

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
 Data File : VN087108.D
 Acq On : 19 Jun 2025 13:59
 Operator : JC\MD
 Sample : Q2333-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 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\82N060625W.M
 Title : SW846 8260

Signal : TIC: VN087108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.353	62	68	74	rVB	9081	15468	1.19%	0.120%
2	2.553	95	102	108	rBV	18631	32297	2.49%	0.250%
3	3.283	216	226	239	rBV2	185956	464065	35.77%	3.589%
4	3.671	284	292	302	rVB3	6339	16618	1.28%	0.129%
5	4.447	412	424	443	rVB2	35216	125007	9.64%	0.967%
6	5.300	553	569	579	rBV2	78179	307740	23.72%	2.380%
7	5.836	647	660	677	rBV2	62754	214490	16.53%	1.659%
8	6.818	815	827	834	rBV4	5840	16684	1.29%	0.129%
9	7.094	866	874	881	rBV7	7160	21132	1.63%	0.163%
10	7.235	888	898	907	rBV2	25667	74554	5.75%	0.577%
11	7.341	907	916	924	rVV2	35007	99028	7.63%	0.766%
12	7.406	924	927	937	rVB5	7539	17433	1.34%	0.135%
13	8.018	1023	1031	1039	rVB5	11088	29634	2.28%	0.229%
14	8.177	1048	1058	1062	rBV	160750	395931	30.52%	3.062%
15	8.229	1062	1067	1078	rVV	274754	657426	50.68%	5.085%
16	8.335	1078	1085	1098	rVB2	68284	172546	13.30%	1.334%
17	8.512	1108	1115	1120	rBV4	9359	24176	1.86%	0.187%
18	8.588	1120	1128	1139	rVV	174450	407230	31.39%	3.150%
19	8.718	1140	1150	1152	rVV3	13816	32438	2.50%	0.251%
20	8.765	1152	1158	1169	rVV4	34278	119981	9.25%	0.928%
21	8.859	1169	1174	1183	rVB2	12609	26634	2.05%	0.206%
22	9.106	1206	1216	1229	rBV	457290	953913	73.53%	7.378%
23	9.341	1249	1256	1260	rBV4	7702	16370	1.26%	0.127%
24	9.388	1260	1264	1273	rVB2	7733	14649	1.13%	0.113%
25	9.565	1287	1294	1304	rBV5	24334	70044	5.40%	0.542%
26	9.865	1336	1345	1353	rBV3	10589	24889	1.92%	0.192%
27	10.018	1365	1371	1379	rVB	29504	59263	4.57%	0.458%
28	10.112	1379	1387	1388	rBV3	11456	19316	1.49%	0.149%
29	10.147	1388	1393	1406	rVB	52219	110719	8.53%	0.856%
30	10.323	1417	1423	1431	rBV5	9193	21849	1.68%	0.169%
31	10.453	1439	1445	1448	rBV2	11777	21693	1.67%	0.168%
32	10.494	1448	1452	1457	rVB2	8631	15057	1.16%	0.116%
33	10.565	1457	1464	1473	rBV	678539	1297245	100.00%	10.033%
34	10.759	1485	1497	1503	rBV4	38395	83946	6.47%	0.649%
35	10.835	1503	1510	1517	rVV2	30821	60085	4.63%	0.465%
36	10.912	1519	1523	1529	rVV6	10732	22580	1.74%	0.175%

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37	10.982	1529	1535	1545	rVV3	29936	78014	6.01%	0.603%
38	11.253	1574	1581	1586	rBV4	14304	30105	2.32%	0.233%
39	11.382	1598	1603	1609	rVB5	7913	13575	1.05%	0.105%
40	11.676	1646	1653	1660	rBV7	6128	13573	1.05%	0.105%
41	11.865	1677	1685	1694	rBV	589159	1073848	82.78%	8.305%
42	11.965	1697	1702	1707	rVB	28007	45879	3.54%	0.355%
43	12.076	1713	1721	1728	rBV	55632	115495	8.90%	0.893%
44	12.394	1767	1775	1784	rBV2	20166	43555	3.36%	0.337%
45	12.570	1796	1805	1813	rBV9	5371	16522	1.27%	0.128%
46	12.694	1821	1826	1832	rVB	54047	87649	6.76%	0.678%
47	12.847	1842	1852	1862	rBV	456007	803655	61.95%	6.216%
48	13.035	1875	1884	1889	rVV	22291	41960	3.23%	0.325%
49	13.094	1889	1894	1897	rVV4	10826	19737	1.52%	0.153%
50	13.129	1897	1900	1904	rVV	21506	33110	2.55%	0.256%
51	13.335	1925	1935	1943	rBV	33296	62156	4.79%	0.481%
52	13.435	1943	1952	1954	rBV7	7054	14073	1.08%	0.109%
53	13.482	1954	1960	1967	rVB	112682	186388	14.37%	1.442%
54	13.600	1973	1980	1988	rBV3	35729	87243	6.73%	0.675%
55	13.788	2004	2012	2024	rBV	616891	1124164	86.66%	8.694%
56	13.888	2024	2029	2033	rVB	18926	32948	2.54%	0.255%
57	13.947	2034	2039	2042	rBV	42822	66169	5.10%	0.512%
58	13.994	2042	2047	2050	rVV	74013	138466	10.67%	1.071%
59	14.029	2050	2053	2060	rVB2	105223	163752	12.62%	1.266%
60	14.111	2062	2067	2073	rVB	54737	92898	7.16%	0.718%
61	14.176	2073	2078	2084	rVB3	16514	30394	2.34%	0.235%
62	14.270	2084	2094	2099	rBV3	40816	93645	7.22%	0.724%
63	14.329	2099	2104	2112	rVB	71997	114745	8.85%	0.887%
64	14.441	2112	2123	2130	rBV	450631	783874	60.43%	6.063%
65	14.511	2130	2135	2136	rBV3	12567	16386	1.26%	0.127%
66	14.576	2142	2146	2150	rVB2	21634	30935	2.38%	0.239%
67	14.653	2150	2159	2164	rBV	121340	215878	16.64%	1.670%
68	14.770	2174	2179	2182	rBV2	25019	42452	3.27%	0.328%
69	14.858	2188	2194	2195	rBV4	8019	14464	1.11%	0.112%
70	14.935	2201	2207	2217	rBV5	42131	90545	6.98%	0.700%
71	15.047	2217	2226	2231	rVB4	21978	58701	4.53%	0.454%
72	15.117	2231	2238	2244	rBV2	52432	101036	7.79%	0.781%
73	15.200	2248	2252	2258	rVV4	17796	27654	2.13%	0.214%
74	15.317	2265	2272	2277	rBV2	46663	92199	7.11%	0.713%
75	15.382	2277	2283	2285	rVV2	25416	52799	4.07%	0.408%
76	15.423	2285	2290	2297	rVB2	83312	163725	12.62%	1.266%

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77	15.535	2304	2309	2313	rBV4	6954	13274	1.02%	0.103%
78	15.594	2313	2319	2322	rVV4	22486	39325	3.03%	0.304%
79	15.635	2322	2326	2331	rVB	30847	51511	3.97%	0.398%
80	15.788	2348	2352	2361	rVB4	23927	62621	4.83%	0.484%
81	15.947	2369	2379	2383	rBV9	14914	41770	3.22%	0.323%
82	16.064	2389	2399	2408	rBV7	17573	67731	5.22%	0.524%
83	16.147	2408	2413	2420	rVV6	12510	32554	2.51%	0.252%
84	16.252	2425	2431	2438	rVV	35114	83333	6.42%	0.645%
85	16.329	2439	2444	2454	rVB4	24254	63799	4.92%	0.493%
86	16.476	2464	2469	2476	rBV4	11698	26173	2.02%	0.202%
87	16.670	2497	2502	2503	rBV3	11694	15043	1.16%	0.116%
88	16.705	2504	2508	2515	rVB	20834	39797	3.07%	0.308%
89	16.770	2515	2519	2522	rBV4	10600	16316	1.26%	0.126%

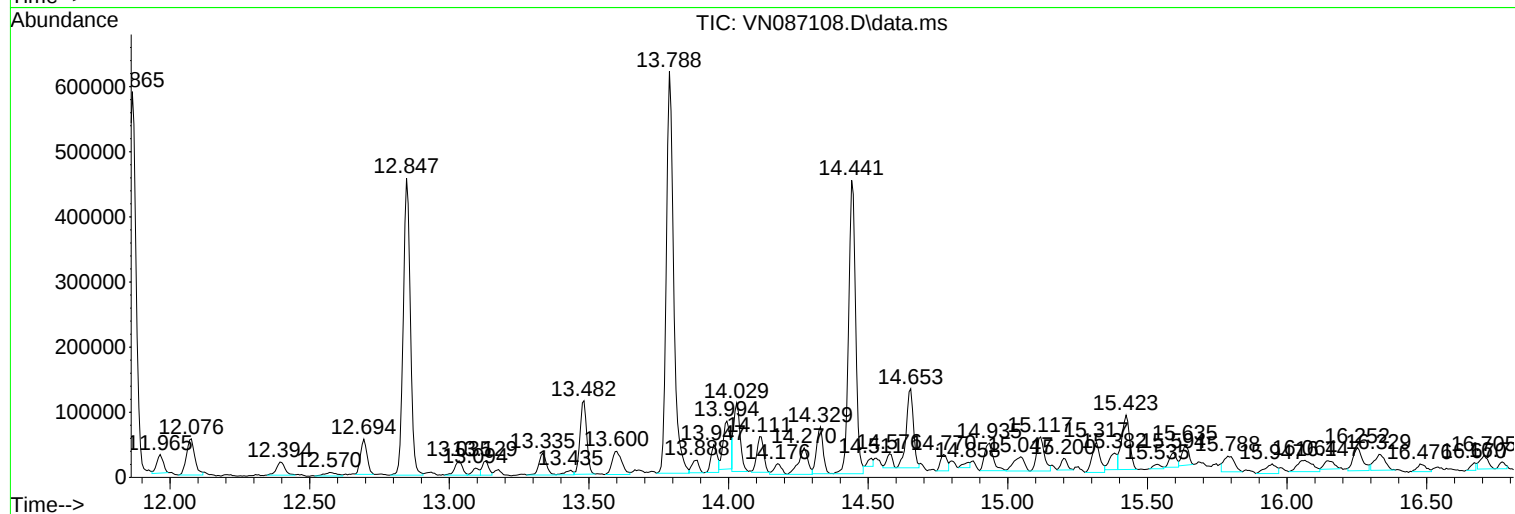
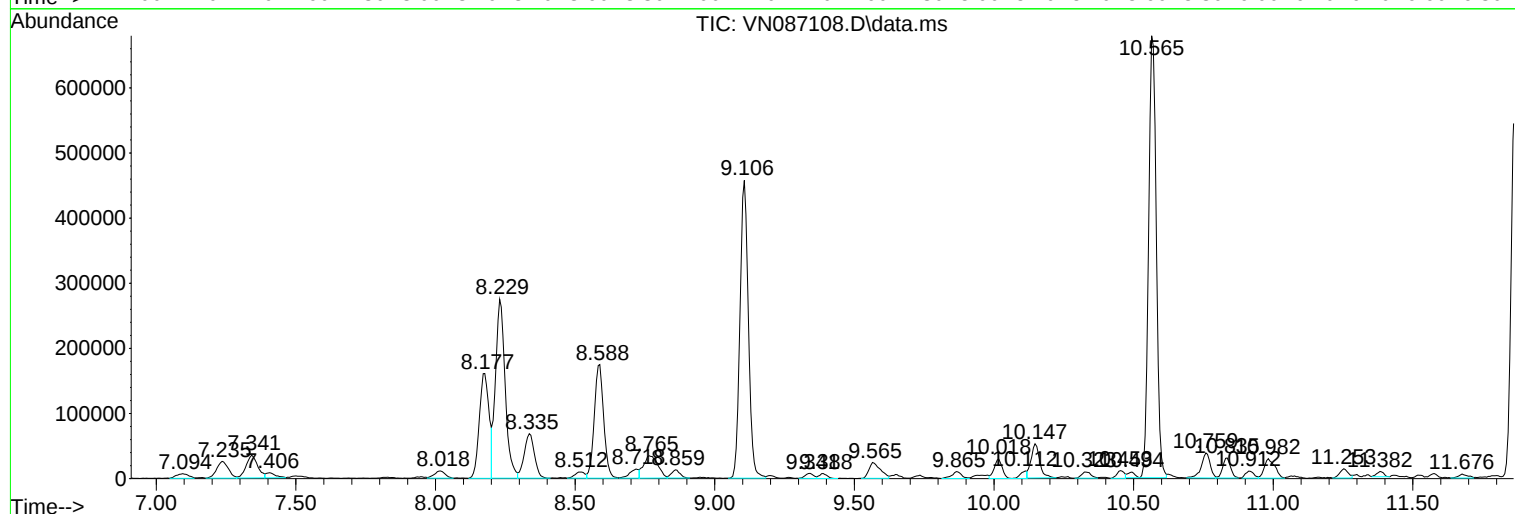
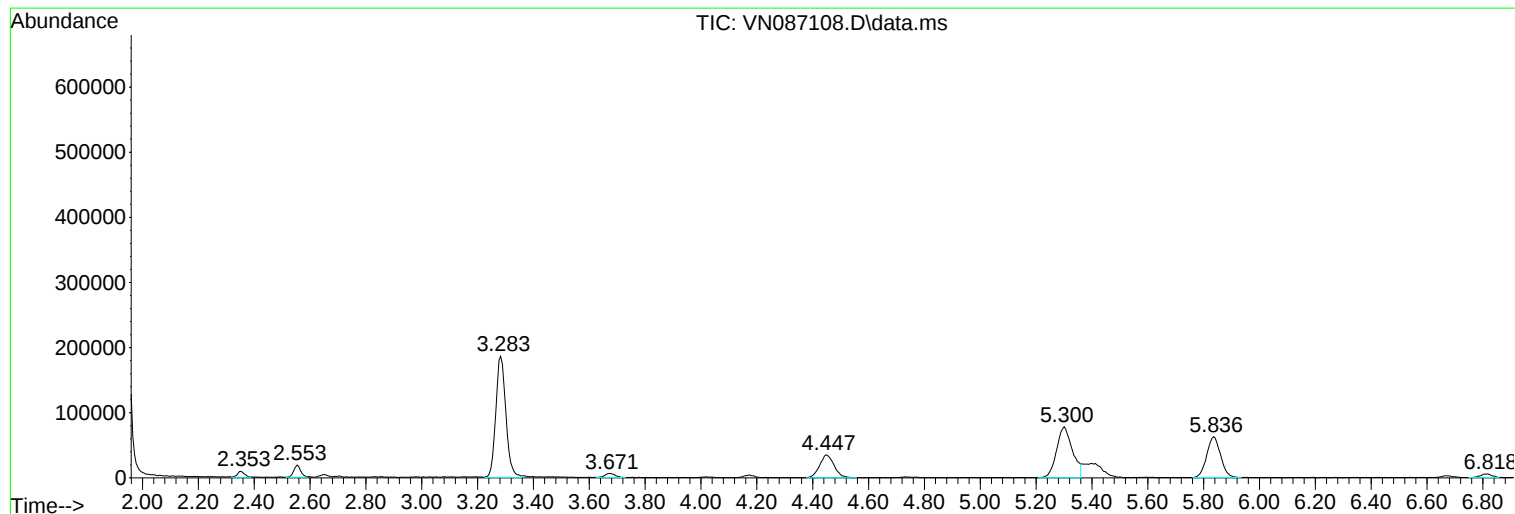
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

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



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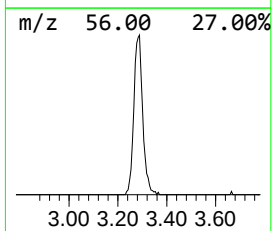
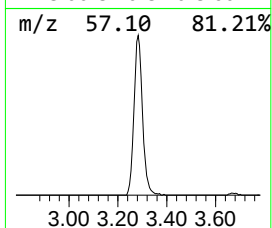
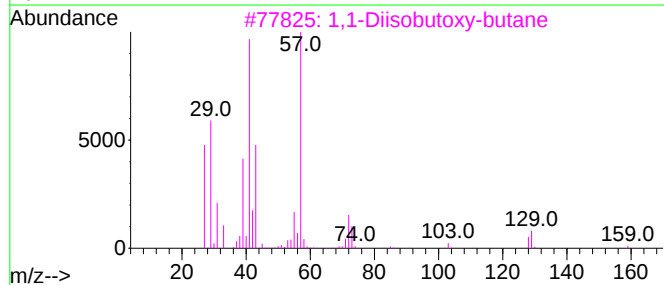
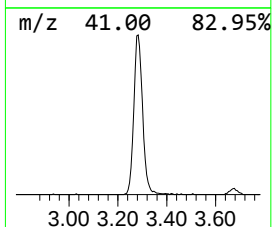
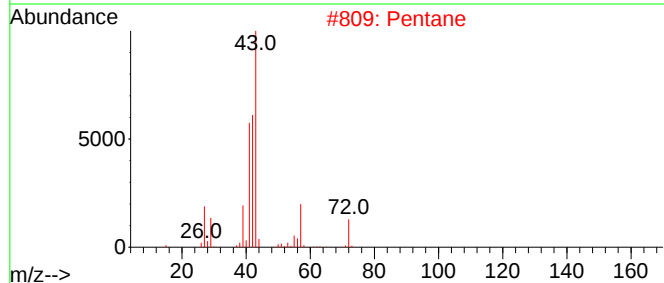
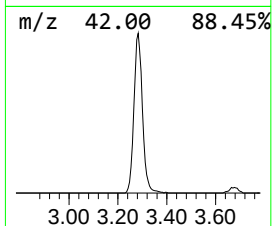
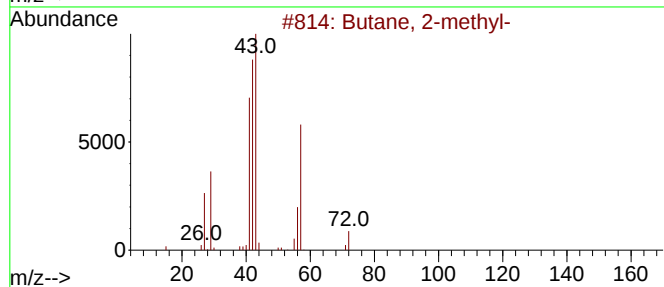
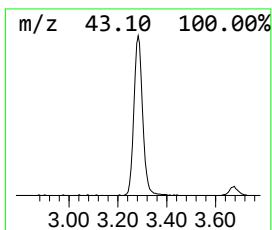
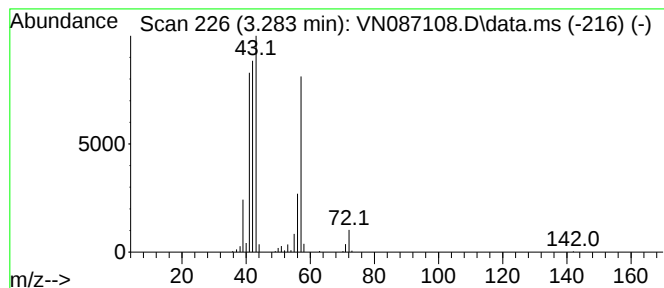
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.283	35.29 ug/l	464065	Pentafluorobenzene	8.229

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			Pentane	72	C5H12	000109-66-0	38
3			1,1-Diisobutoxy-butane	202	C12H26O2	013002-16-9	27
4			3-Buten-1-ol	72	C4H8O	000627-27-0	25
5			Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	22



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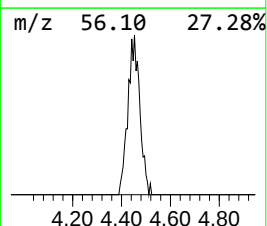
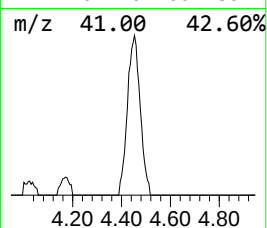
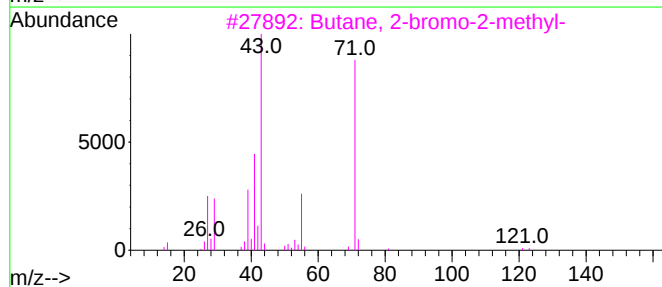
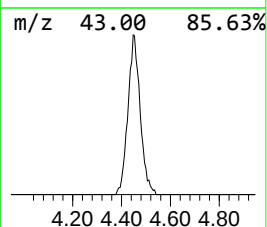
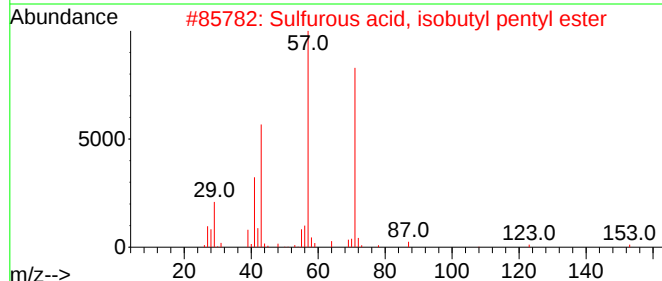
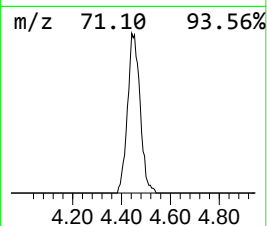
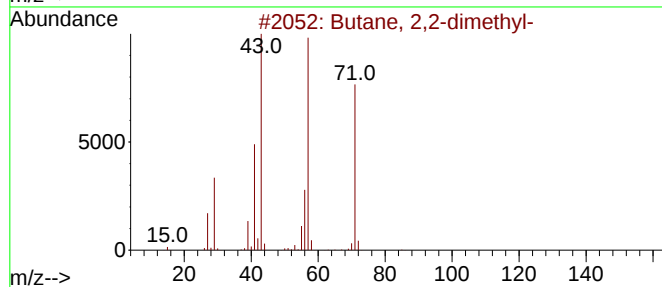
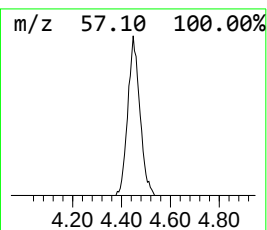
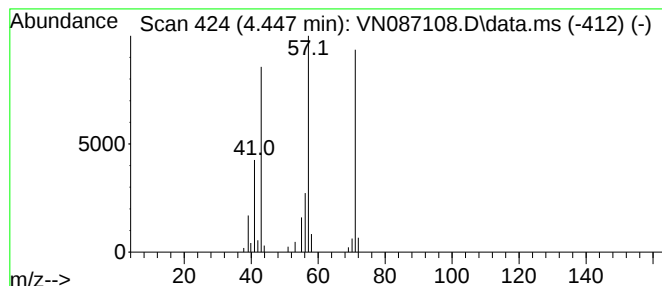
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

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

Peak Number 2 Butane, 2,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.447	9.51 ug/l	125007	Pentafluorobenzene	8.229

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	83
2			Sulfurous acid, isobutyl pentyl ...	208	C9H20O3S	1000309-13-8	50
3			Butane, 2-bromo-2-methyl-	150	C5H11Br	000507-36-8	42
4			Butane, 2-iodo-3-methyl-	198	C5H11I	018295-27-7	42
5			Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	40



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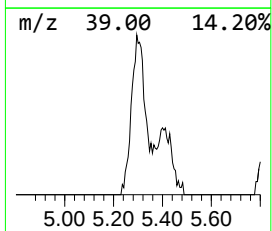
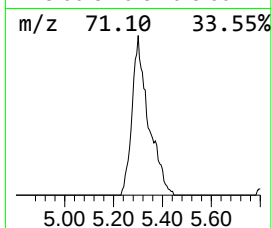
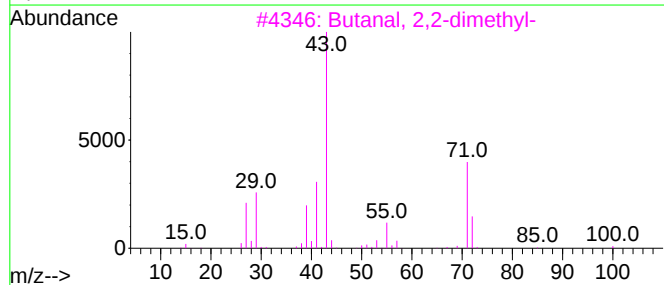
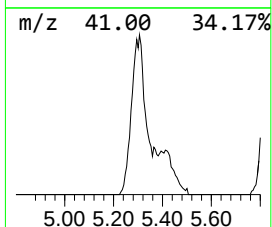
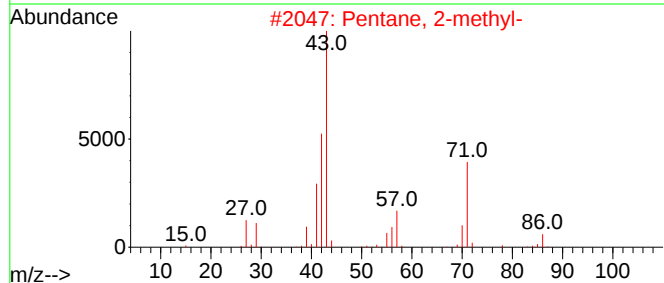
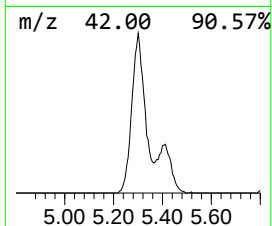
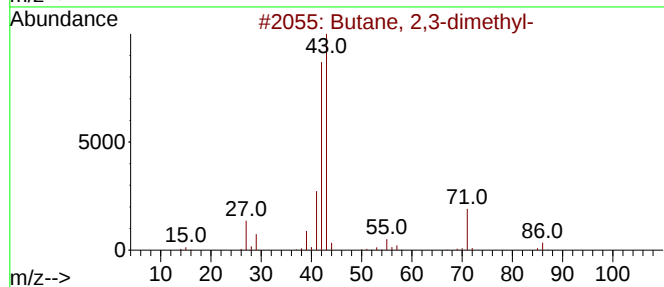
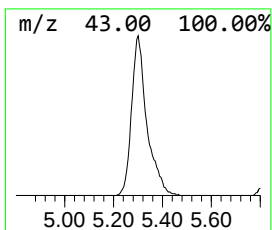
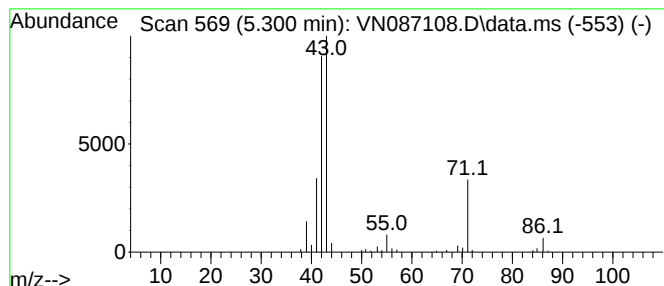
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.300	23.40 ug/l	307740	Pentafluorobenzene	8.229

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	90
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3			Butanal, 2,2-dimethyl-	100	C6H12O	002094-75-9	22
4			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9
5			3-Aminopyrrolidine	86	C4H10N2	079286-79-6	9



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Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

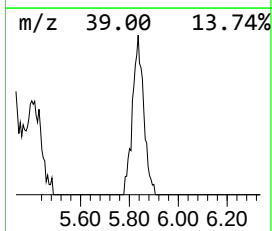
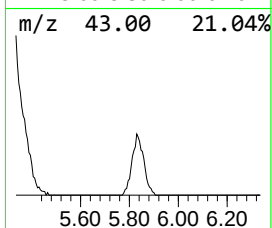
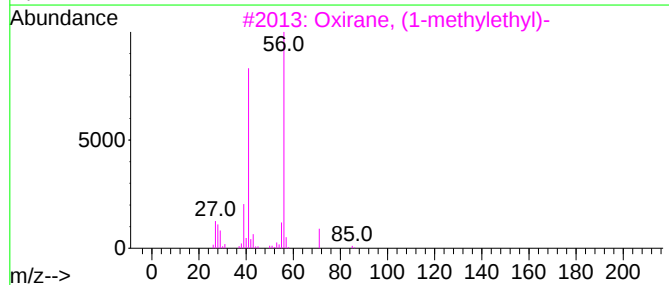
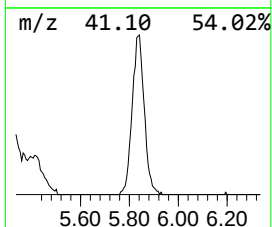
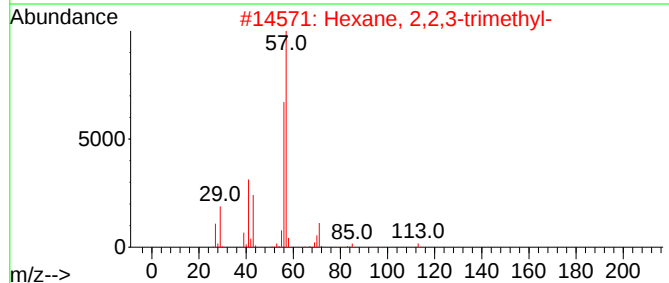
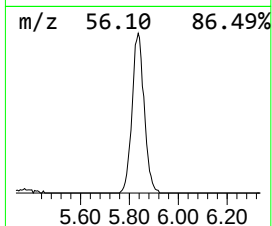
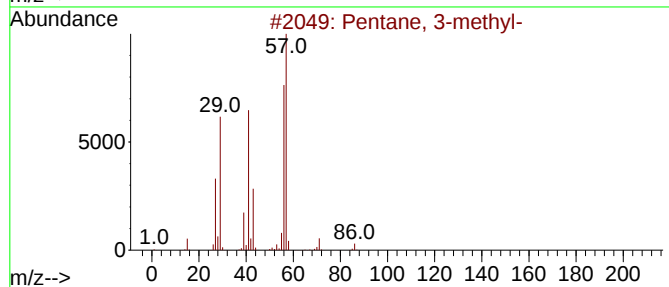
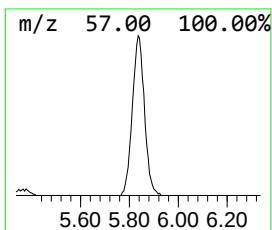
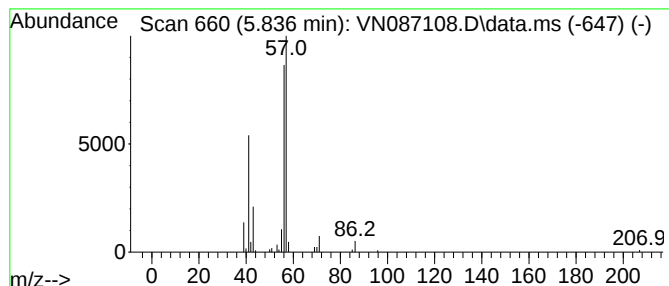
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.836	16.31 ug/l	214490	Pentafluorobenzene	8.229

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
3			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	64
4			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	56
5			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

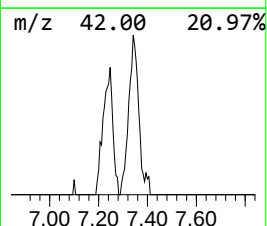
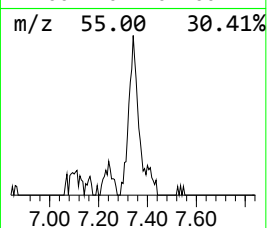
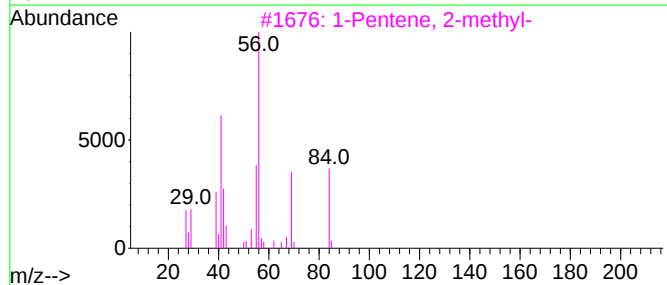
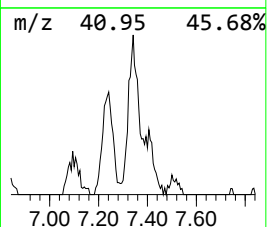
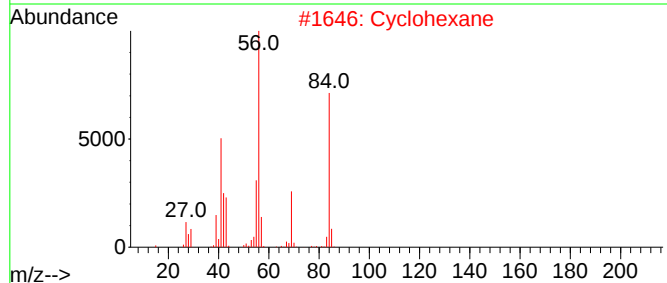
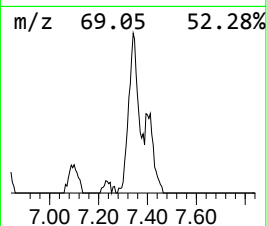
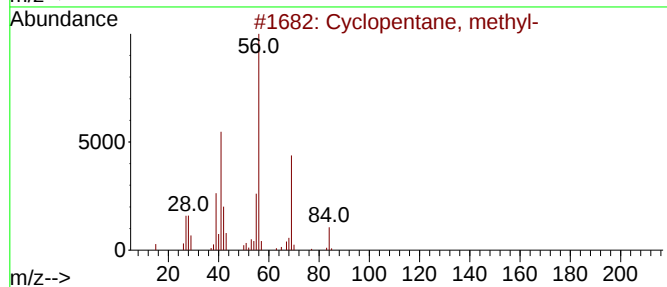
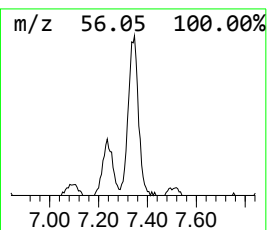
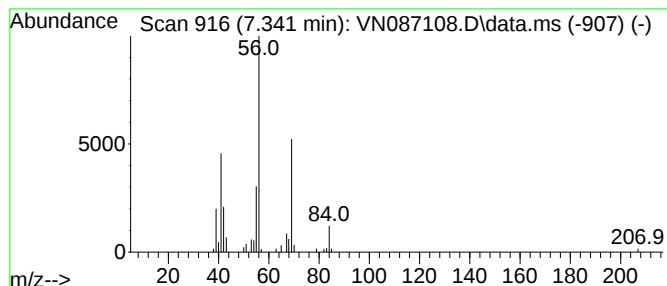
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.341	7.53 ug/l	99028	Pentafluorobenzene	8.229

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	90
3			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
4			Tetramethylene diisocyanate	140	C6H8N2O2	004538-37-8	64
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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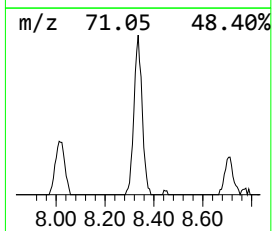
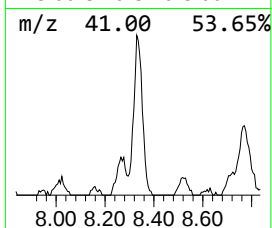
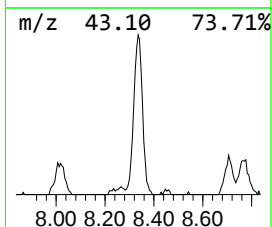
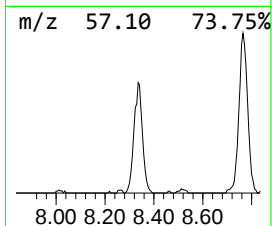
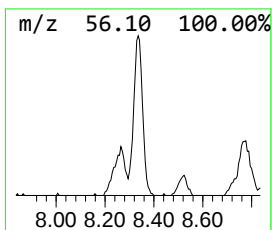
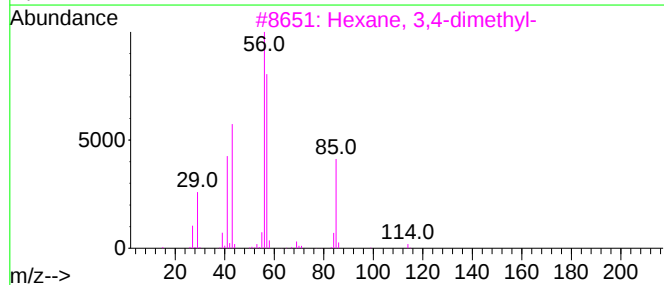
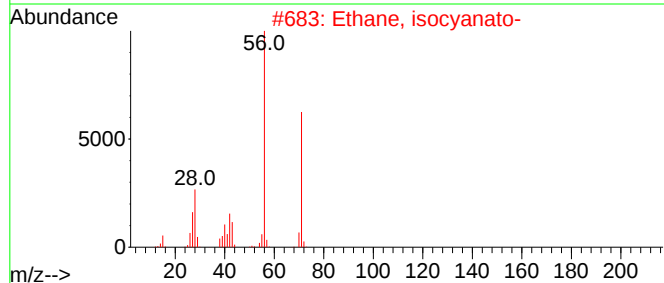
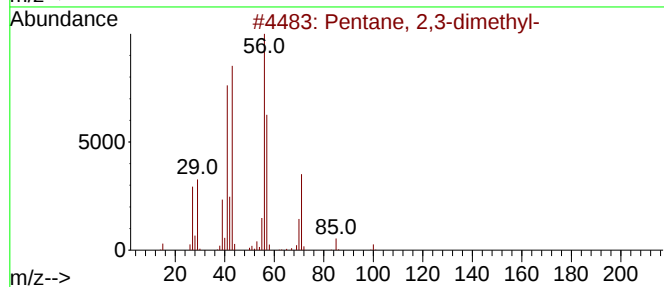
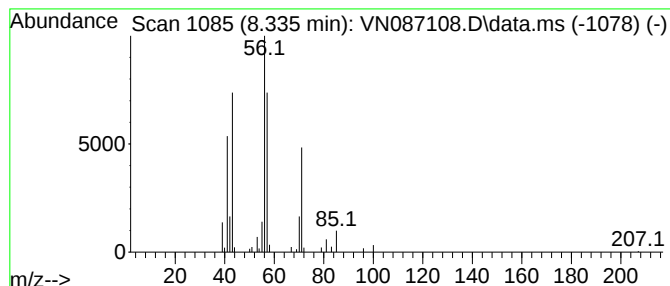
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Quant Title : SW846 8260

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

Peak Number 6 Pentane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.335	13.12 ug/l	172546	Pentafluorobenzene	8.229

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			Ethane, isocyanato-	71	C3H5NO	000109-90-0	50
3			Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	50
4			Hydroxylamine, O-(2-methylpropyl)-	89	C4H11NO	005618-62-2	47
5			Heptane	100	C7H16	000142-82-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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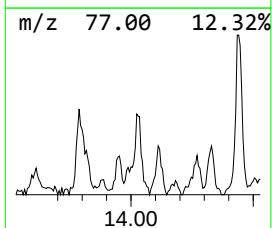
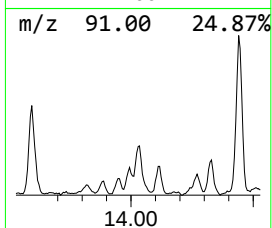
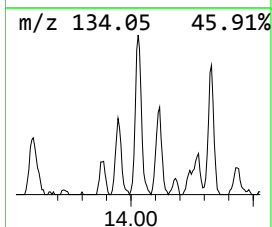
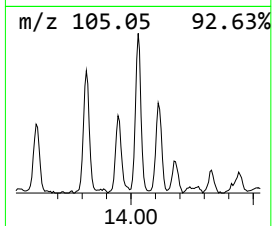
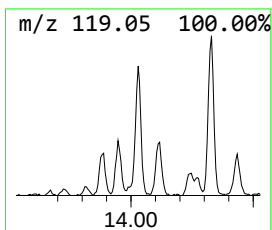
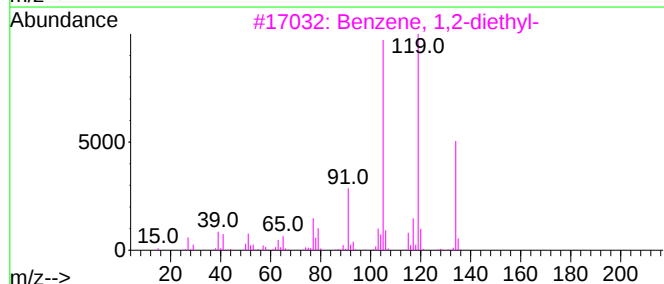
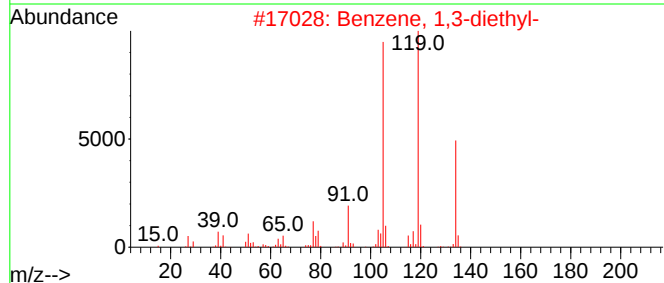
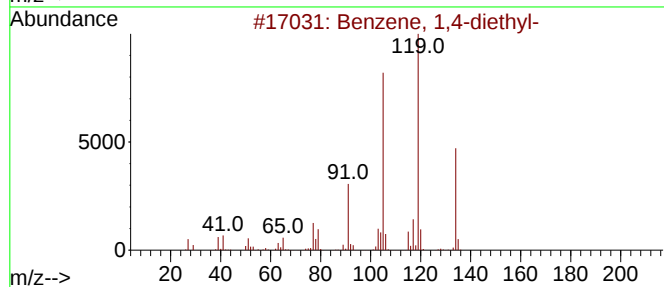
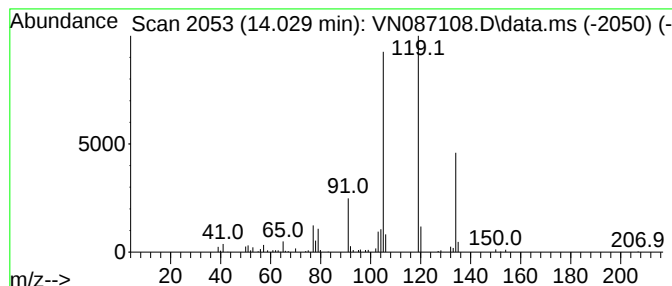
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.029	7.28 ug/l	163752	1,4-Dichlorobenzene-d4	13.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	91
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	80
5			Benzene, tert-butyl-	134	C10H14	000098-06-6	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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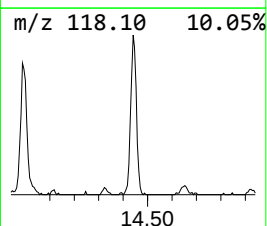
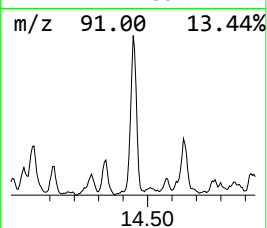
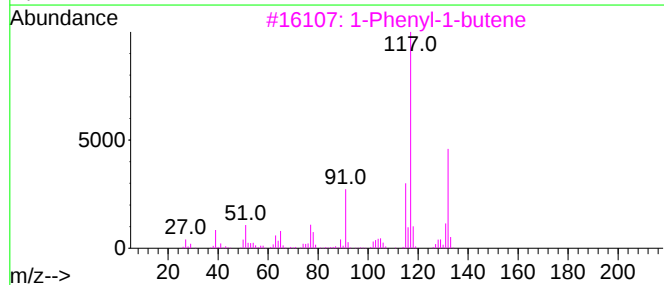
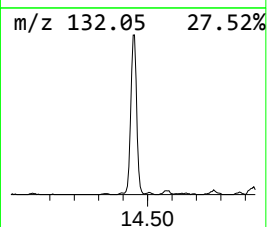
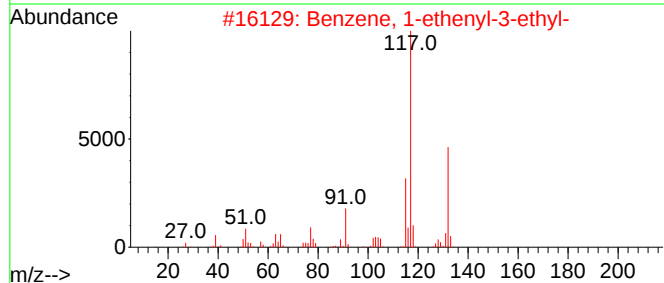
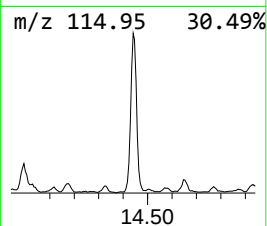
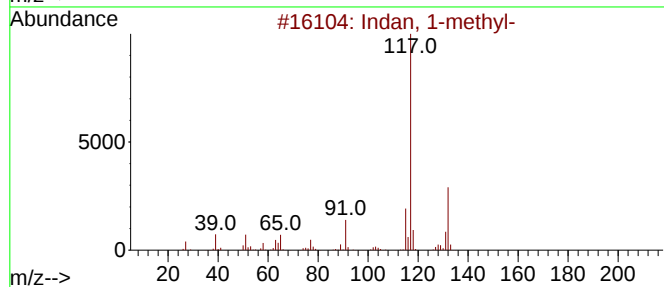
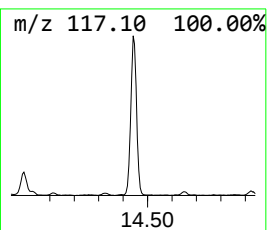
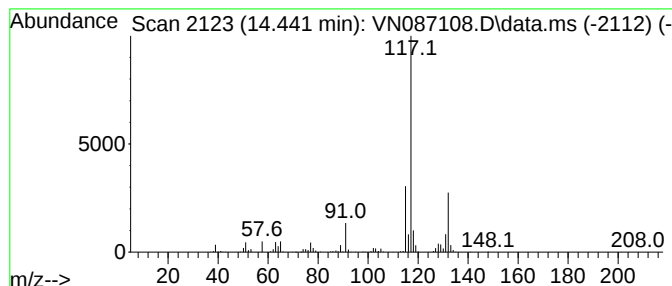
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Quant Title : SW846 8260

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

Peak Number 8 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.441	34.86 ug/l	783874	1,4-Dichlorobenzene-d4	13.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

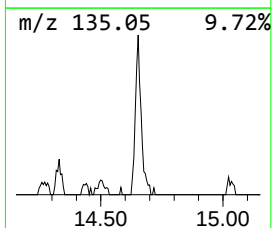
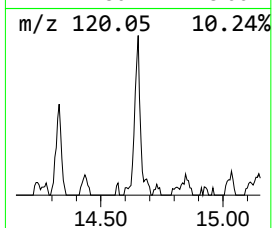
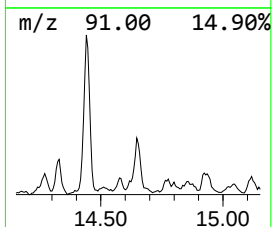
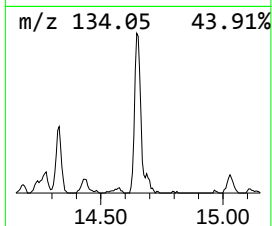
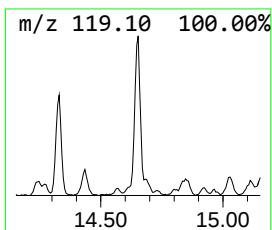
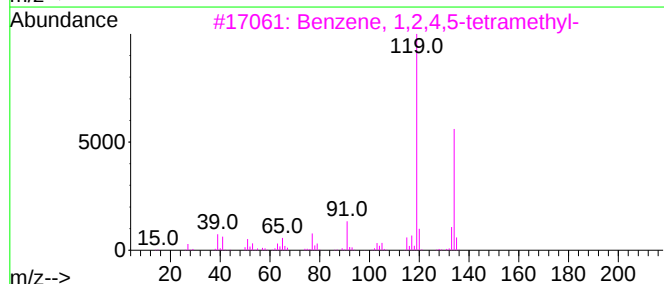
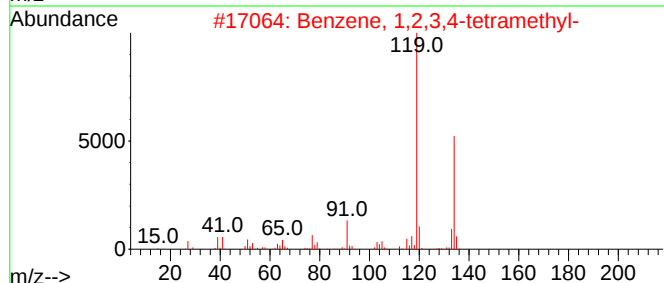
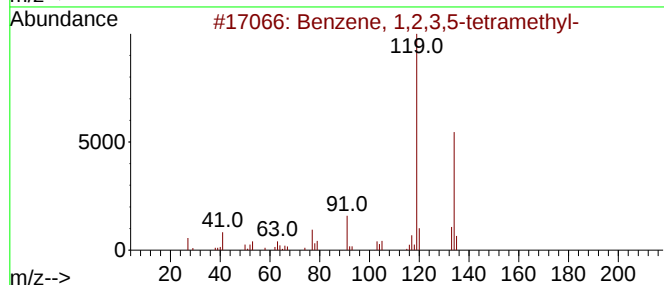
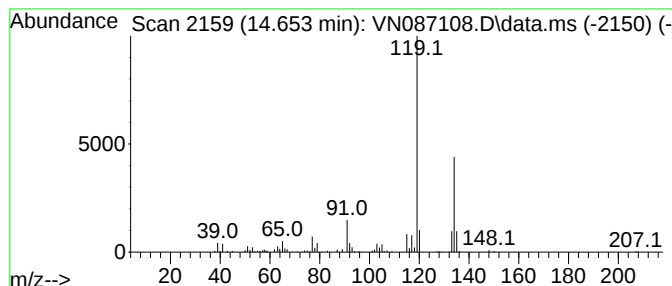
Instrument :
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ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.653	9.60 ug/l	215878	1,4-Dichlorobenzene-d4	13.788	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	91
5	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

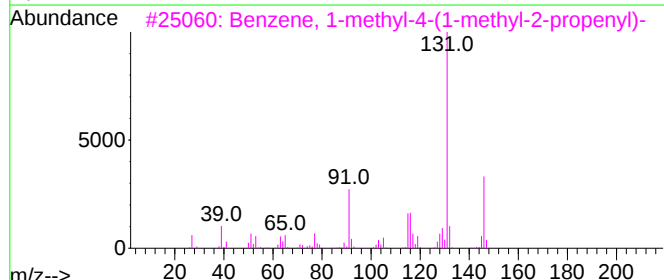
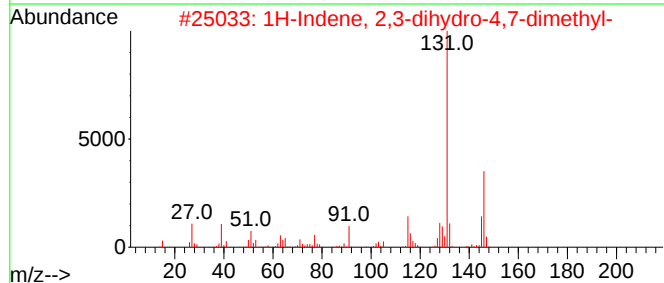
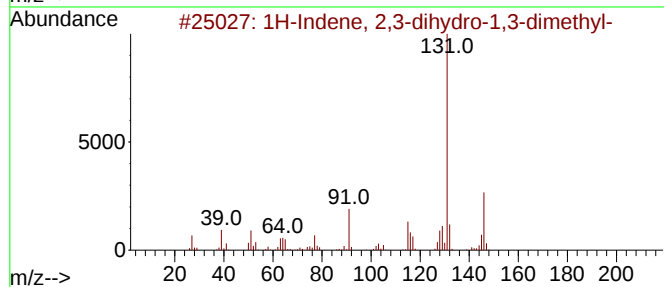
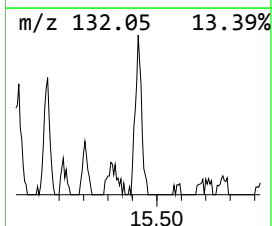
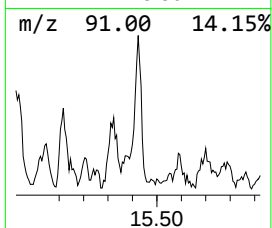
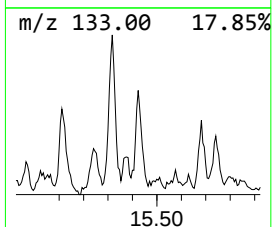
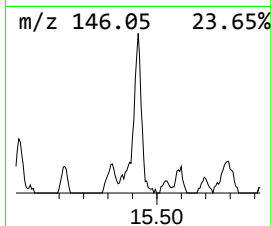
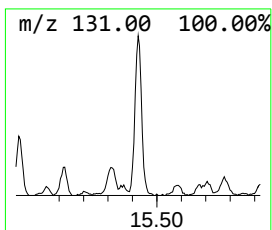
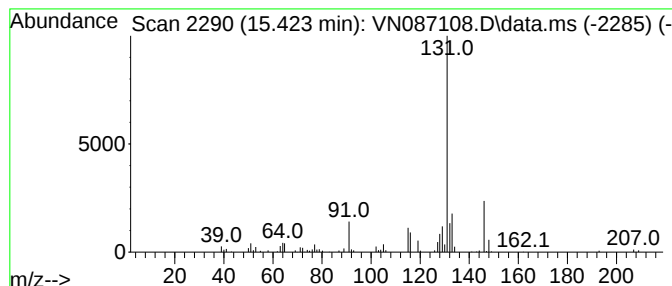
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.423	7.28 ug/l	163725	1,4-Dichlorobenzene-d4	13.788

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
3			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087108.D
Acq On : 19 Jun 2025 13:59
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.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-	3.283	35.3	ug/l	464065	1	8.229	657426	50.0
Butane, 2,2-dim...	4.447	9.5	ug/l	125007	1	8.229	657426	50.0
Butane, 2,3-dim...	5.300	23.4	ug/l	307740	1	8.229	657426	50.0
Pentane, 3-methyl-	5.836	16.3	ug/l	214490	1	8.229	657426	50.0
Cyclopentane, m...	7.341	7.5	ug/l	99028	1	8.229	657426	50.0
Pentane, 2,3-di...	8.335	13.1	ug/l	172546	1	8.229	657426	50.0
Benzene, 1,4-di...	14.029	7.3	ug/l	163752	4	13.788	1124160	50.0
Indan, 1-methyl-	14.441	34.9	ug/l	783874	4	13.788	1124160	50.0
Benzene, 1,2,3,...	14.653	9.6	ug/l	215878	4	13.788	1124160	50.0
1H-Indene, 2,3-...	15.423	7.3	ug/l	163725	4	13.788	1124160	50.0