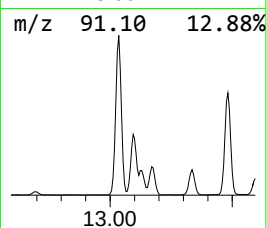
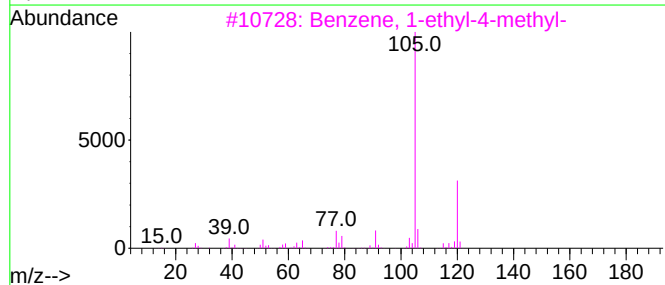
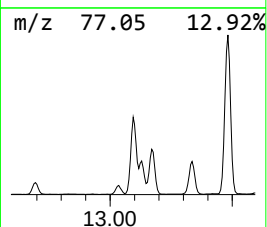
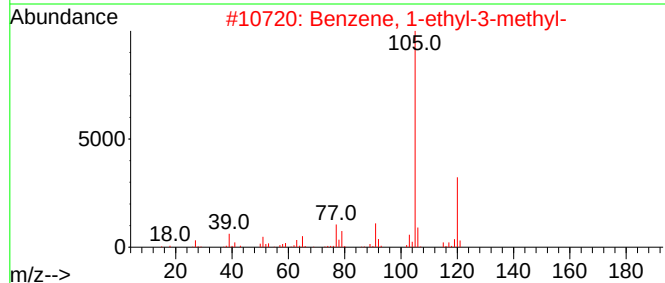
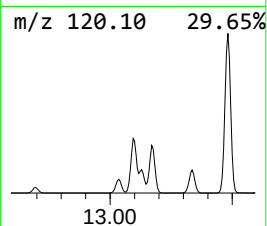
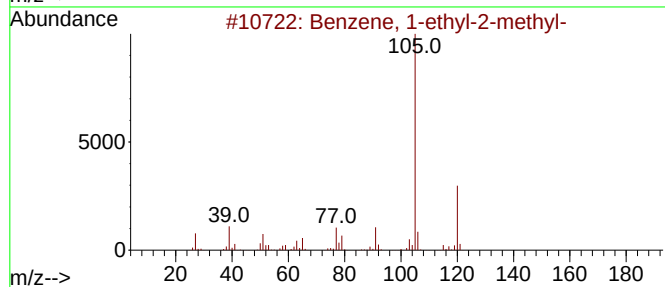
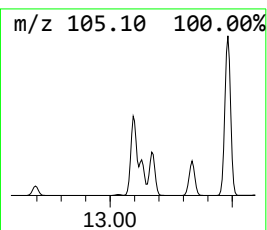
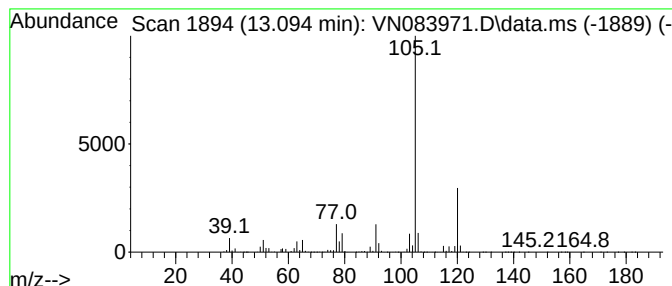
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.094	42.79 ug/l	6149510	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	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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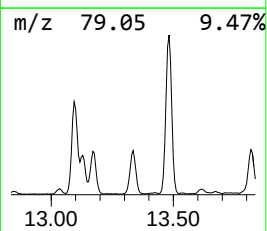
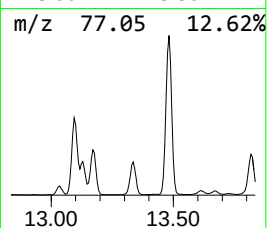
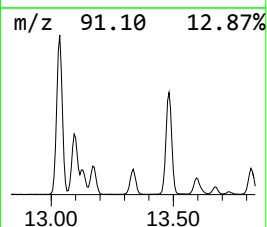
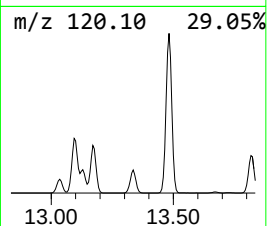
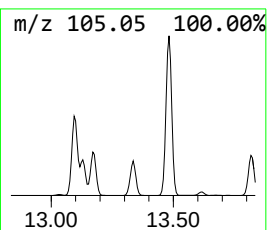
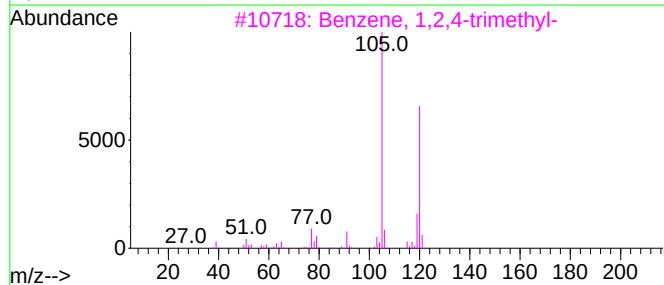
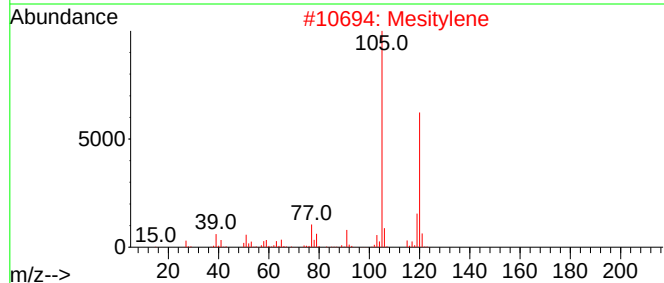
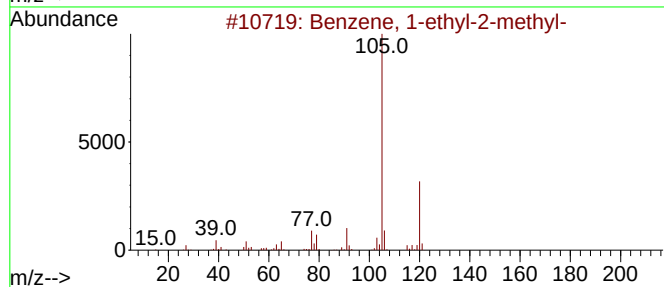
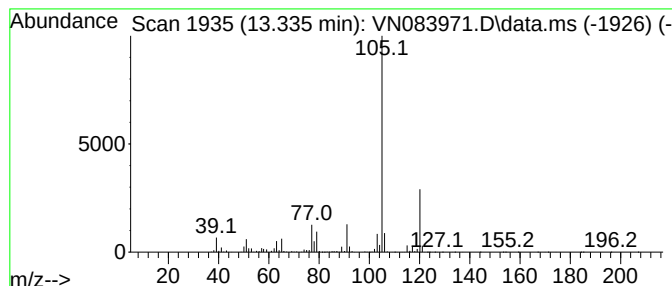
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.335	19.33 ug/l	2778120	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			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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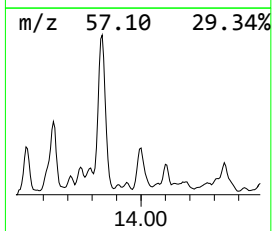
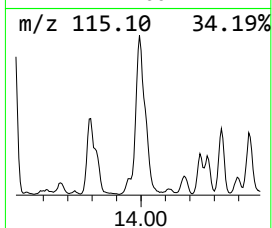
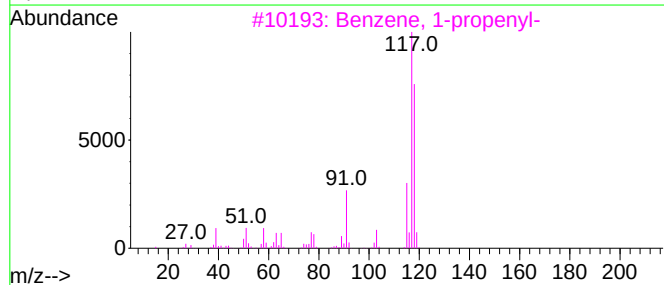
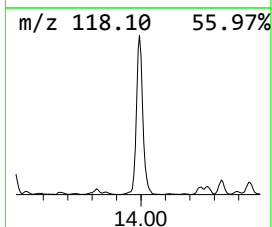
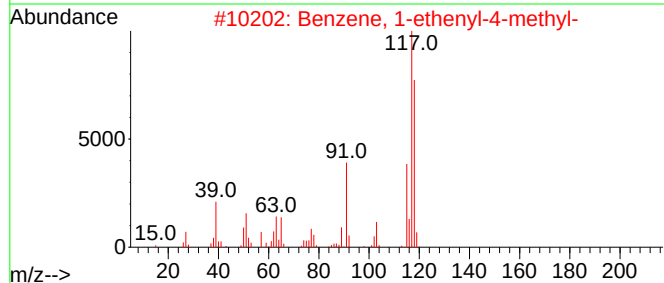
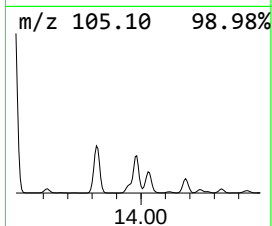
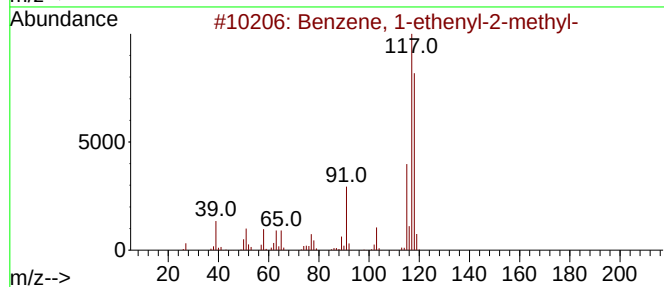
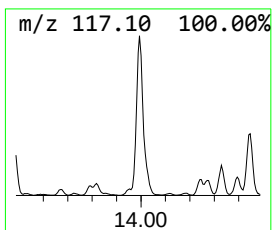
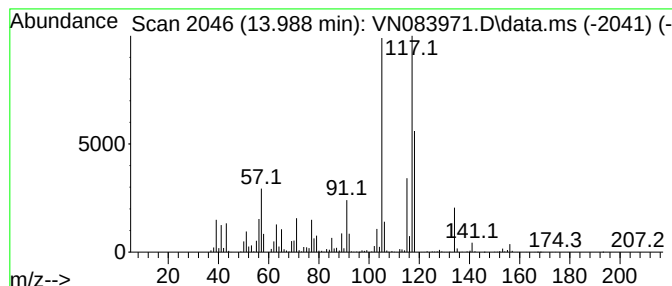
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.988	23.99 ug/l	3447640	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
2			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

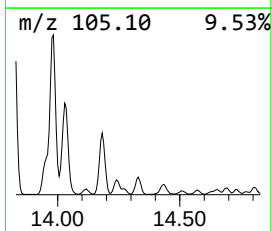
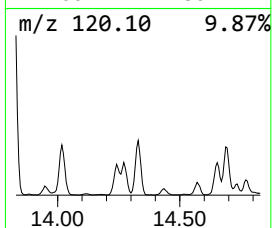
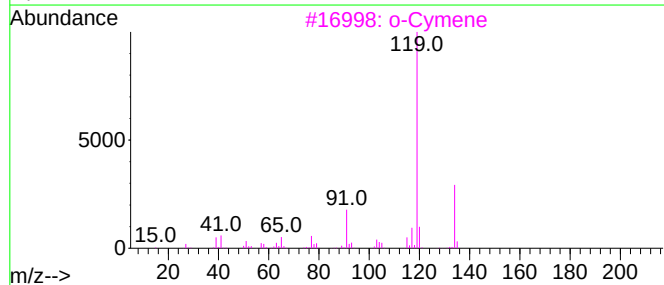
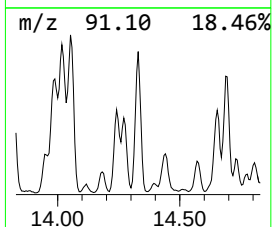
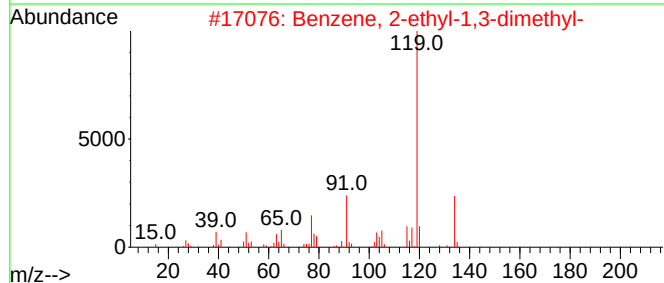
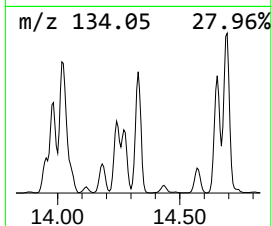
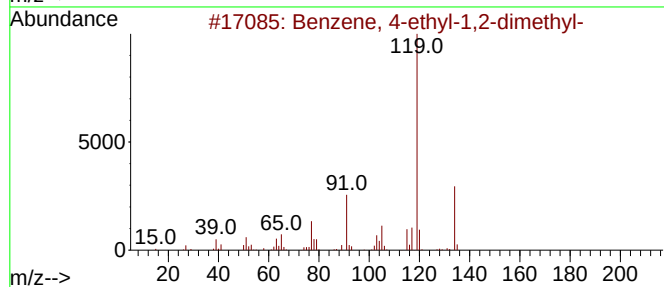
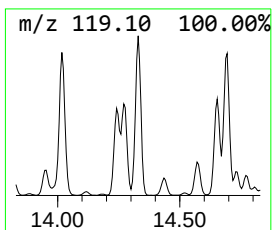
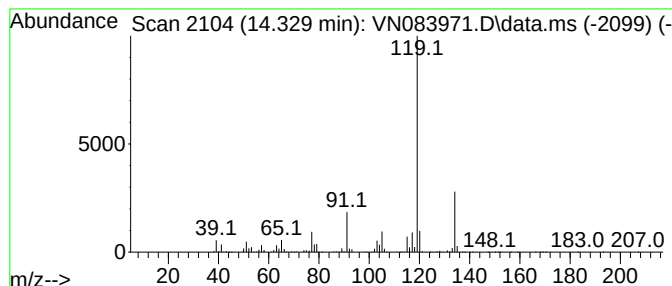
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.329	20.03 ug/l	2878640	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	96
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
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-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083971.D
Acq On : 18 Sep 2024 08:44
Operator : JC\MD
Sample : P3964-10
Misc : 3.26g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CONVEYANA-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.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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Hexane, 2,5-dim...	9.583	21.5	ug/l	366060	2	9.100	849716	50.0
Pentane, 2,3,4-...	10.012	41.6	ug/l	706786	2	9.100	849716	50.0
Pentane, 3-ethy...	10.141	55.6	ug/l	945172	2	9.100	849716	50.0
Heptane, 3-methyl-	10.335	16.9	ug/l	286853	2	9.100	849716	50.0
Hexane, 2,2,5-t...	10.488	30.2	ug/l	754281	3	11.865	1249300	50.0
Benzene, 1-ethy...	13.094	42.8	ug/l	6149510	4	13.794	7186280	50.0
Benzene, 1-ethy...	13.335	19.3	ug/l	2778120	4	13.794	7186280	50.0
Benzene, 1-ethe...	13.988	24.0	ug/l	3447640	4	13.794	7186280	50.0
Benzene, 4-ethy...	14.329	20.0	ug/l	2878640	4	13.794	7186280	50.0