

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
 Data File : VN088135.D
 Acq On : 28 Oct 2025 13:38
 Operator : JC\MD
 Sample : Q3468-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-10

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\82N102325W.M
 Title : SW846 8260

Signal : TIC: VN088135.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.342	59	66	71	rBV	35151	60585	3.39%	0.254%
2	2.536	89	99	108	rBV2	91128	169658	9.48%	0.712%
3	2.636	108	116	122	rVV2	25841	52032	2.91%	0.218%
4	3.265	212	223	240	rBV	463875	1120797	62.65%	4.701%
5	3.659	283	290	297	rBV5	12846	33288	1.86%	0.140%
6	4.418	406	419	436	rVB2	174249	627911	35.10%	2.634%
7	4.730	461	472	481	rVB3	8457	26972	1.51%	0.113%
8	5.265	545	563	578	rBV2	372518	1602438	89.58%	6.721%
9	5.800	638	654	667	rBV	120437	423315	23.66%	1.775%
10	7.053	857	867	875	rBV3	20417	61326	3.43%	0.257%
11	7.206	882	893	903	rBV2	94000	278251	15.55%	1.167%
12	7.312	903	911	919	rVV2	77083	211019	11.80%	0.885%
13	7.371	919	921	928	rVB3	14868	28525	1.59%	0.120%
14	7.477	937	939	953	rVB8	14147	35198	1.97%	0.148%
15	7.806	986	995	1004	rVB5	6964	19745	1.10%	0.083%
16	7.918	1007	1014	1019	rBV3	11228	29262	1.64%	0.123%
17	7.988	1019	1026	1035	rVB3	35817	100748	5.63%	0.423%
18	8.153	1042	1054	1058	rBV2	232656	571558	31.95%	2.397%
19	8.206	1058	1063	1074	rVV	371895	970845	54.27%	4.072%
20	8.312	1074	1081	1093	rVV2	238721	605399	33.84%	2.539%
21	8.424	1094	1100	1105	rVV4	9040	19899	1.11%	0.083%
22	8.494	1105	1112	1116	rVV3	36748	86140	4.82%	0.361%
23	8.559	1116	1123	1137	rVV	257716	628685	35.14%	2.637%
24	8.688	1137	1145	1149	rVV4	42321	118017	6.60%	0.495%
25	8.747	1149	1155	1164	rVV3	76849	261805	14.63%	1.098%
26	8.835	1164	1170	1180	rVB2	63212	148946	8.33%	0.625%
27	9.082	1202	1212	1224	rBV	672708	1420681	79.42%	5.959%
28	9.177	1224	1228	1235	rVB4	18503	36927	2.06%	0.155%
29	9.318	1245	1252	1256	rBV3	26821	54838	3.07%	0.230%
30	9.365	1256	1260	1268	rVB3	21303	41882	2.34%	0.176%
31	9.547	1281	1291	1295	rBV4	77680	204688	11.44%	0.859%
32	9.624	1302	1304	1313	rVB6	16635	33588	1.88%	0.141%
33	9.712	1313	1319	1322	rBV5	9020	18703	1.05%	0.078%
34	9.847	1332	1342	1348	rBV4	31990	73193	4.09%	0.307%
35	9.994	1360	1367	1376	rVB	57377	126897	7.09%	0.532%
36	10.088	1376	1383	1384	rBV3	25093	36471	2.04%	0.153%

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Integrator: RTE

Smoothing : ON

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Max Peaks: 100

Stop Thrs : 0

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

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37	10.124	1384	1389	1401	rVV	92800	215375	12.04%	0.903%
38	10.306	1413	1420	1427	rBV4	24666	54230	3.03%	0.227%
39	10.435	1435	1442	1445	rBV4	31819	59655	3.33%	0.250%
40	10.465	1445	1447	1453	rVB4	19331	32024	1.79%	0.134%
41	10.547	1453	1461	1478	rBV	918064	1788910	100.00%	7.503%
42	10.741	1481	1494	1500	rVV2	100802	224744	12.56%	0.943%
43	10.812	1500	1506	1514	rVV2	87445	161063	9.00%	0.676%
44	10.894	1514	1520	1524	rVV5	24976	49045	2.74%	0.206%
45	10.965	1524	1532	1541	rVV3	68528	189863	10.61%	0.796%
46	11.041	1541	1545	1555	rVB9	12914	33340	1.86%	0.140%
47	11.235	1570	1578	1582	rBV3	33246	68749	3.84%	0.288%
48	11.365	1595	1600	1604	rVB4	16450	27479	1.54%	0.115%
49	11.559	1627	1633	1642	rVB4	17897	34681	1.94%	0.145%
50	11.847	1673	1682	1695	rVV	910580	1731118	96.77%	7.261%
51	11.941	1695	1698	1709	rVV8	23358	56934	3.18%	0.239%
52	12.053	1709	1717	1723	rVV3	26430	68511	3.83%	0.287%
53	12.553	1793	1802	1804	rBV8	12372	24457	1.37%	0.103%
54	12.676	1817	1823	1830	rVB	82425	145048	8.11%	0.608%
55	12.829	1839	1849	1861	rBV	635496	1208846	67.57%	5.070%
56	12.906	1861	1862	1870	rVB6	14346	23984	1.34%	0.101%
57	13.017	1875	1881	1887	rVB2	41329	77301	4.32%	0.324%
58	13.153	1901	1904	1910	rVB7	11391	21061	1.18%	0.088%
59	13.312	1926	1931	1940	rBV2	40497	78965	4.41%	0.331%
60	13.576	1968	1976	1982	rBV2	67478	173959	9.72%	0.730%
61	13.770	2002	2009	2021	rBV	922569	1757859	98.26%	7.373%
62	13.859	2021	2024	2030	rVB	43381	68105	3.81%	0.286%
63	13.929	2030	2036	2038	rBV2	77625	129231	7.22%	0.542%
64	13.970	2038	2043	2048	rBV	659114	1105507	61.80%	4.637%
65	14.094	2059	2064	2070	rVB	82689	139146	7.78%	0.584%
66	14.164	2071	2076	2083	rVB3	18972	38240	2.14%	0.160%
67	14.247	2085	2090	2096	rBV2	45988	82653	4.62%	0.347%
68	14.306	2096	2100	2107	rVB2	40076	68253	3.82%	0.286%
69	14.370	2107	2111	2113	rBV3	22951	33661	1.88%	0.141%
70	14.423	2113	2120	2127	rVV	881113	1468088	82.07%	6.158%
71	14.482	2128	2130	2138	rVB8	22038	51214	2.86%	0.215%
72	14.559	2139	2143	2146	rVB	37995	51296	2.87%	0.215%
73	14.629	2147	2155	2161	rVB3	64513	125909	7.04%	0.528%
74	14.717	2165	2170	2172	rBV4	30536	52956	2.96%	0.222%
75	14.747	2172	2175	2179	rBV2	28909	43409	2.43%	0.182%
76	14.906	2196	2202	2208	rVB4	102936	195118	10.91%	0.818%

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Title : SW846 8260

77	15.023	2214	2222	2229	rBV2	174063	424167	23.71%	1.779%
78	15.094	2229	2234	2240	rVB4	79740	138047	7.72%	0.579%
79	15.182	2246	2249	2255	rBV8	11754	23628	1.32%	0.099%
80	15.288	2262	2267	2274	rBV5	49756	115727	6.47%	0.485%
81	15.400	2274	2286	2293	rVV4	119384	339821	19.00%	1.425%
82	15.470	2293	2298	2309	rVB9	47283	146608	8.20%	0.615%
83	15.576	2310	2316	2319	rBV4	35786	62435	3.49%	0.262%
84	15.664	2328	2331	2332	rBV3	16738	18338	1.03%	0.077%
85	16.017	2387	2391	2392	rBV4	14517	17891	1.00%	0.075%
86	16.223	2423	2426	2433	rVB3	37181	69747	3.90%	0.293%
87	16.311	2433	2441	2450	rVB5	42915	112659	6.30%	0.473%
88	16.682	2501	2504	2510	rVB2	42457	72811	4.07%	0.305%

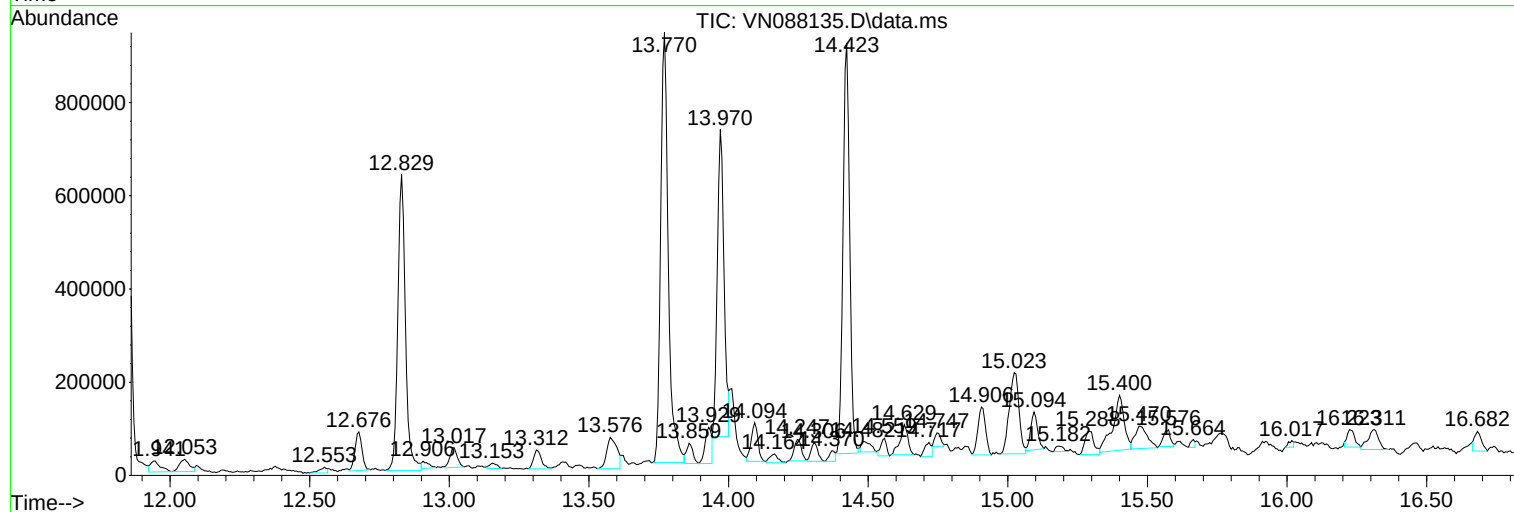
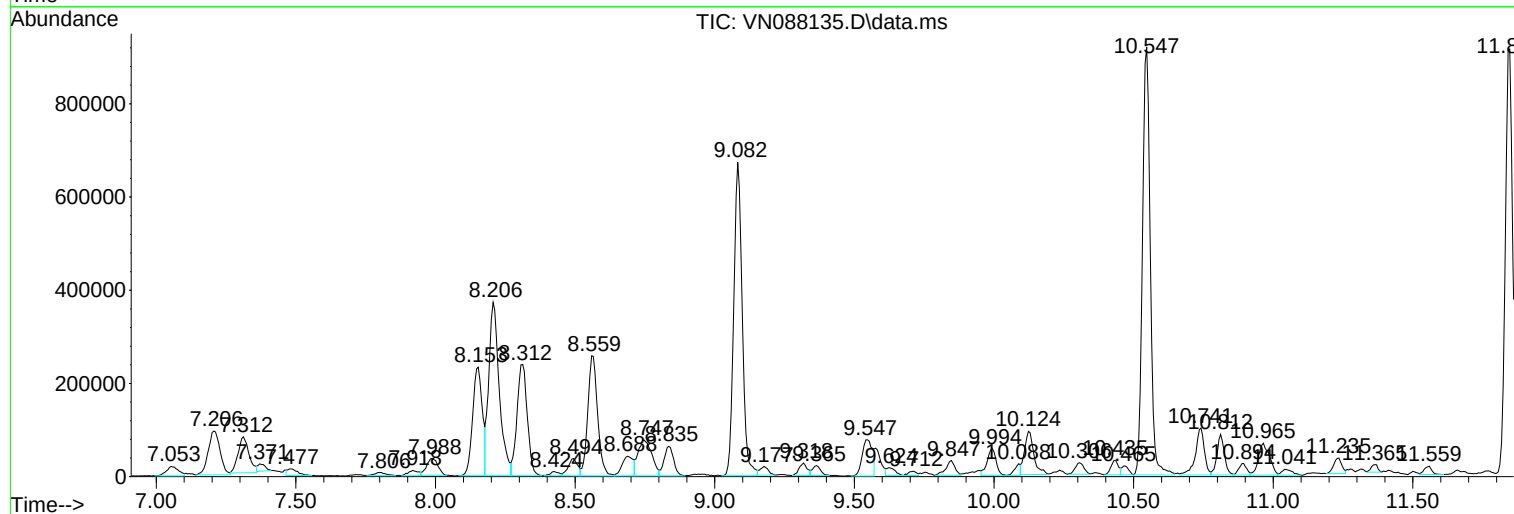
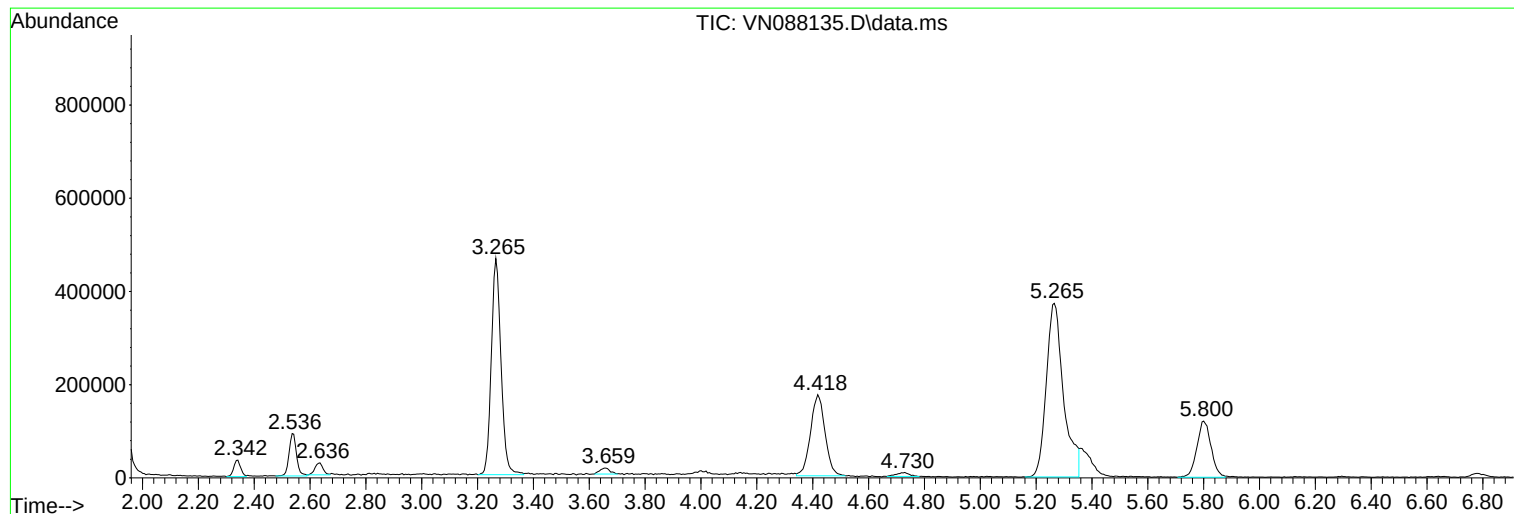
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

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



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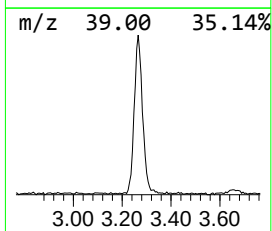
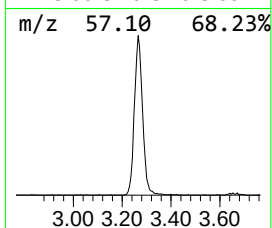
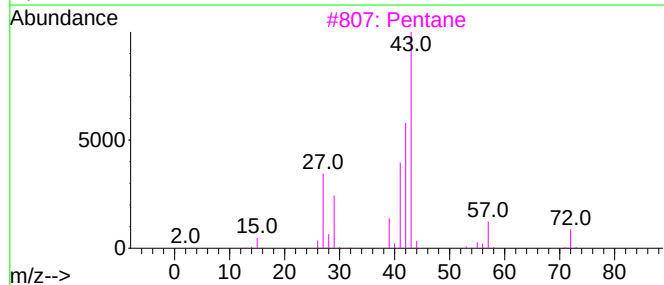
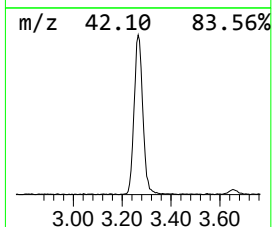
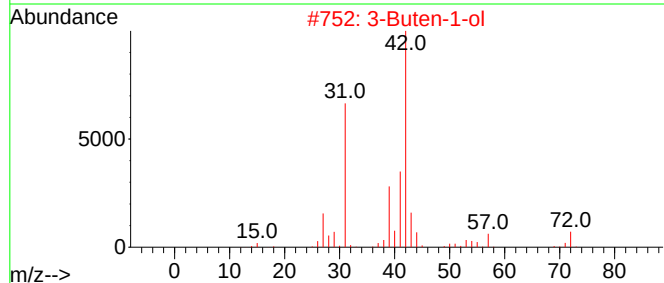
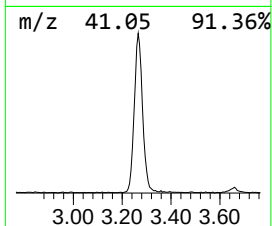
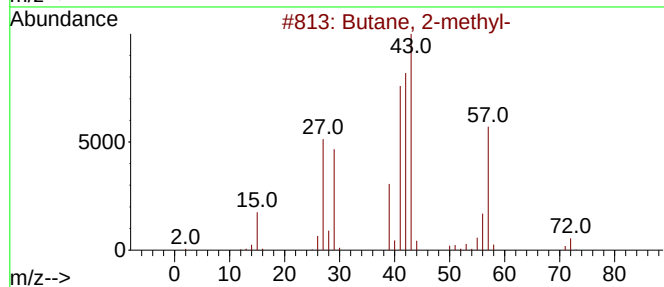
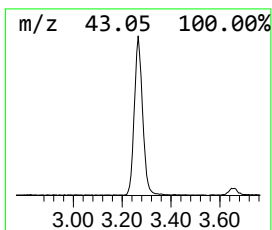
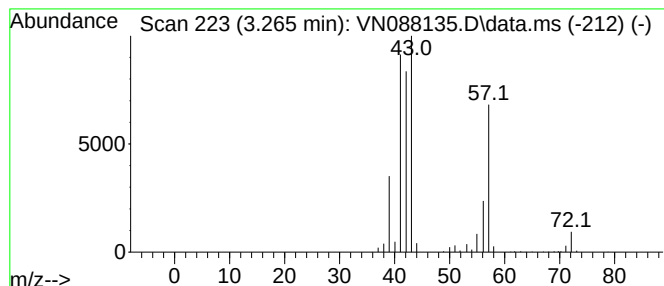
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.265	57.72 ug/l	1120800	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			3-Buten-1-ol	72	C4H8O	000627-27-0	47
3			Pentane	72	C5H12	000109-66-0	45
4			1-Propene, 2-methyl-	56	C4H8	000115-11-7	14
5			1-Butene	56	C4H8	000106-98-9	9



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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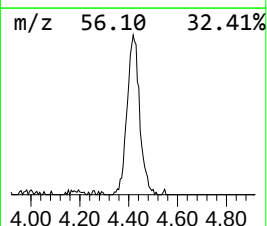
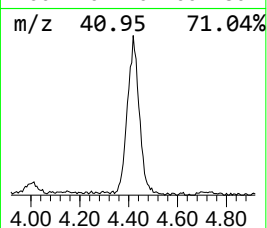
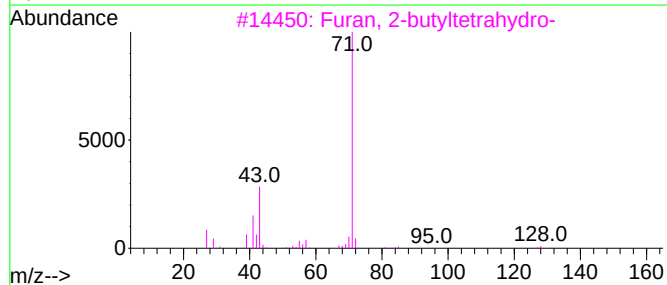
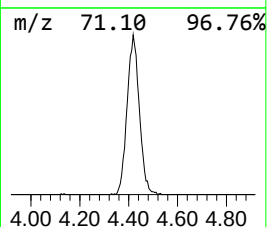
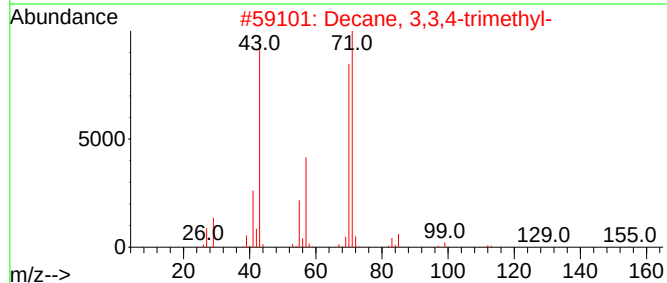
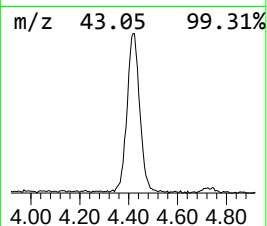
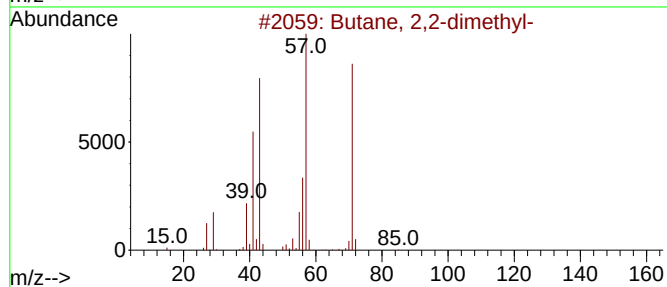
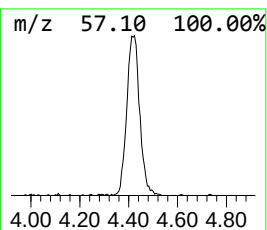
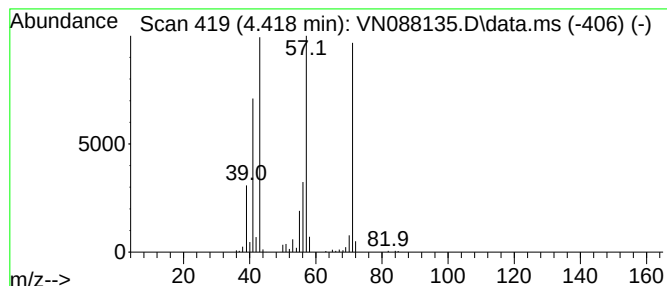
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.418	32.34 ug/l	627911	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	90
2			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	42
3			Furan, 2-butyltetrahydro-	128	C8H16O	001004-29-1	38
4			Pentane, 3-bromo-	150	C5H11Br	001809-10-5	36
5			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	36



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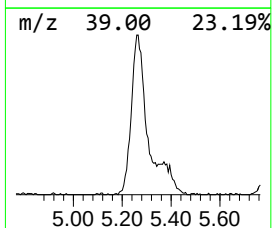
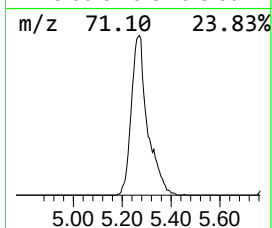
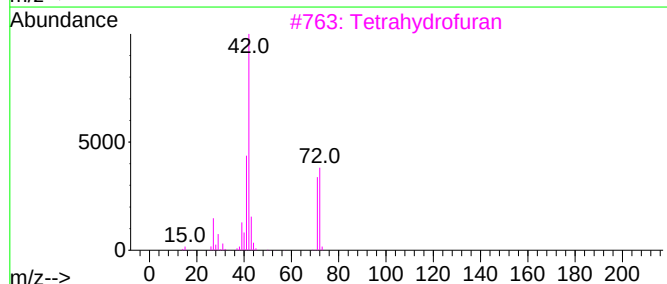
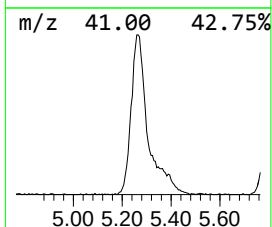
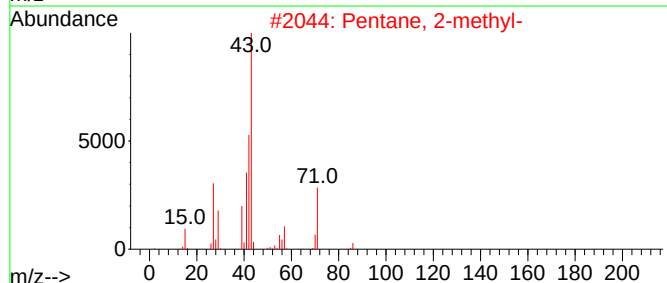
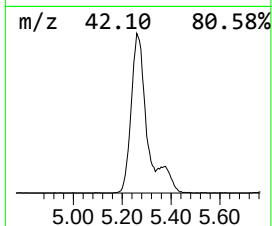
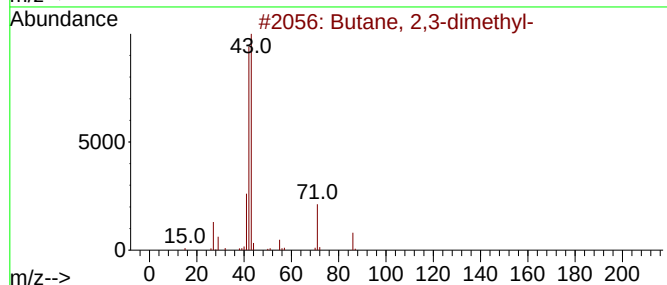
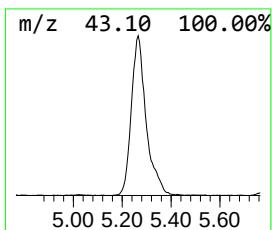
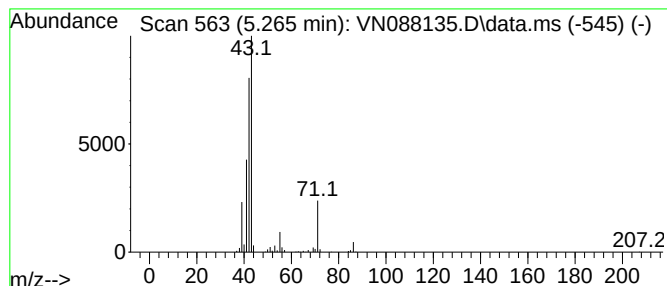
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.265	82.53 ug/l	1602440	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3			Tetrahydrofuran	72	C4H8O	000109-99-9	38
4			Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	12
5			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9



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ClientSampleId :
MW-10

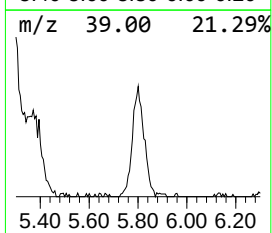
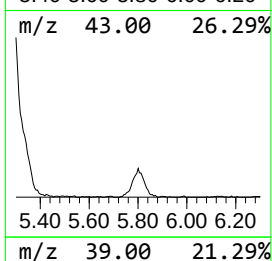
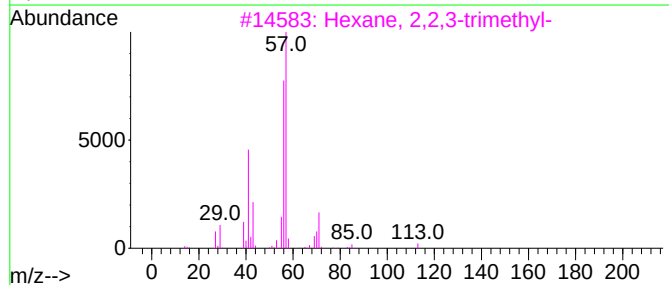
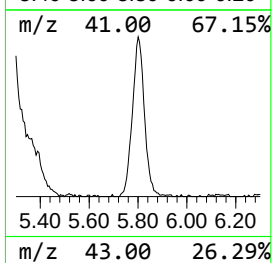
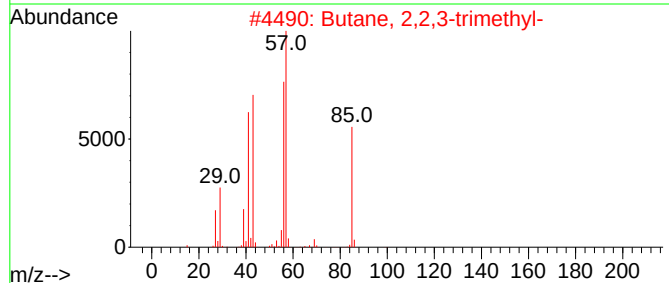
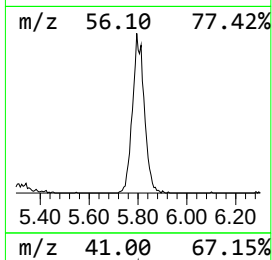
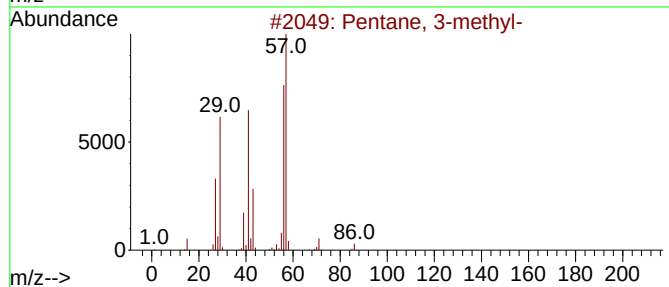
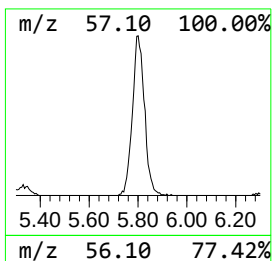
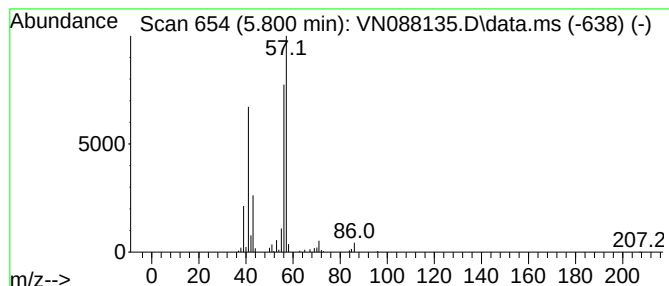
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.800	21.80 ug/l	423315	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56
4			3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	42
5			1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

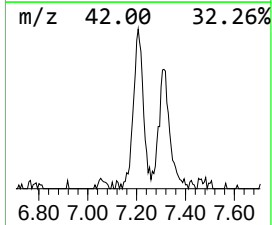
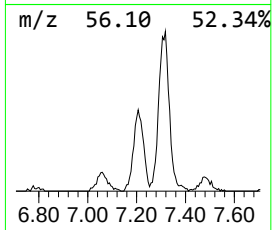
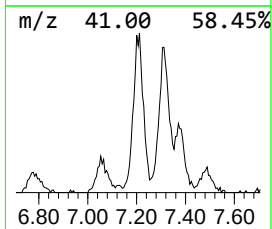
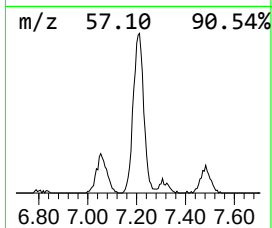
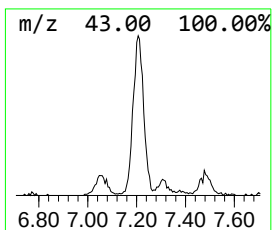
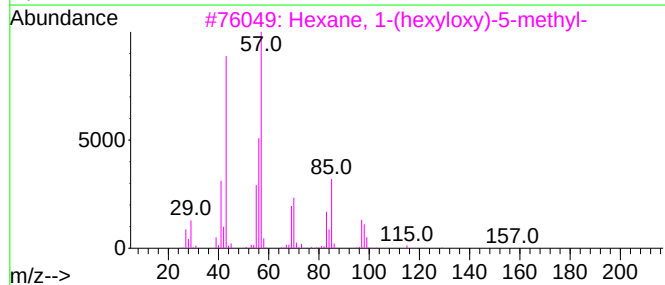
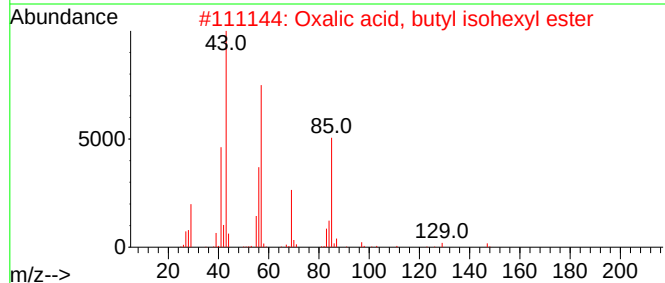
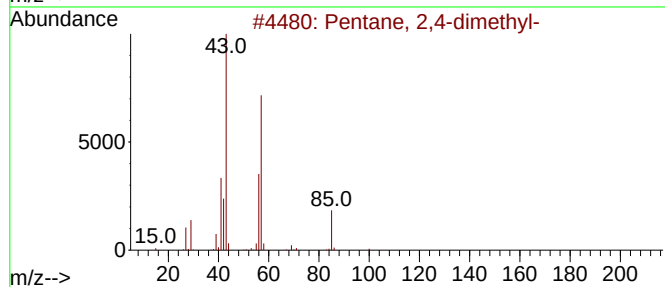
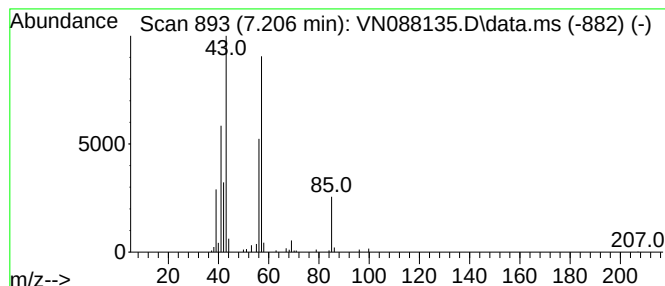
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.206	14.33 ug/l	278251	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2			Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	64
3			Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	50
4			2,2,6,6-Tetramethylheptane	156	C11H24	040117-45-1	50
5			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

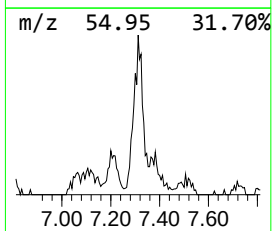
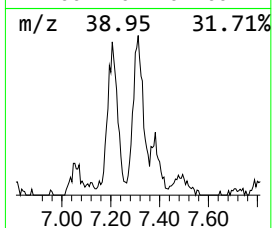
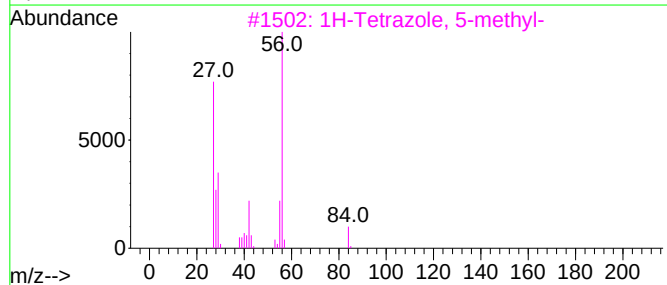
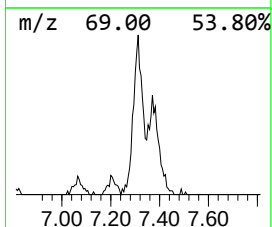
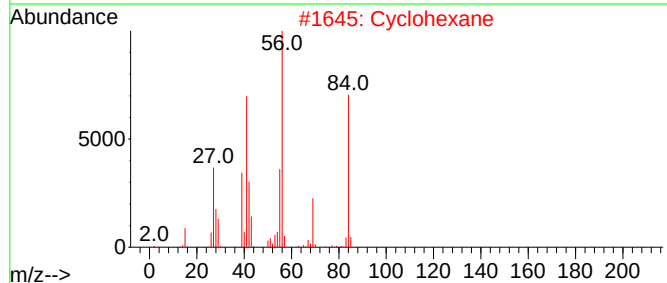
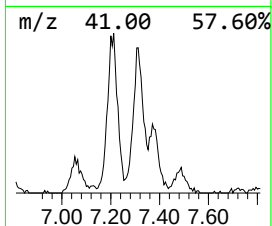
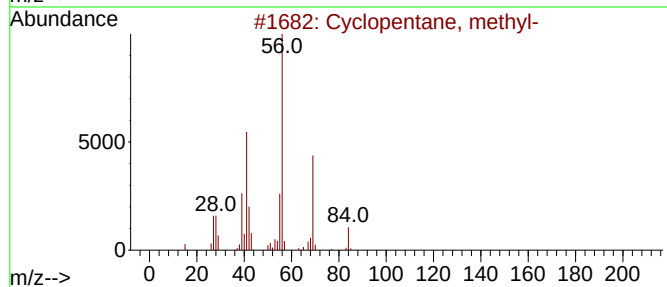
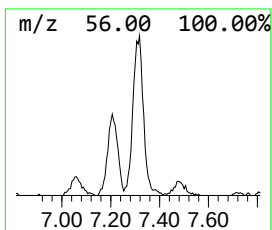
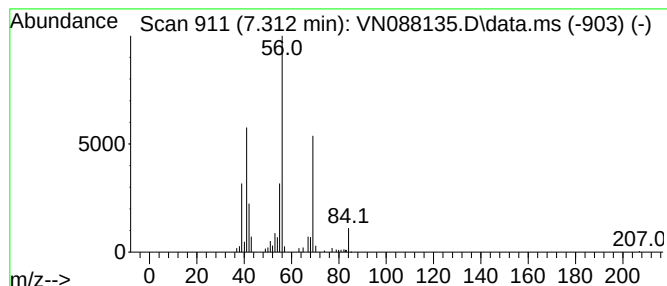
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.312	10.87 ug/l	211019	Pentafluorobenzene	8.206

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclohexane	84	C6H12	000110-82-7	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4		1-Heptene, 2-methyl-	112	C8H16	015870-10-7	64
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

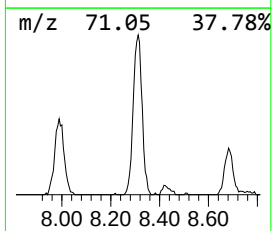
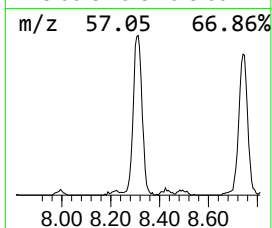
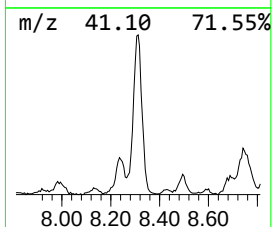
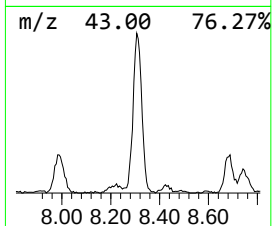
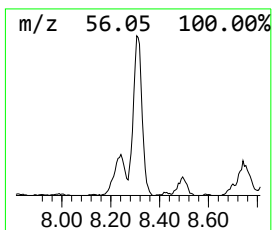
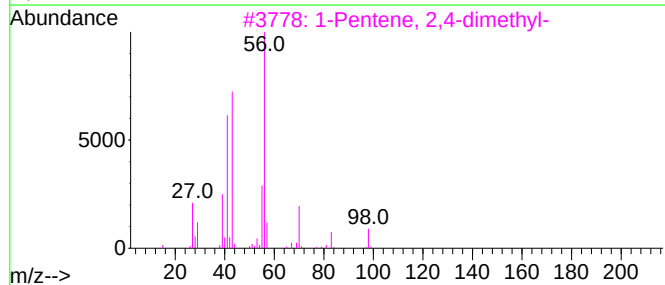
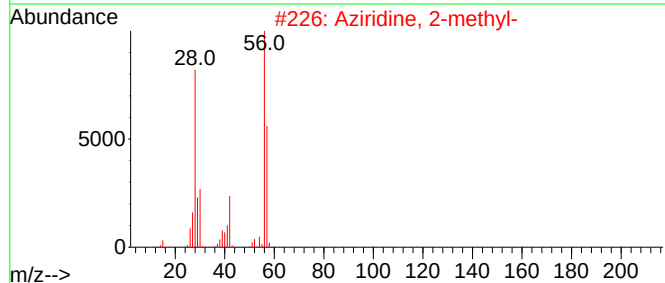
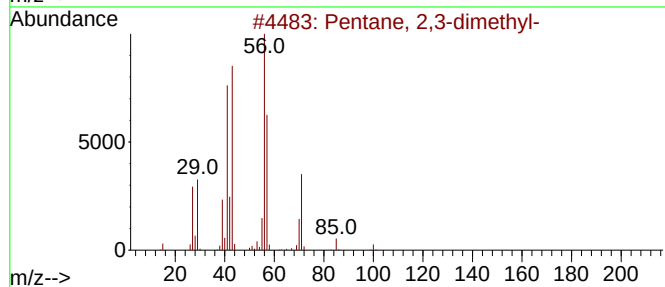
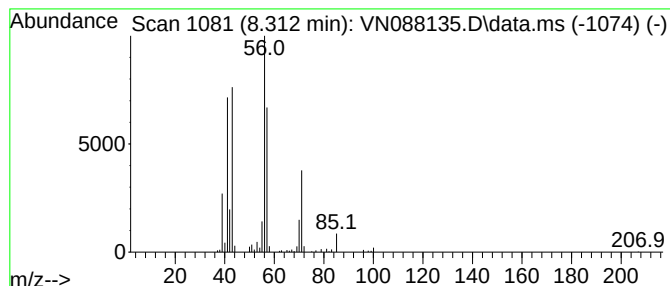
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Quant Title : SW846 8260

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

Peak Number 7 Pentane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.312	31.18 ug/l	605399	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	95
2			Aziridine, 2-methyl-	57	C3H7N	000075-55-8	49
3			1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	42
4			2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	40
5			Oxirane, 2-methyl-3-(1-methyleth...	100	C6H12O	001192-31-0	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

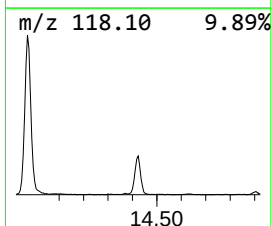
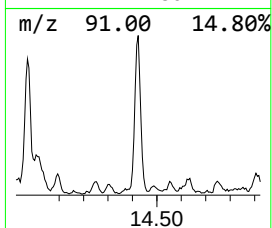
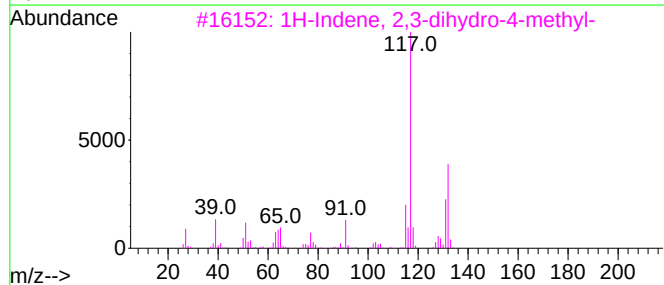
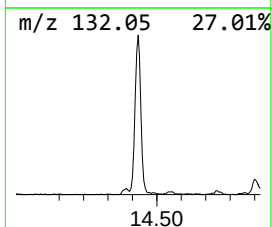
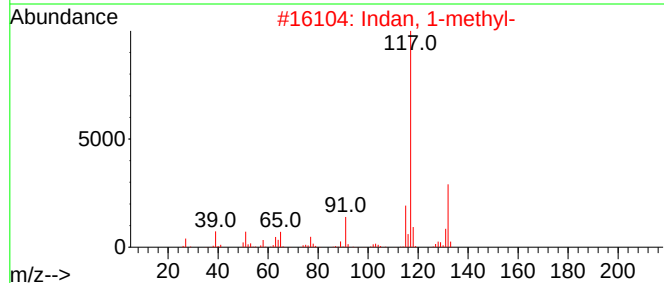
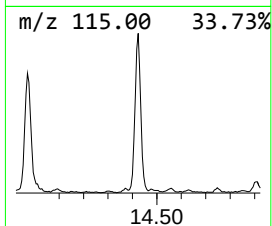
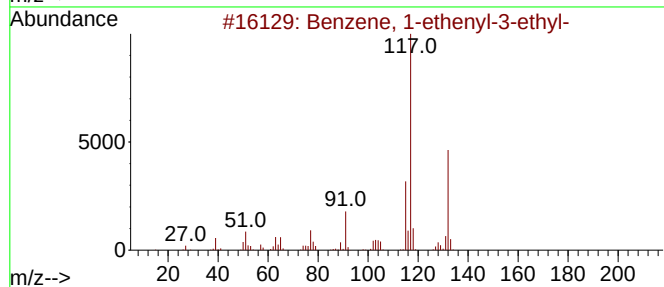
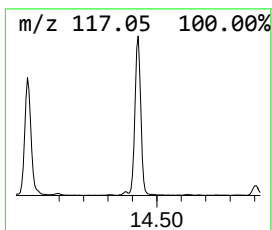
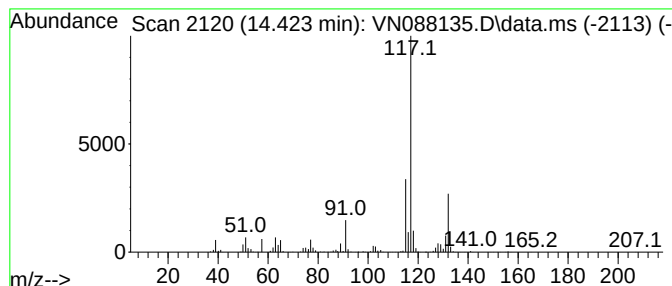
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.423	41.76 ug/l	1468090	1,4-Dichlorobenzene-d4	13.770

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2			Indan, 1-methyl-	132	C10H12	000767-58-8	90
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
5			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

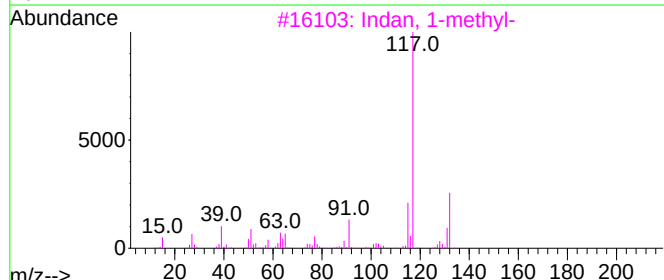
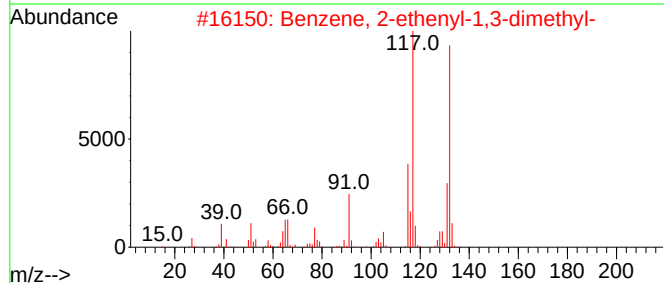
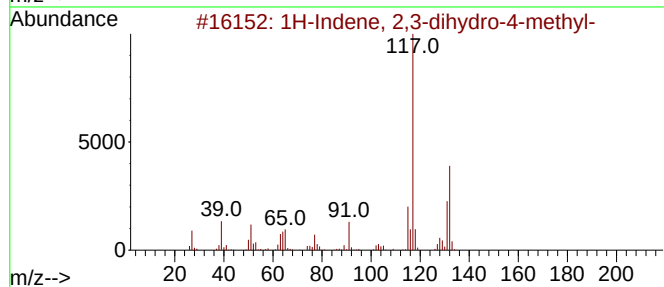
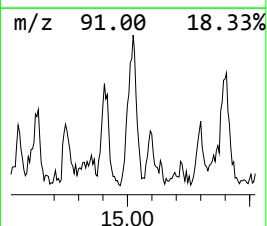
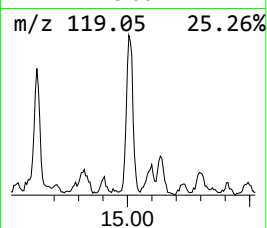
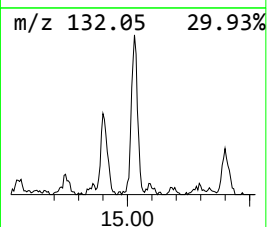
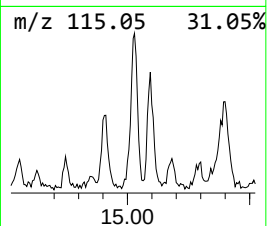
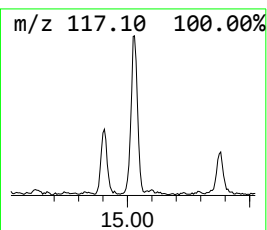
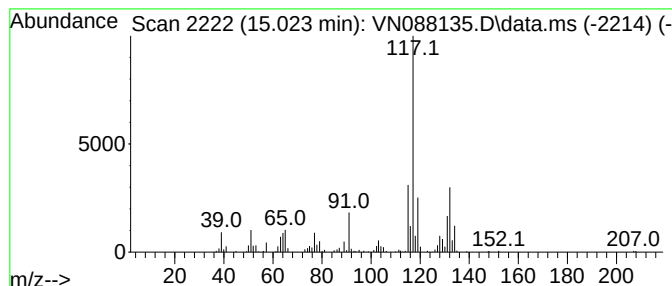
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.023	12.06 ug/l	424167	1,4-Dichlorobenzene-d4	13.770

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	90
3		Indan, 1-methyl-	132	C10H12	000767-58-8	76
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	74
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

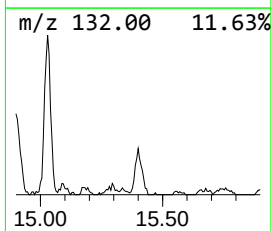
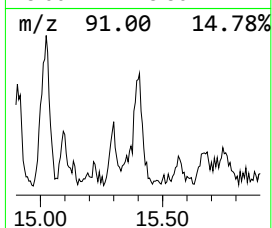
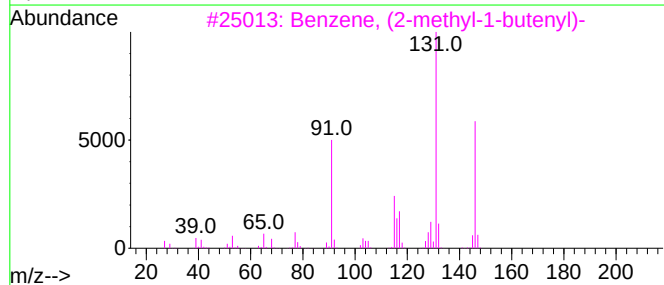
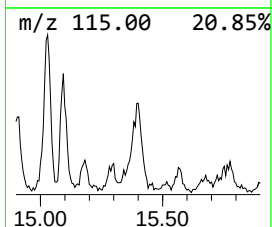
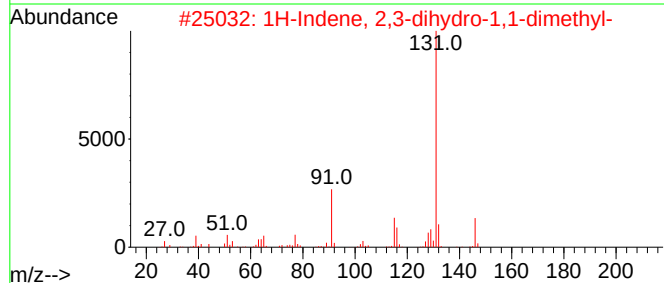
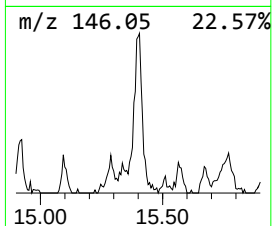
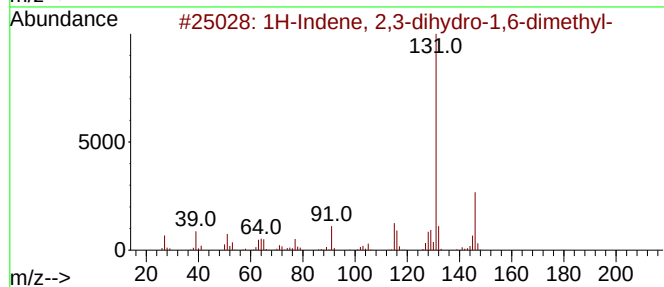
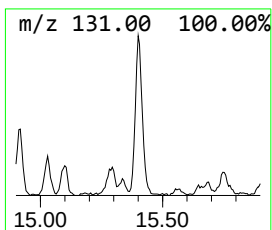
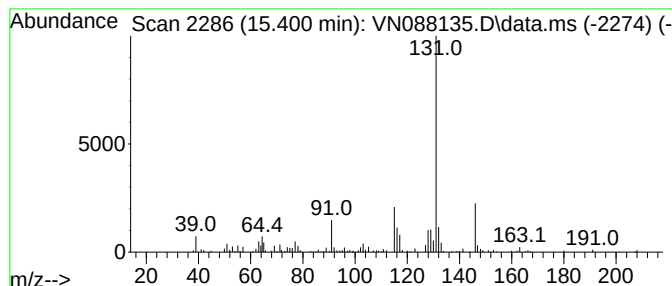
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.400	9.67 ug/l	339821	1,4-Dichlorobenzene-d4	13.770

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94		
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	90		
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90		
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90		
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102825\
Data File : VN088135.D
Acq On : 28 Oct 2025 13:38
Operator : JC\MD
Sample : Q3468-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.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-	3.265	57.7	ug/l	1120800	1	8.206	970845	50.0
Butane, 2,2-dim...	4.418	32.3	ug/l	627911	1	8.206	970845	50.0
Butane, 2,3-dim...	5.265	82.5	ug/l	1602440	1	8.206	970845	50.0
Pentane, 3-methyl-	5.800	21.8	ug/l	423315	1	8.206	970845	50.0
Pentane, 2,4-di...	7.206	14.3	ug/l	278251	1	8.206	970845	50.0
Cyclopentane, m...	7.312	10.9	ug/l	211019	1	8.206	970845	50.0
Pentane, 2,3-di...	8.312	31.2	ug/l	605399	1	8.206	970845	50.0
Benzene, 1-ethe...	14.423	41.8	ug/l	1468090	4	13.770	1757860	50.0
1H-Indene, 2,3-...	15.023	12.1	ug/l	424167	4	13.770	1757860	50.0
1H-Indene, 2,3-...	15.400	9.7	ug/l	339821	4	13.770	1757860	50.0