

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
 Data File : VN087853.D
 Acq On : 16 Sep 2025 16:29
 Operator : JC\MD
 Sample : Q3108-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW2

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\82N082125W.M
 Title : SW846 8260

Signal : TIC: VN087853.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.542	93	100	106	rBV	71483	125749	1.96%	0.168%
2	3.265	212	223	241	rBV	629159	1530180	23.79%	2.043%
3	3.653	279	289	299	rBV2	303285	776866	12.08%	1.037%
4	3.865	317	325	336	rVB4	44690	118734	1.85%	0.158%
5	4.048	346	356	365	rBV2	21659	64372	1.00%	0.086%
6	4.142	365	372	385	rVB2	35518	105241	1.64%	0.140%
7	4.412	405	418	434	rVB4	61404	223033	3.47%	0.298%
8	5.265	550	563	566	rBV2	138561	437988	6.81%	0.585%
9	5.330	567	574	605	rVB4	321107	1480864	23.02%	1.977%
10	5.800	639	654	669	rBV2	333371	1145337	17.81%	1.529%
11	6.300	728	739	752	rVB5	77625	244074	3.79%	0.326%
12	6.642	791	797	806	rVB3	29428	78563	1.22%	0.105%
13	6.777	809	820	828	rBV5	21572	65180	1.01%	0.087%
14	7.059	859	868	876	rBV5	28500	85202	1.32%	0.114%
15	7.206	882	893	901	rBV4	49451	149855	2.33%	0.200%
16	7.312	901	911	930	rVB	606394	1738717	27.03%	2.321%
17	7.994	1019	1027	1033	rVB4	31643	78831	1.23%	0.105%
18	8.147	1043	1053	1058	rBV	212946	516725	8.03%	0.690%
19	8.212	1058	1064	1075	rVV3	388223	1594242	24.79%	2.128%
20	8.306	1075	1080	1092	rVB2	168205	447378	6.96%	0.597%
21	8.430	1092	1101	1107	rBV2	118494	268518	4.17%	0.358%
22	8.494	1107	1112	1116	rVV4	38647	79300	1.23%	0.106%
23	8.559	1116	1123	1133	rVV	223505	534240	8.31%	0.713%
24	8.706	1133	1148	1149	rVV6	158068	364678	5.67%	0.487%
25	8.741	1150	1154	1165	rVV2	230660	761172	11.83%	1.016%
26	8.835	1165	1170	1179	rVB2	84447	189994	2.95%	0.254%
27	8.947	1179	1189	1199	rVB5	21680	65355	1.02%	0.087%
28	9.083	1202	1212	1225	rBV	540875	1186738	18.45%	1.584%
29	9.577	1281	1296	1301	rBV4	315795	934266	14.53%	1.247%
30	9.624	1301	1304	1313	rVB	139261	243016	3.78%	0.324%
31	9.759	1321	1327	1332	rVB2	47812	86965	1.35%	0.116%
32	9.847	1332	1342	1348	rVB2	54381	135253	2.10%	0.181%
33	9.994	1361	1367	1375	rVB	165810	340648	5.30%	0.455%
34	10.082	1375	1382	1385	rBV2	129043	248428	3.86%	0.332%
35	10.124	1385	1389	1395	rVB	189491	352427	5.48%	0.470%
36	10.218	1401	1405	1412	rVB2	67645	116229	1.81%	0.155%

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37	10.312	1412	1421	1425	rBV5	90863	242656	3.77%	0.324%
38	10.430	1436	1441	1445	rBV3	35579	70114	1.09%	0.094%
39	10.471	1445	1448	1453	rVB3	46441	76810	1.19%	0.103%
40	10.547	1453	1461	1467	rBV	866917	1646289	25.59%	2.198%
41	10.606	1467	1471	1478	rVB	180077	322714	5.02%	0.431%
42	10.888	1514	1519	1525	rBV3	43776	79891	1.24%	0.107%
43	10.971	1525	1533	1541	rVV7	85193	262768	4.09%	0.351%
44	11.247	1570	1580	1583	rBV6	26252	64576	1.00%	0.086%
45	11.447	1609	1614	1619	rVB3	55763	94190	1.46%	0.126%
46	11.653	1645	1649	1657	rVB6	31365	68032	1.06%	0.091%
47	11.847	1674	1682	1690	rBV	735578	1365026	21.22%	1.822%
48	11.941	1691	1698	1708	rVV	1672766	2889548	44.92%	3.857%
49	12.047	1708	1716	1727	rVV	2200625	4511212	70.14%	6.022%
50	12.376	1761	1772	1786	rBV	1075396	1939372	30.15%	2.589%
51	12.671	1816	1822	1834	rVB	501657	891832	13.87%	1.190%
52	12.829	1841	1849	1860	rBV	561660	1026454	15.96%	1.370%
53	13.012	1873	1880	1886	rBV	1624299	2676631	41.61%	3.573%
54	13.106	1886	1896	1900	rVV	1817344	3406014	52.95%	4.547%
55	13.153	1900	1904	1911	rVB	1050889	1727838	26.86%	2.306%
56	13.312	1924	1931	1942	rBV	2337631	3890514	60.49%	5.193%
57	13.459	1949	1956	1970	rBV	3989424	6432087	100.00%	8.586%
58	13.588	1970	1978	1984	rVB3	112054	276311	4.30%	0.369%
59	13.706	1994	1998	2002	rBV3	42686	66680	1.04%	0.089%
60	13.794	2002	2013	2021	rBV2	2320136	5090137	79.14%	6.795%
61	13.929	2030	2036	2039	rBV	491667	828006	12.87%	1.105%
62	13.970	2039	2043	2047	rBV	856701	1387991	21.58%	1.853%
63	14.094	2059	2064	2069	rVB2	116335	190442	2.96%	0.254%
64	14.159	2069	2075	2080	rBV2	299932	503244	7.82%	0.672%
65	14.217	2080	2085	2089	rBV	332955	569002	8.85%	0.760%
66	14.306	2094	2100	2106	rVV	1241195	1957312	30.43%	2.613%
67	14.370	2106	2111	2114	rVV	282375	464268	7.22%	0.620%
68	14.417	2114	2119	2127	rVV2	967356	1710238	26.59%	2.283%
69	14.547	2136	2141	2147	rVB	233225	371883	5.78%	0.496%
70	14.629	2147	2155	2159	rBV	733558	1242580	19.32%	1.659%
71	14.670	2159	2162	2166	rVV	450958	662262	10.30%	0.884%
72	14.706	2166	2168	2172	rVB2	81990	100795	1.57%	0.135%
73	14.900	2196	2201	2207	rBV	787404	1325575	20.61%	1.769%
74	15.023	2213	2222	2228	rBV2	1218173	2619839	40.73%	3.497%
75	15.094	2228	2234	2243	rVB3	162014	377466	5.87%	0.504%
76	15.182	2243	2249	2256	rBV3	202169	378036	5.88%	0.505%

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77	15.294	2256	2268	2273	rBV3	248603	633801	9.85%	0.846%
78	15.412	2280	2288	2295	rVB2	190229	452268	7.03%	0.604%
79	15.612	2315	2322	2330	rBV	530783	1012391	15.74%	1.351%
80	15.723	2335	2341	2346	rVV3	77894	165599	2.57%	0.221%
81	15.770	2346	2349	2361	rVB5	55552	137457	2.14%	0.183%
82	15.917	2366	2374	2382	rBV	112037	253737	3.94%	0.339%
83	16.100	2391	2405	2414	rBV4	68363	239566	3.72%	0.320%
84	16.223	2421	2426	2432	rBV5	30825	68591	1.07%	0.092%
85	16.306	2433	2440	2453	rVB3	50877	169096	2.63%	0.226%
86	16.682	2492	2504	2508	rBV7	25010	70063	1.09%	0.094%
87	16.747	2508	2515	2522	rBV	433020	990843	15.40%	1.323%

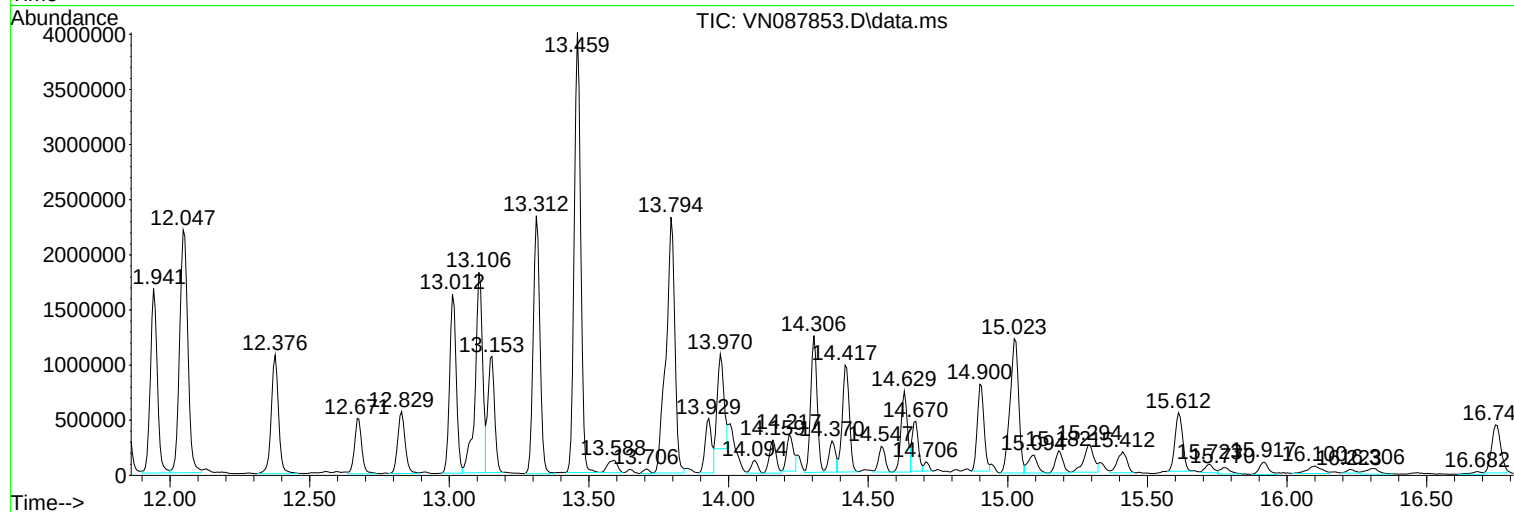
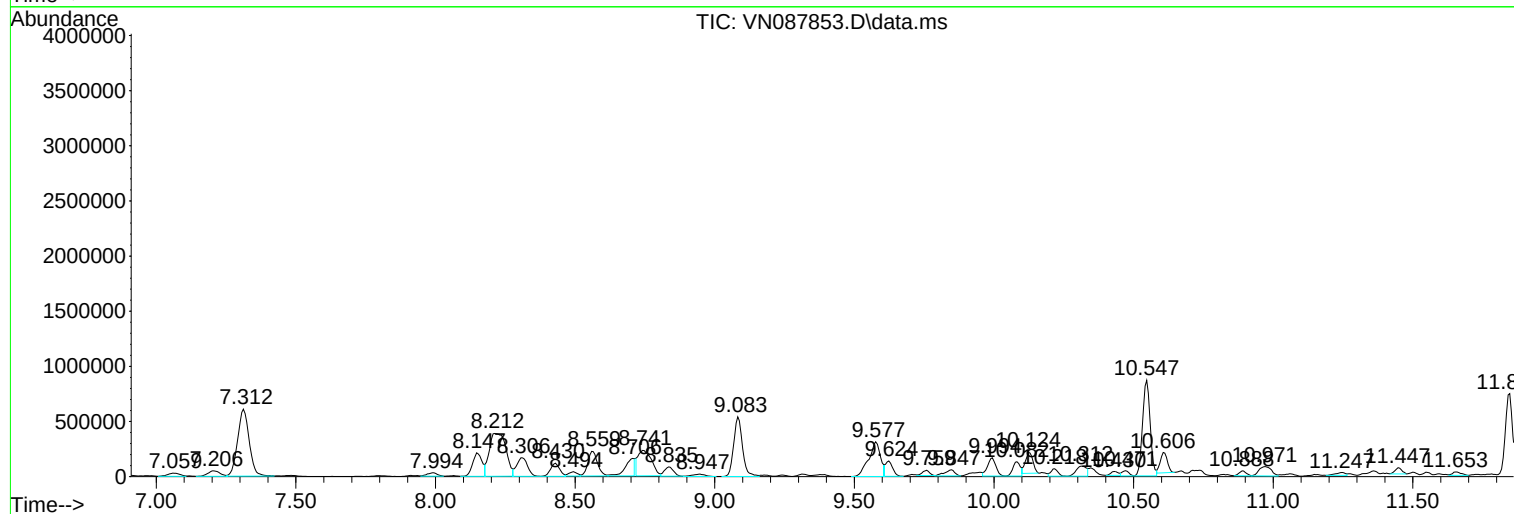
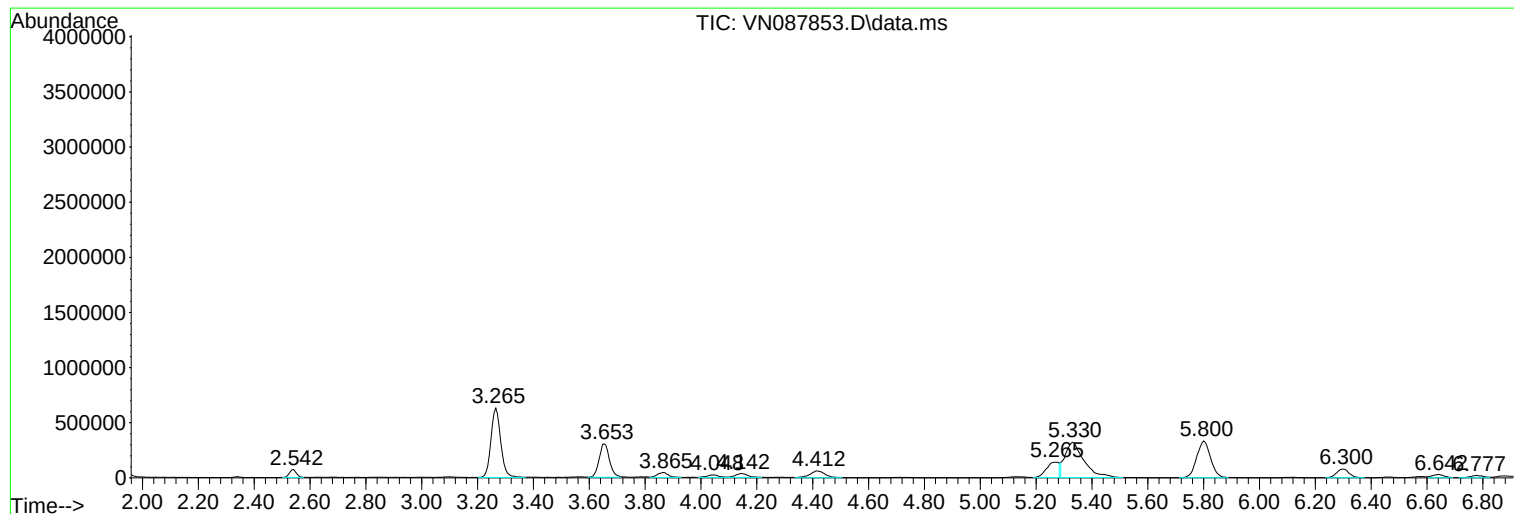
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082125W.M
Quant Title : SW846 8260

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



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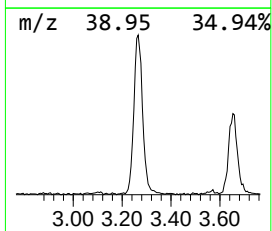
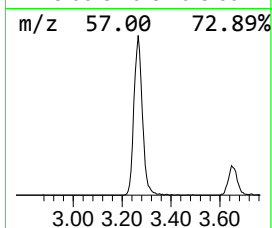
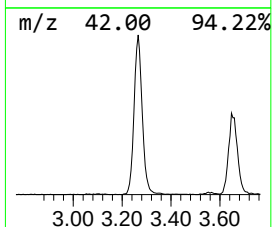
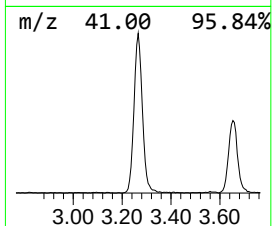
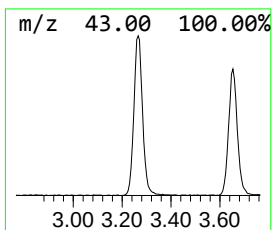
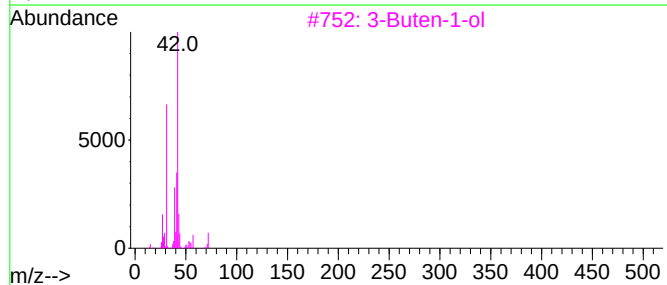
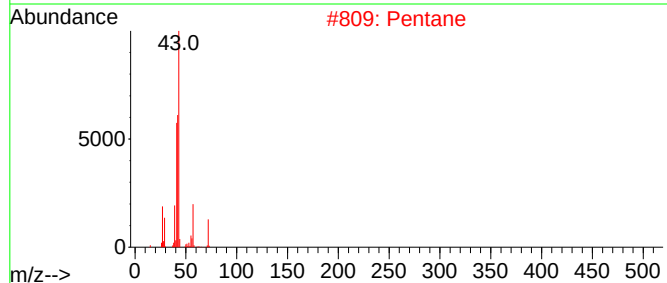
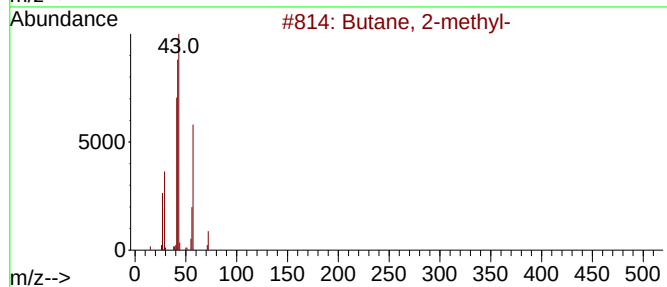
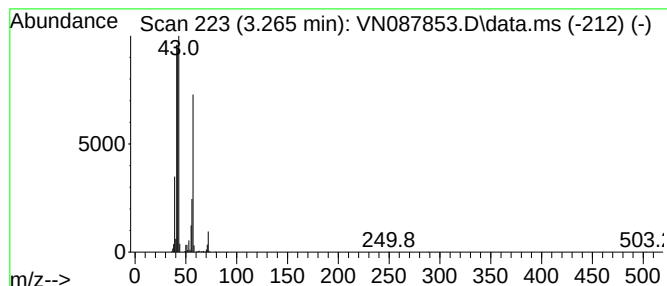
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082125W.M
Quant Title : SW846 8260

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	47.99 ug/l	1530180	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	86
2			Pentane	72	C5H12	000109-66-0	42
3			3-Buten-1-ol	72	C4H8O	000627-27-0	38
4			Propane, 2-methyl-1-nitro-	103	C4H9NO2	000625-74-1	10
5			1-Butene	56	C4H8	000106-98-9	10



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Instrument :
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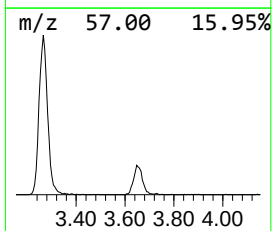
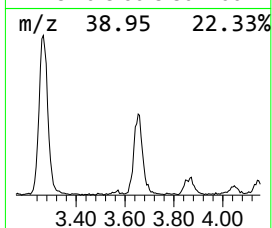
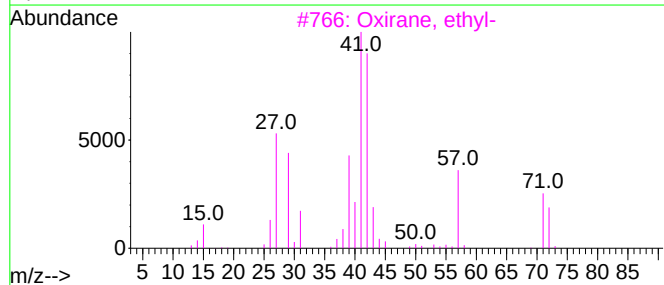
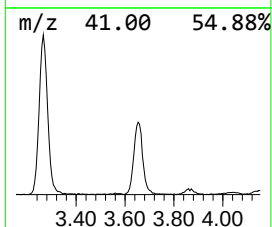
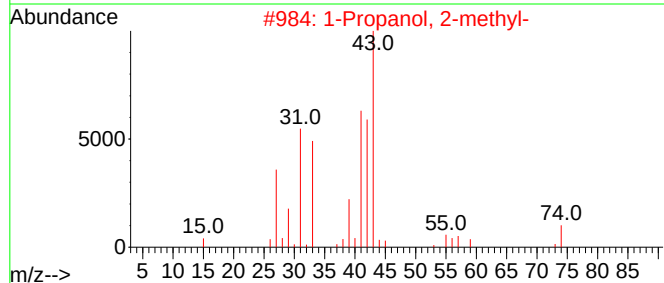
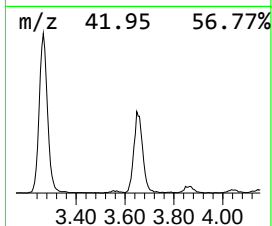
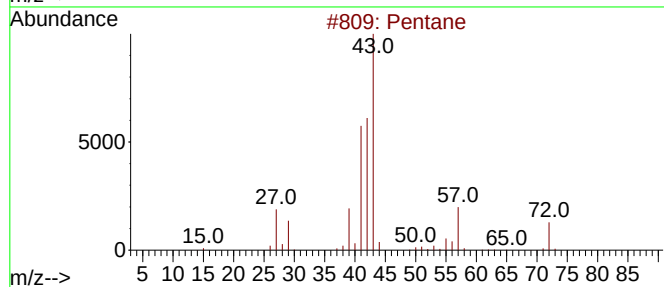
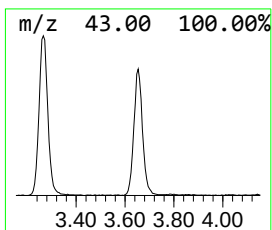
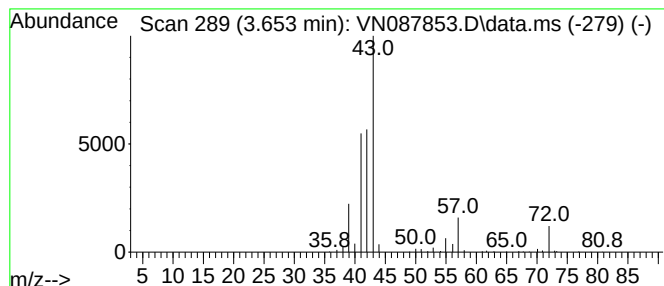
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082125W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.653	24.36 ug/l	776866	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane	72	C5H12	000109-66-0	91
2			1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	64
3			Oxirane, ethyl-	72	C4H8O	000106-88-7	47
4			Butane, 2-methyl-	72	C5H12	000078-78-4	42
5			2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	25



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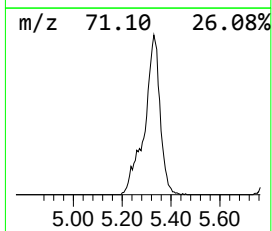
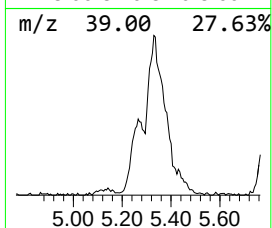
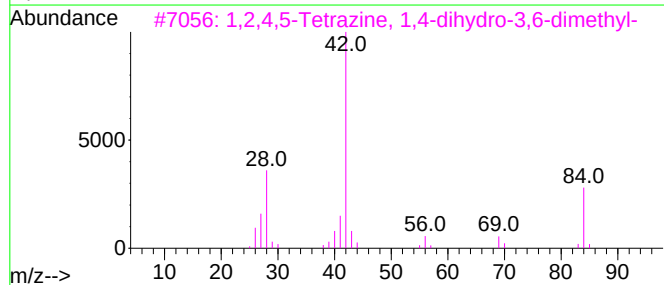
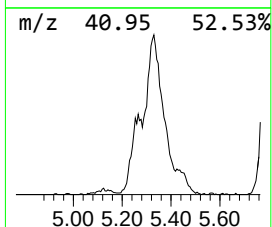
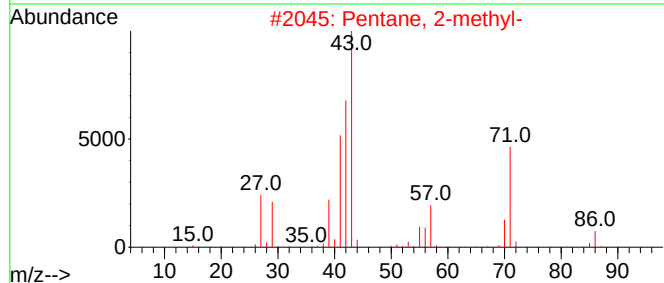
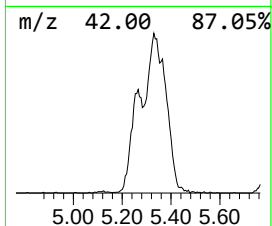
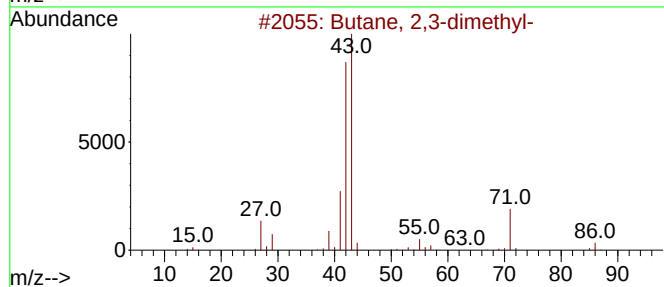
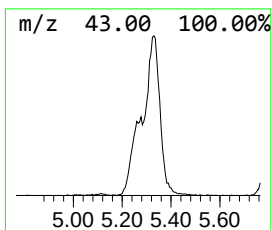
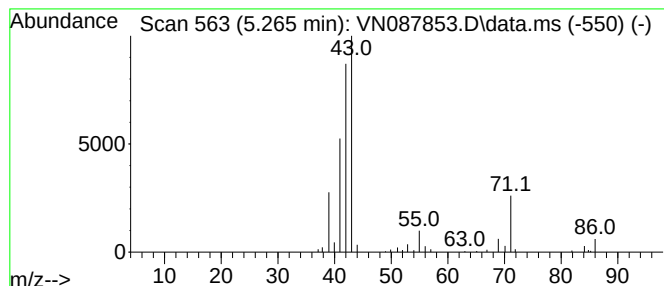
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082125W.M
Quant Title : SW846 8260

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

Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.265	13.74 ug/l	437988	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	72
3			1,2,4,5-Tetrazine, 1,4-dihydro-3...	112	C4H8N4	037454-64-1	28
4			Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	22
5			Pentane, 2-bromo-	150	C5H11Br	000107-81-3	17



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MW2

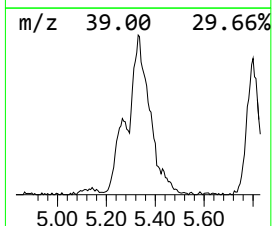
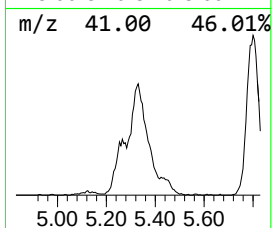
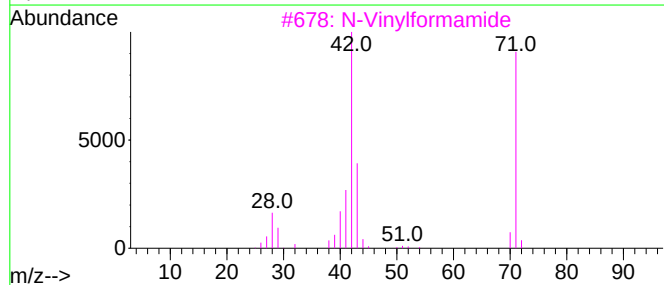
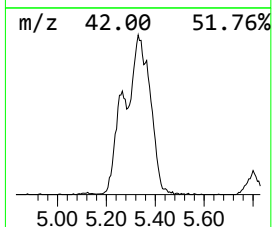
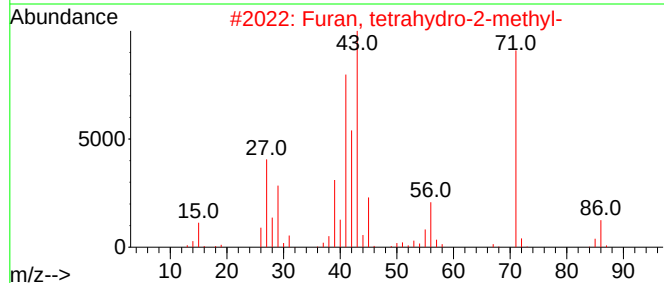
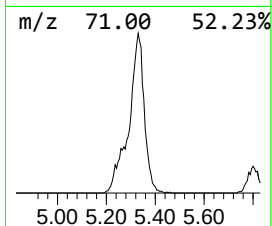
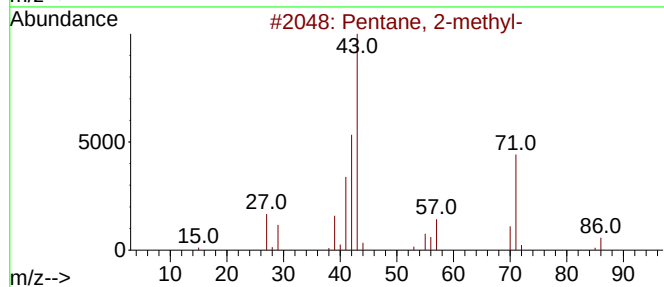
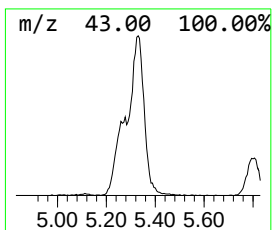
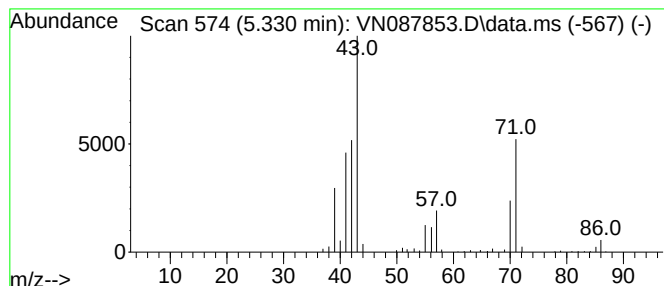
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.330	46.44 ug/l	1480860	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	90
2			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	52
3			N-Vinylformamide	71	C3H5NO	013162-05-5	50
4			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
5			Heptane, 4-methyl-	114	C8H18	000589-53-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW2

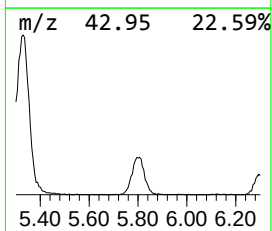
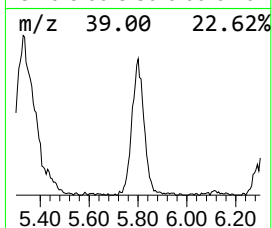
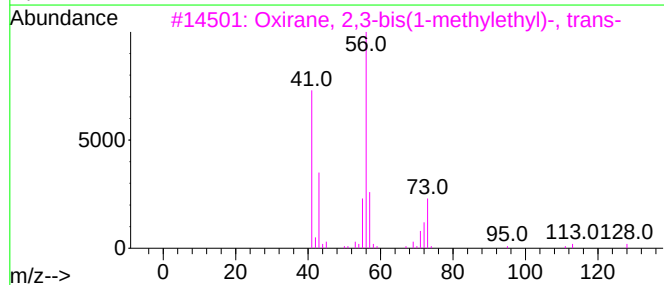
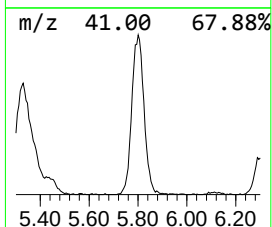
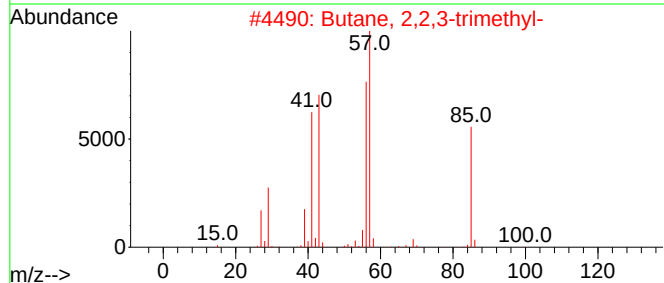
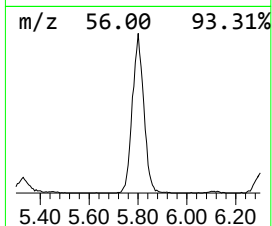
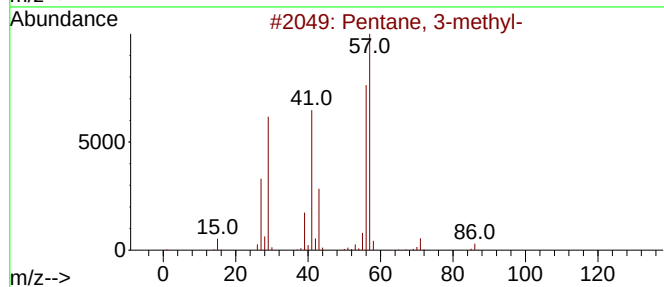
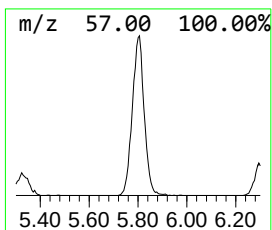
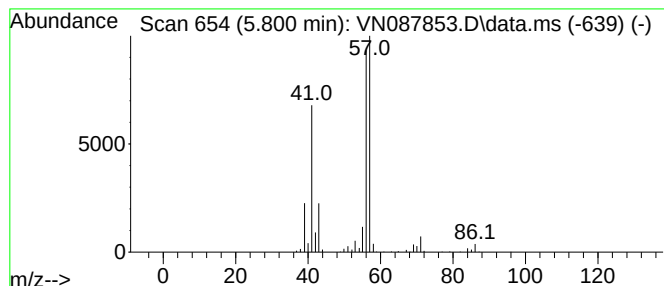
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.800	35.92 ug/l	1145340	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
3			Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	50
4			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	43
5			Cyclobutane, ethyl-	84	C6H12	004806-61-5	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW2

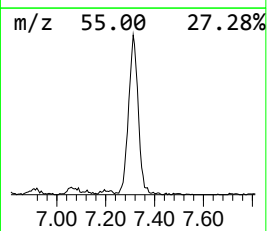
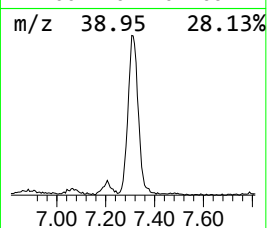
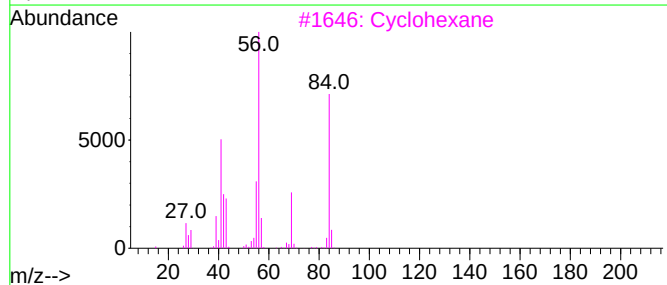
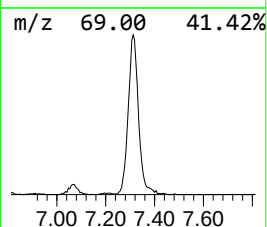
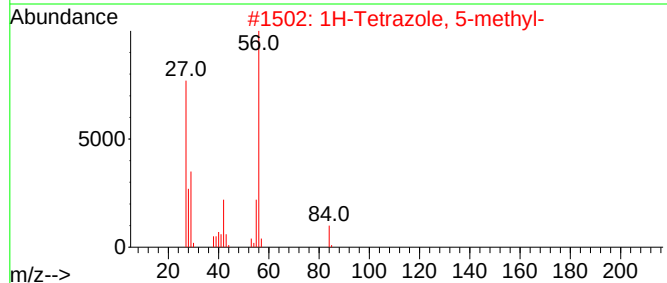
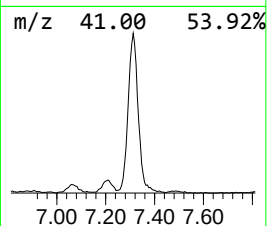
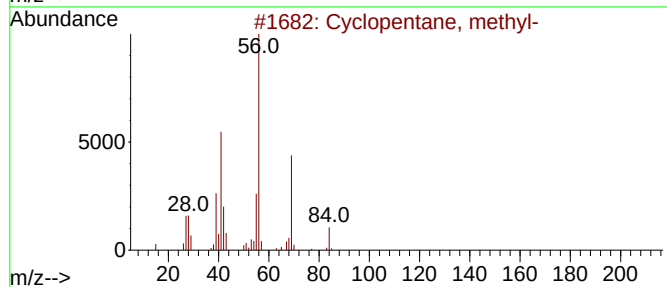
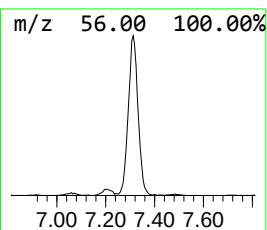
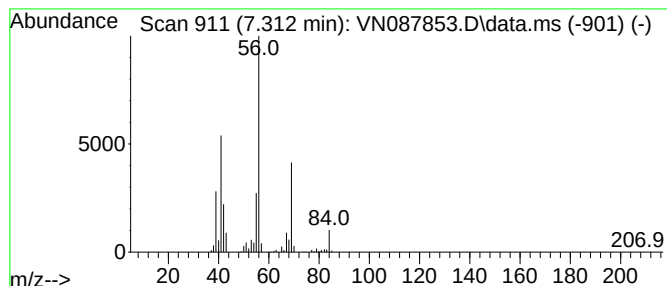
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.312	54.53 ug/l	1738720	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
3			Cyclohexane	84	C6H12	000110-82-7	64
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW2

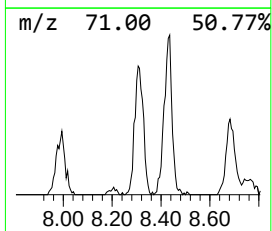
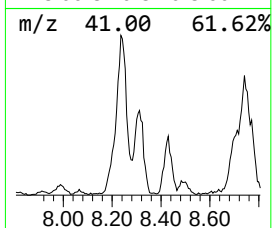
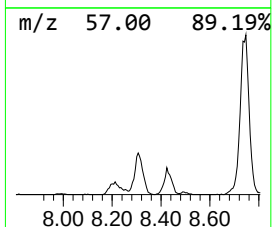
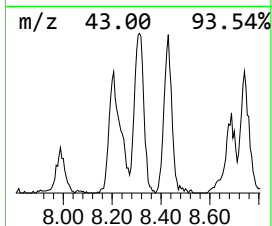
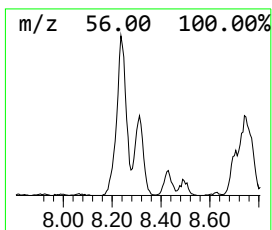
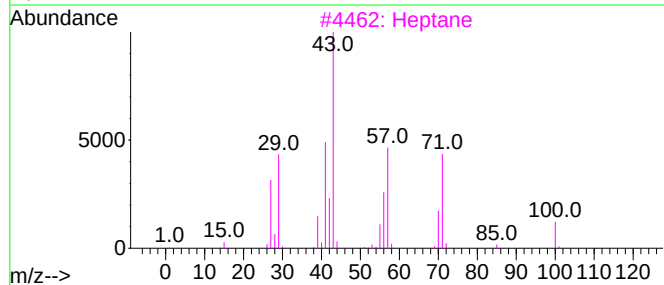
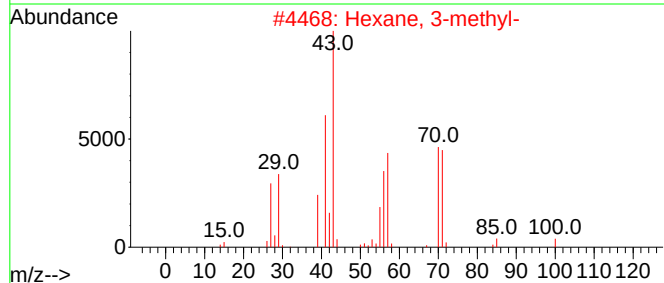
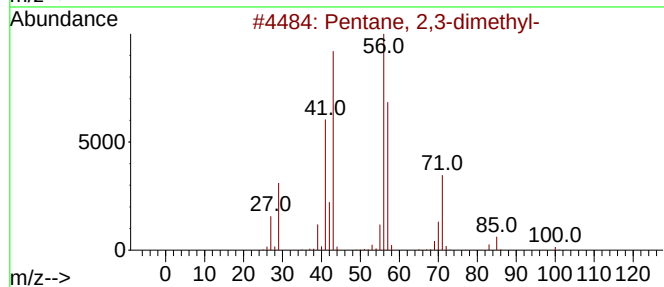
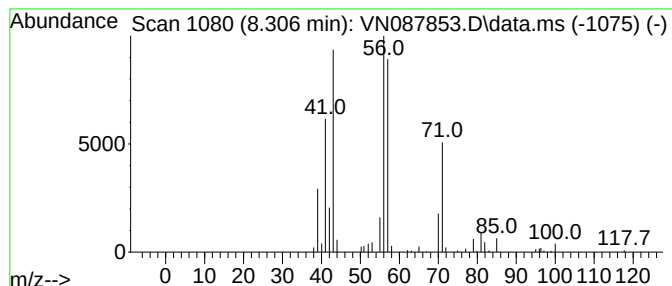
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Quant Title : SW846 8260

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

Peak Number 7 Pentane, 2,3-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.306	14.03 ug/l	447378	Pentafluorobenzene	8.206

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	86
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	46
3		Heptane	100	C7H16	000142-82-5	43
4		Aziridine, 2-methyl-	57	C3H7N	000075-55-8	43
5		Ethane, isocyanato-	71	C3H5NO	000109-90-0	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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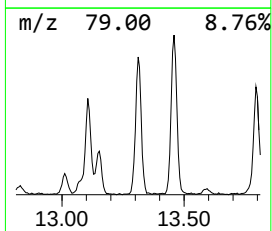
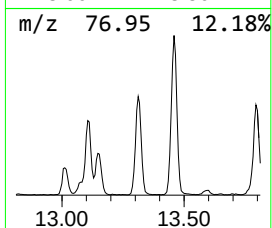
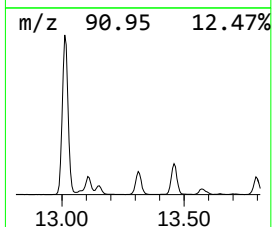
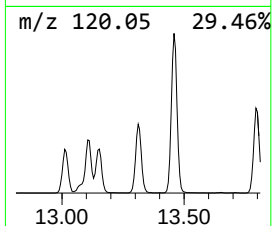
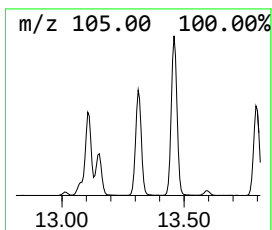
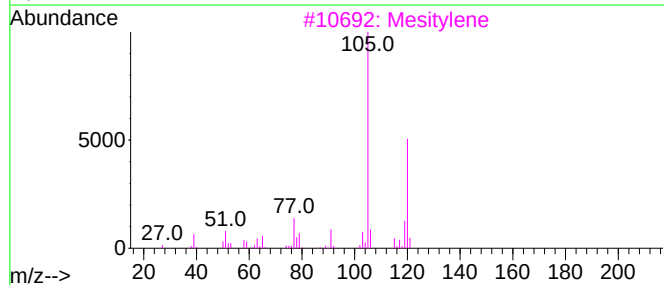
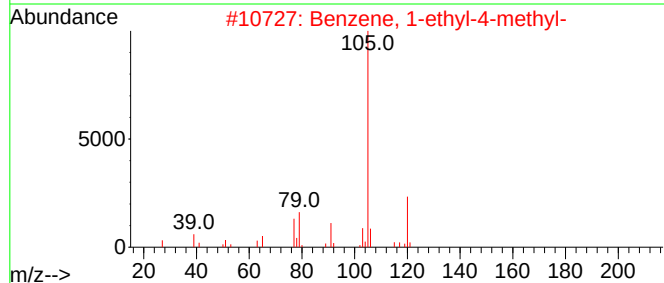
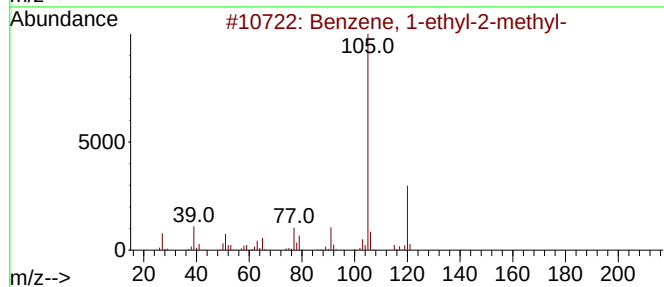
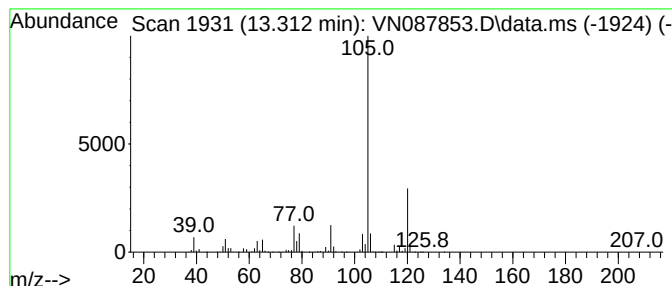
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082125W.M
Quant Title : SW846 8260

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

Peak Number 12 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.312	38.22 ug/l	3890510	1,4-Dichlorobenzene-d4	13.771

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-4-methyl-	120	C9H12	000622-96-8	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW2

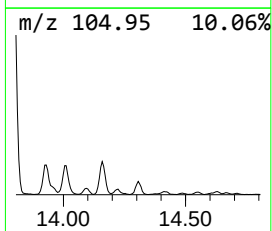
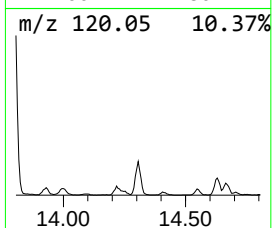
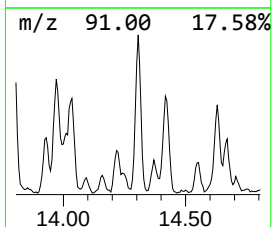
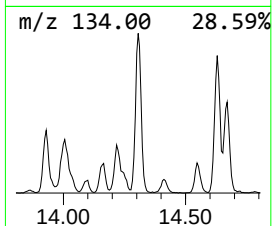
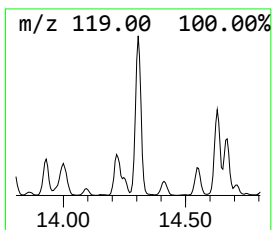
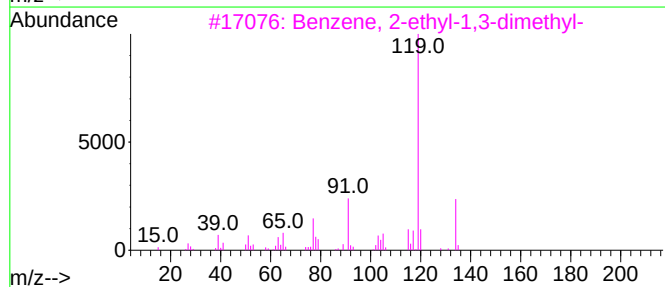
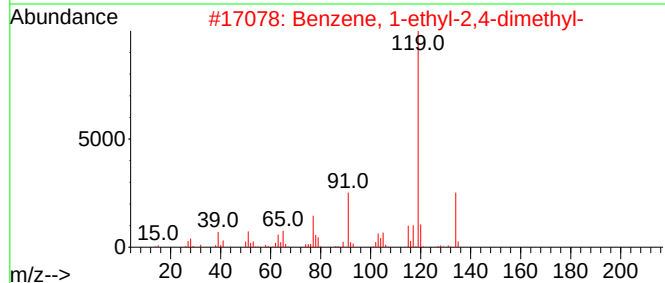
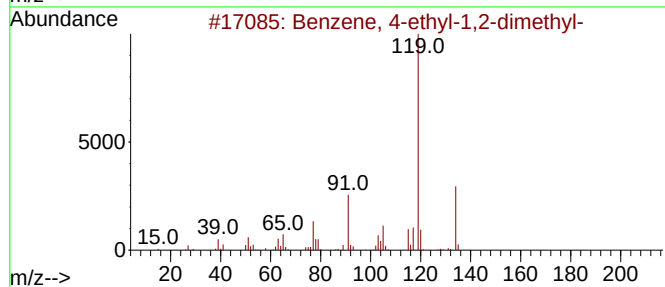
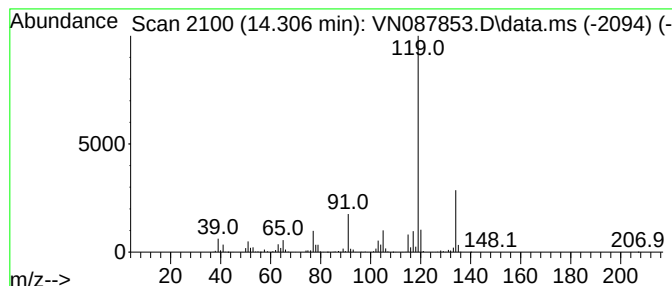
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Quant Title : SW846 8260

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

Peak Number 13 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.306	19.23 ug/l	1957310	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW2

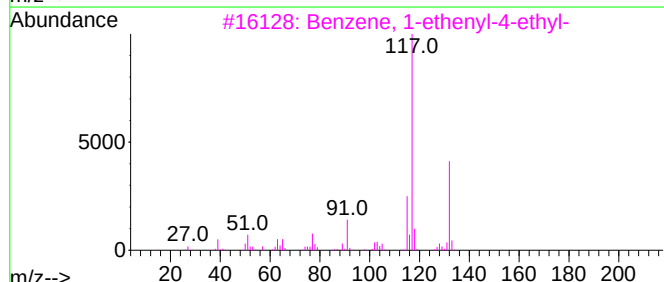
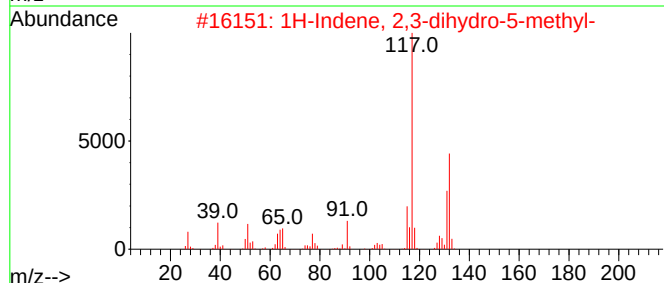
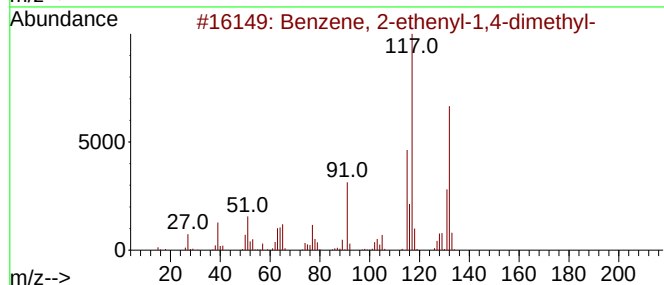
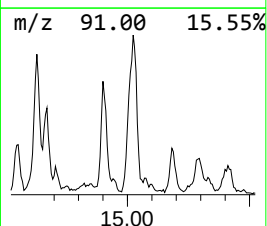
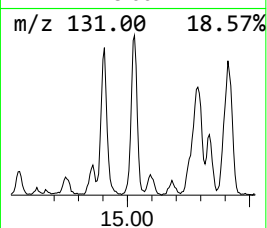
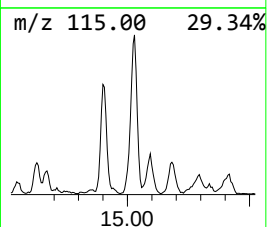
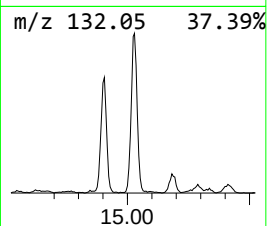
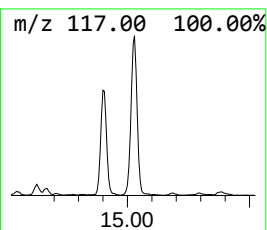
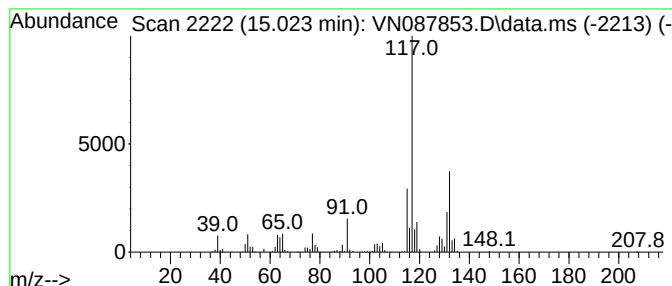
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Quant Title : SW846 8260

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

Peak Number 15 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.023	25.73 ug/l	2619840	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	93
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091625\
Data File : VN087853.D
Acq On : 16 Sep 2025 16:29
Operator : JC\MD
Sample : Q3108-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082125W.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	48.0	ug/l	1530180	1	8.206	1594240	50.0
Pentane	3.653	24.4	ug/l	776866	1	8.206	1594240	50.0
Butane, 2,3-dim...	5.265	13.7	ug/l	437988	1	8.206	1594240	50.0
Pentane, 2-methyl-	5.330	46.4	ug/l	1480860	1	8.206	1594240	50.0
Pentane, 3-methyl-	5.800	35.9	ug/l	1145340	1	8.206	1594240	50.0
Cyclopentane, m...	7.312	54.5	ug/l	1738720	1	8.206	1594240	50.0
Pentane, 2,3-di...	8.306	14.0	ug/l	447378	1	8.206	1594240	50.0
Benzene, 1-ethy...	13.312	38.2	ug/l	3890510	4	13.771	5090140	50.0
Benzene, 4-ethy...	14.306	19.2	ug/l	1957310	4	13.771	5090140	50.0
Benzene, 2-ethe...	15.023	25.7	ug/l	2619840	4	13.771	5090140	50.0