

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
 Data File : VN088118.D
 Acq On : 27 Oct 2025 16:37
 Operator : JC\MD
 Sample : Q3426-02 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TWP3

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: VN088118.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.265	215	223	232	rBV4	16091	41186	1.56%	0.152%
2	3.659	282	290	291	rBV4	14594	28655	1.09%	0.106%
3	4.148	362	373	384	rBV3	38001	119158	4.52%	0.441%
4	5.147	532	543	551	rBV7	11828	38042	1.44%	0.141%
5	5.353	568	578	579	rBV6	16123	43526	1.65%	0.161%
6	5.800	642	654	666	rVB4	17262	60515	2.29%	0.224%
7	6.106	693	706	717	rBV7	8893	29178	1.11%	0.108%
8	6.306	729	740	750	rBV3	14228	42052	1.59%	0.156%
9	6.636	785	796	809	rBV4	35010	108844	4.13%	0.403%
10	6.783	809	821	830	rBV6	21683	65403	2.48%	0.242%
11	7.065	860	869	880	rBV4	29542	84325	3.20%	0.312%
12	7.312	900	911	920	rBV	108022	321713	12.20%	1.191%
13	7.418	926	929	942	rVB	14037	33843	1.28%	0.125%
14	7.988	1018	1026	1036	rVB2	66814	158650	6.01%	0.587%
15	8.147	1043	1053	1058	rBV	218355	539359	20.45%	1.997%
16	8.206	1058	1063	1078	rVV2	365520	1171117	44.40%	4.335%
17	8.306	1078	1080	1090	rVB6	13353	31602	1.20%	0.117%
18	8.430	1094	1101	1106	rBV5	12608	31229	1.18%	0.116%
19	8.588	1115	1128	1140	rBV2	964871	2637932	100.00%	9.766%
20	8.718	1140	1150	1156	rBV5	43532	126626	4.80%	0.469%
21	8.841	1165	1171	1179	rVB4	21872	50105	1.90%	0.185%
22	8.941	1179	1188	1198	rVB4	10303	31253	1.18%	0.116%
23	9.083	1203	1212	1225	rBV	656855	1408229	53.38%	5.213%
24	9.365	1255	1260	1270	rVB2	19334	41478	1.57%	0.154%
25	9.577	1280	1296	1312	rBV	159908	396542	15.03%	1.468%
26	9.759	1320	1327	1334	rBV5	17497	31482	1.19%	0.117%
27	9.924	1348	1355	1356	rBV3	14442	27090	1.03%	0.100%
28	10.082	1375	1382	1386	rBV2	54134	101204	3.84%	0.375%
29	10.130	1387	1390	1396	rVB4	16990	29788	1.13%	0.110%
30	10.318	1414	1422	1429	rBV7	13734	35763	1.36%	0.132%
31	10.429	1435	1441	1447	rBV2	45107	87397	3.31%	0.324%
32	10.547	1453	1461	1467	rBV	889774	1643329	62.30%	6.084%
33	10.606	1467	1471	1477	rVB	75752	144719	5.49%	0.536%
34	10.741	1485	1494	1499	rVB3	69000	130364	4.94%	0.483%
35	10.812	1500	1506	1513	rVB2	65189	118902	4.51%	0.440%
36	10.965	1525	1532	1533	rBV5	20710	32615	1.24%	0.121%

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37	11.394	1601	1605	1610	rVB4	17625	31881	1.21%	0.118%
38	11.841	1674	1681	1692	rBV	858052	1612824	61.14%	5.971%
39	11.941	1692	1698	1708	rVB	304738	548449	20.79%	2.030%
40	12.053	1708	1717	1729	rBV	952298	1841107	69.79%	6.816%
41	12.376	1762	1772	1782	rBV3	39866	106597	4.04%	0.395%
42	12.559	1790	1803	1808	rBV8	13358	44028	1.67%	0.163%
43	12.671	1816	1822	1832	rBV	81517	144703	5.49%	0.536%
44	12.829	1842	1849	1860	rBV	593561	1099065	41.66%	4.069%
45	13.012	1873	1880	1886	rBV	165016	296072	11.22%	1.096%
46	13.076	1886	1891	1892	rVV	62914	91831	3.48%	0.340%
47	13.106	1892	1896	1900	rVV	187882	359460	13.63%	1.331%
48	13.147	1900	1903	1910	rVB	365226	593273	22.49%	2.196%
49	13.312	1923	1931	1939	rBV	310136	549770	20.84%	2.035%
50	13.459	1949	1956	1970	rBV	636895	1039860	39.42%	3.850%
51	13.594	1970	1979	1983	rVB4	23666	48589	1.84%	0.180%
52	13.653	1983	1989	1993	rBV3	36363	60926	2.31%	0.226%
53	13.706	1993	1998	2002	rBV4	17879	30942	1.17%	0.115%
54	13.770	2002	2009	2024	rBV2	927077	2244380	85.08%	8.309%
55	13.929	2031	2036	2037	rBV2	50430	68493	2.60%	0.254%
56	13.970	2037	2043	2058	rVB2	609445	1376953	52.20%	5.097%
57	14.094	2058	2064	2070	rBV7	39051	65633	2.49%	0.243%
58	14.159	2070	2075	2080	rBV	54393	87969	3.33%	0.326%
59	14.223	2080	2086	2088	rBV	61185	97679	3.70%	0.362%
60	14.306	2095	2100	2107	rBV	155299	233873	8.87%	0.866%
61	14.370	2107	2111	2114	rVV3	42414	65160	2.47%	0.241%
62	14.417	2114	2119	2127	rVB2	186887	330781	12.54%	1.225%
63	14.547	2136	2141	2146	rBV4	43994	79789	3.02%	0.295%
64	14.629	2146	2155	2158	rVV	99351	210971	8.00%	0.781%
65	14.665	2158	2161	2166	rVV	114334	191241	7.25%	0.708%
66	14.706	2166	2168	2172	rVV4	39670	56276	2.13%	0.208%
67	14.753	2172	2176	2183	rVB5	21437	48519	1.84%	0.180%
68	14.853	2190	2193	2196	rVB4	22328	27841	1.06%	0.103%
69	14.906	2196	2202	2213	rVB2	156301	335575	12.72%	1.242%
70	15.029	2213	2223	2230	rBV2	340157	804965	30.52%	2.980%
71	15.182	2243	2249	2256	rBV	200447	380766	14.43%	1.410%
72	15.294	2262	2268	2272	rBV3	45192	86604	3.28%	0.321%
73	15.412	2280	2288	2296	rVB3	69457	186266	7.06%	0.690%
74	15.612	2316	2322	2328	rVV	275030	523955	19.86%	1.940%
75	15.670	2328	2332	2336	rVB2	36718	56694	2.15%	0.210%
76	15.723	2336	2341	2343	rBV3	28715	48919	1.85%	0.181%

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Title : SW846 8260

77	15.917	2365	2374	2380	rBV2	48318	123821	4.69%	0.458%
78	16.100	2398	2405	2406	rBV2	33099	58001	2.20%	0.215%
79	16.135	2406	2411	2421	rVB2	110351	235938	8.94%	0.873%
80	16.306	2436	2440	2446	rVB4	22222	50661	1.92%	0.188%
81	16.470	2459	2468	2481	rBV4	71931	193798	7.35%	0.717%
82	16.670	2495	2502	2509	rBV5	42268	107108	4.06%	0.397%
83	16.747	2509	2515	2522	rVV3	47926	111434	4.22%	0.413%

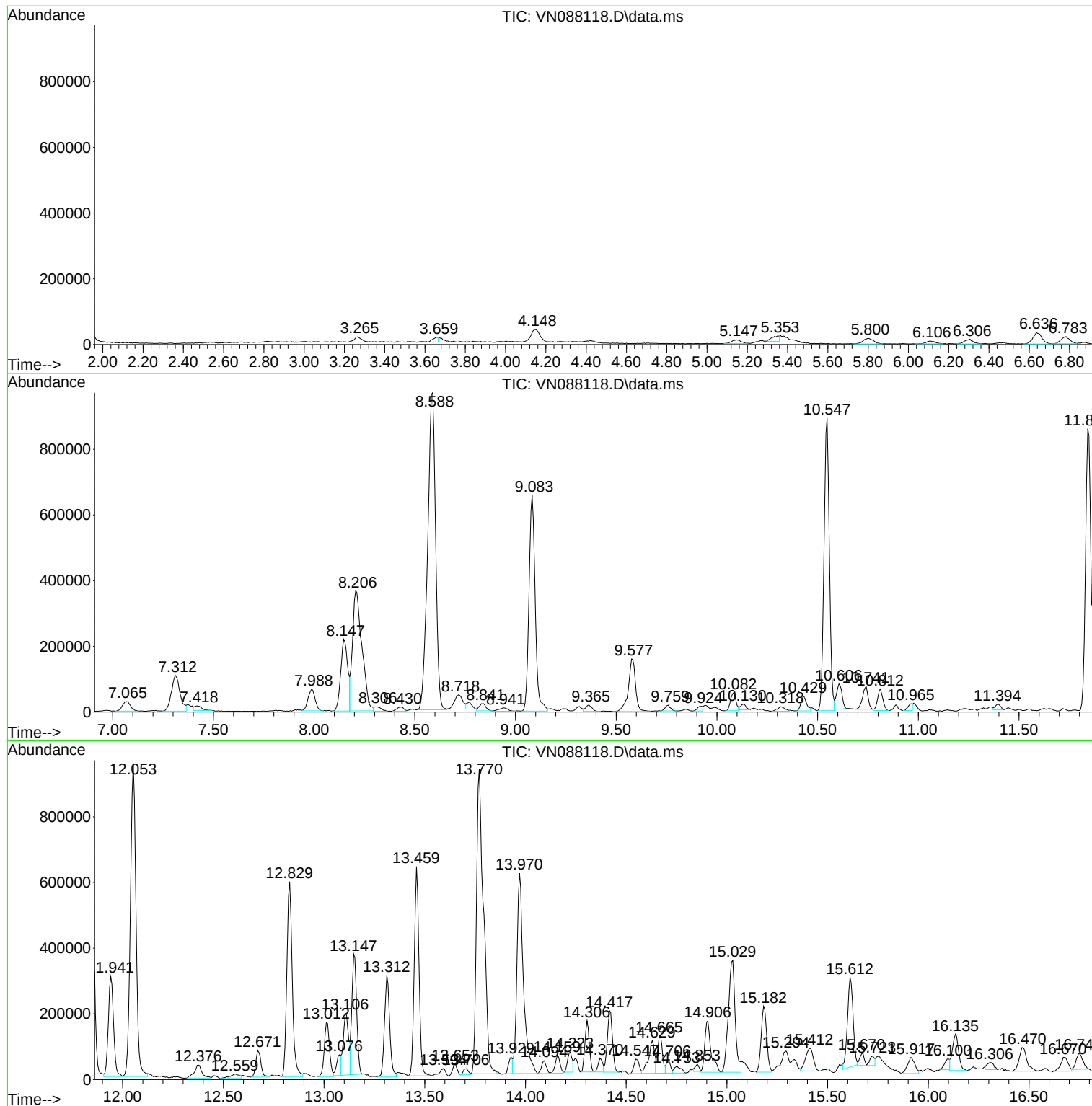
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
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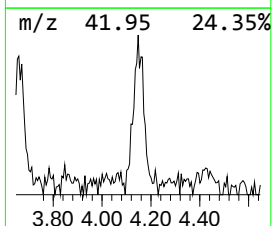
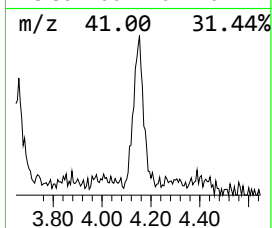
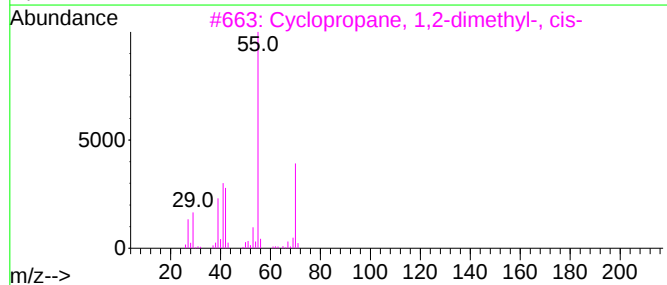
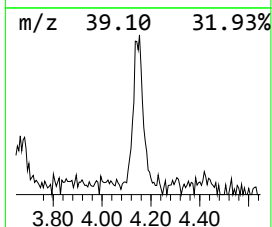
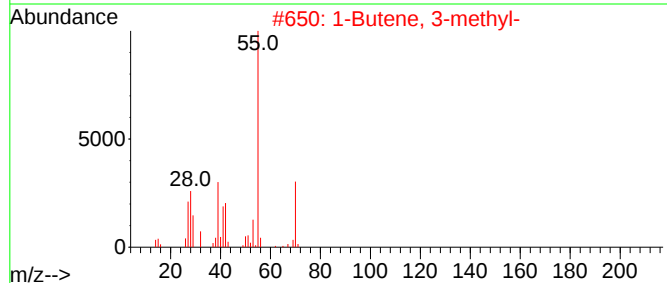
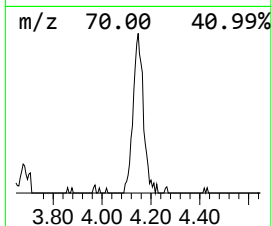
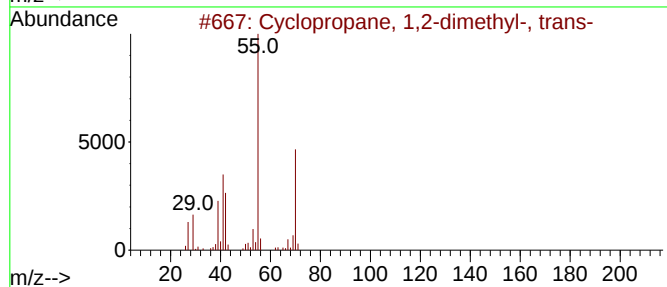
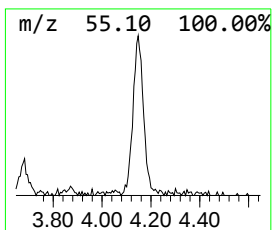
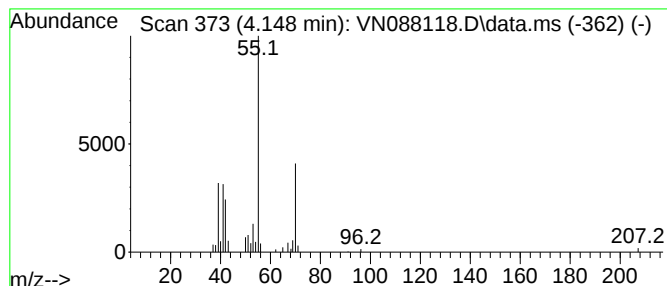
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopropane, 1,2-dimethyl-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.148	5.09 ug/l	119158	Pentafluorobenzene	8.206

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87
2			1-Butene, 3-methyl-	70	C5H10	000563-45-1	86
3			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	86
4			2-Butene, 2-methyl-	70	C5H10	000513-35-9	80
5			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80



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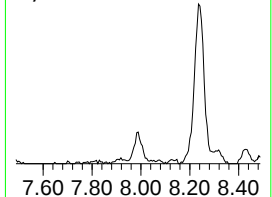
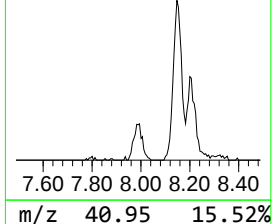
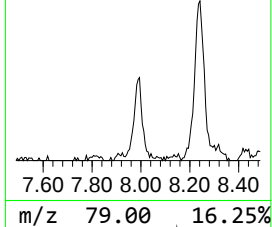
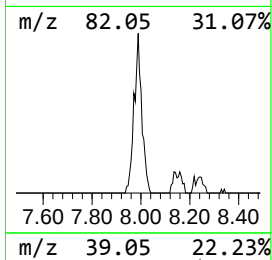
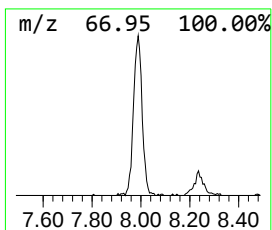
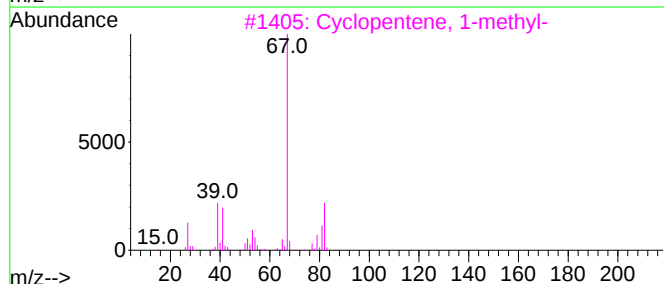
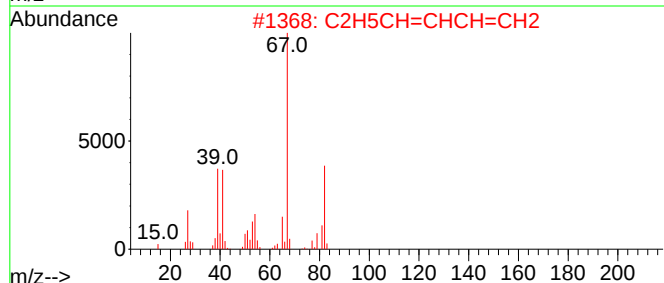
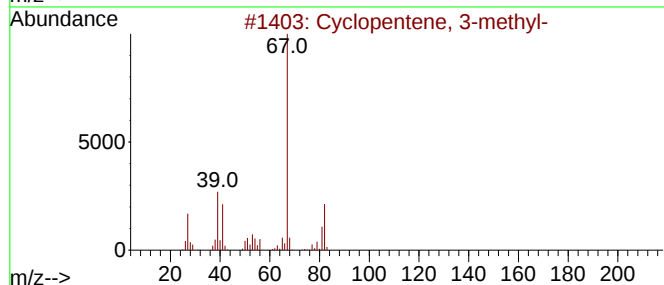
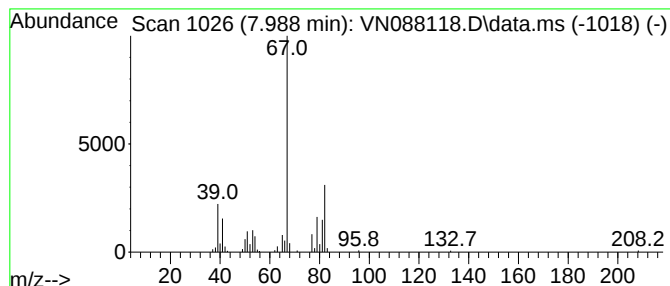
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Cyclopentene, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.988	6.77 ug/l	158650	Pentafluorobenzene	8.206

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	86
2		C2H5CH=CHCH=CH2	82	C6H10	000592-48-3	83
3		Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	83
4		Cyclohexene	82	C6H10	000110-83-8	80
5		Cyclopentane, methylene-	82	C6H10	001528-30-9	80



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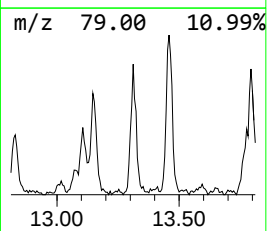
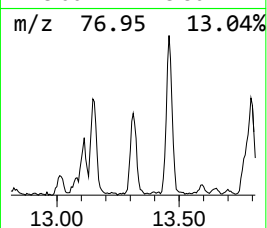
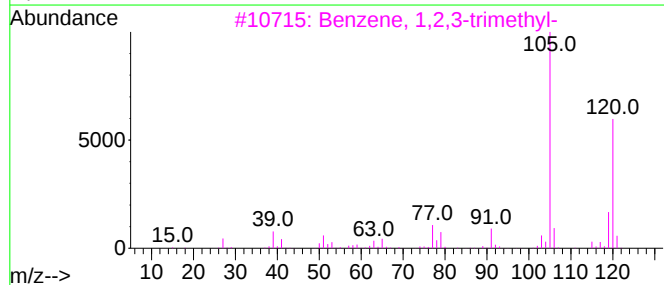
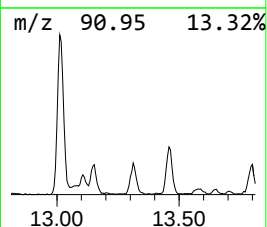
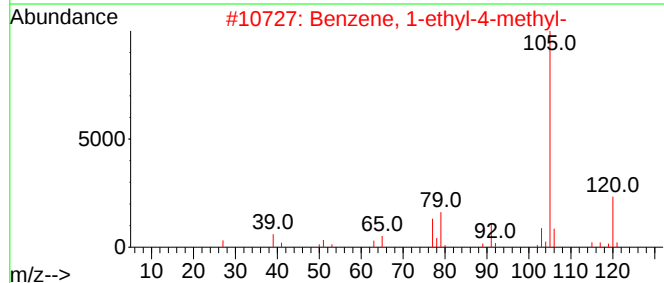
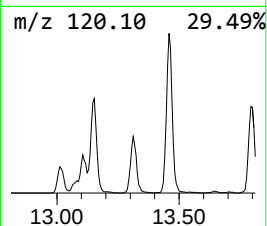
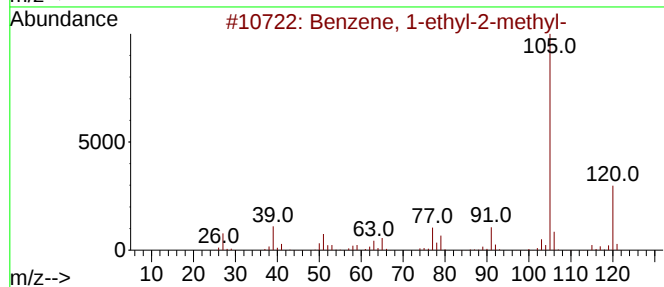
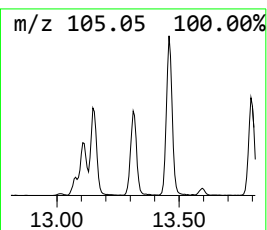
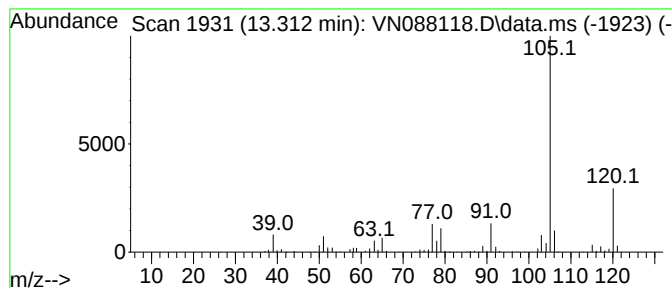
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.312	12.25 ug/l	549770	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	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



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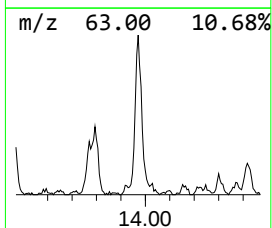
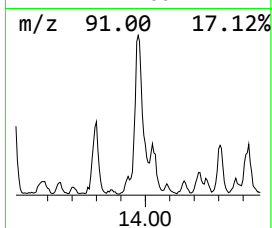
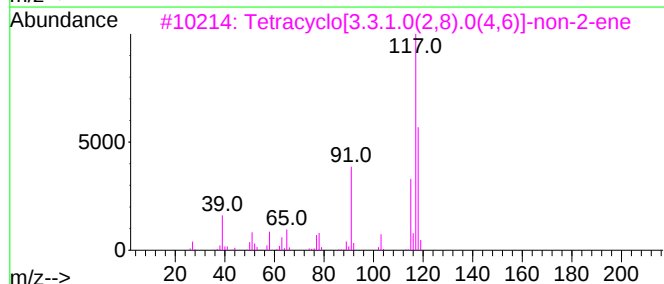
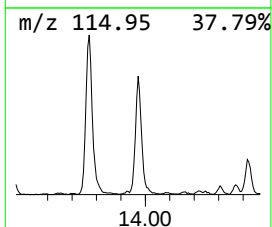
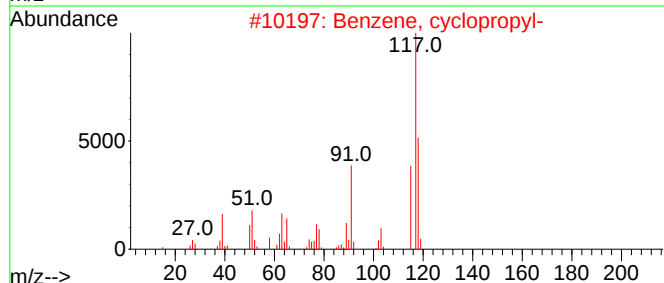
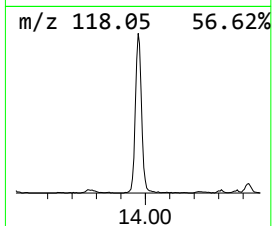
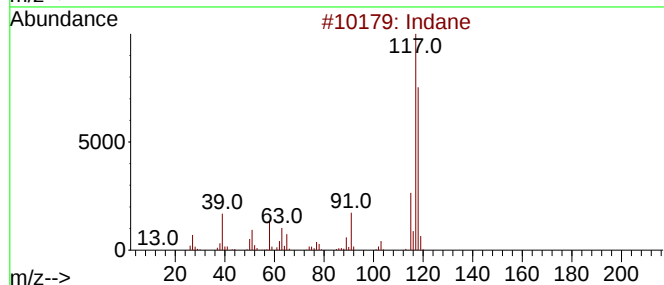
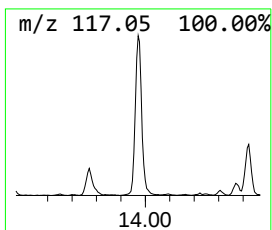
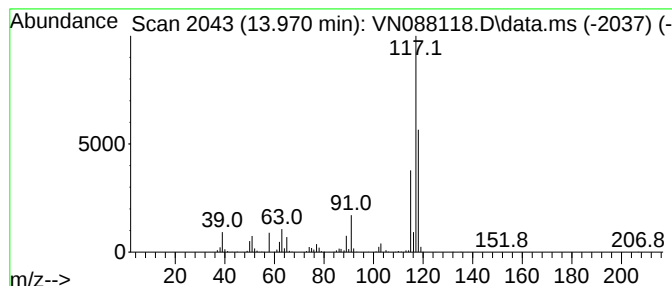
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Quant Title : SW846 8260

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

Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.970	30.68 ug/l	1376950	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	91
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	68
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
Data File : VN088118.D
Acq On : 27 Oct 2025 16:37
Operator : JC\MD
Sample : Q3426-02 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TWP3

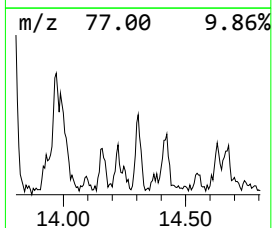
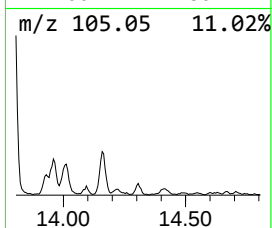
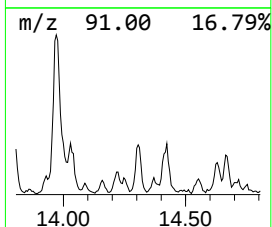
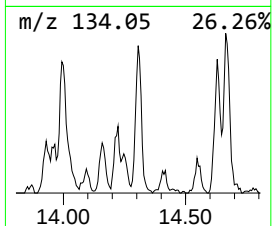
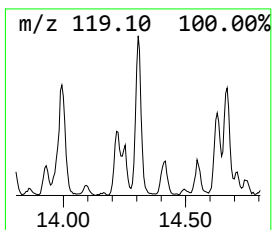
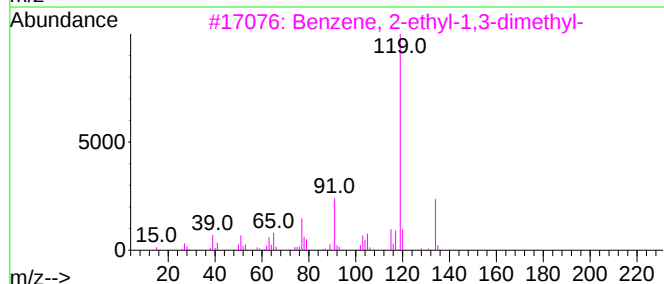
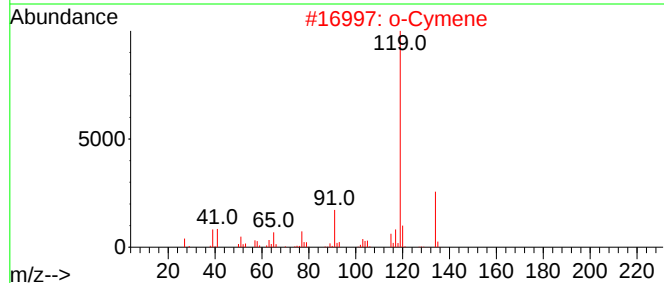
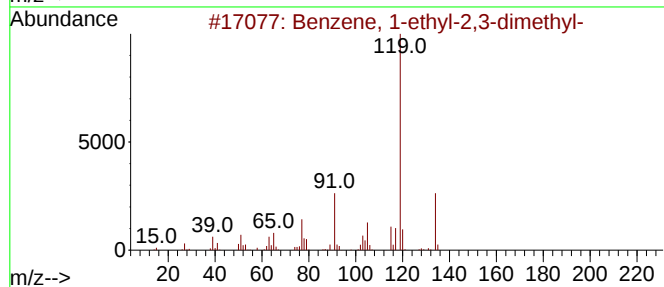
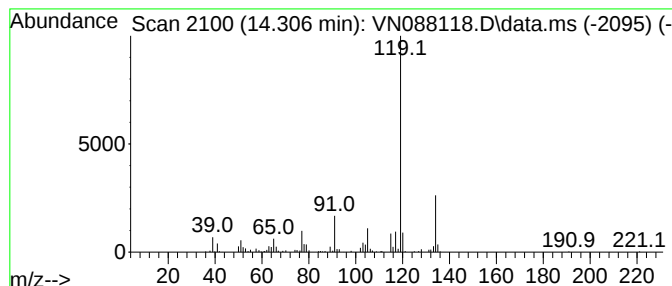
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.306	5.21 ug/l	233873	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2			o-Cymene	134	C10H14	000527-84-4	94
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
Data File : VN088118.D
Acq On : 27 Oct 2025 16:37
Operator : JC\MD
Sample : Q3426-02 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TWP3

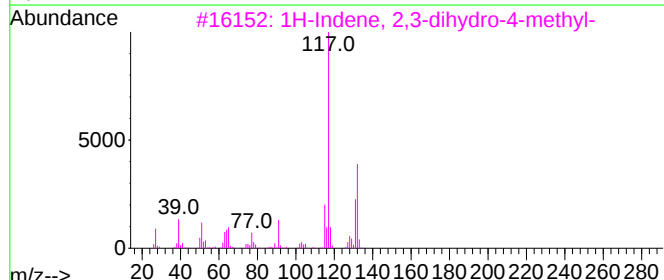
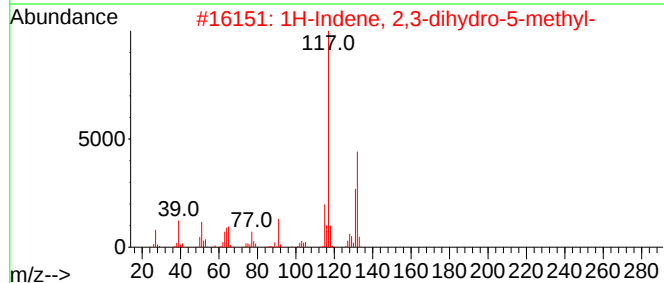
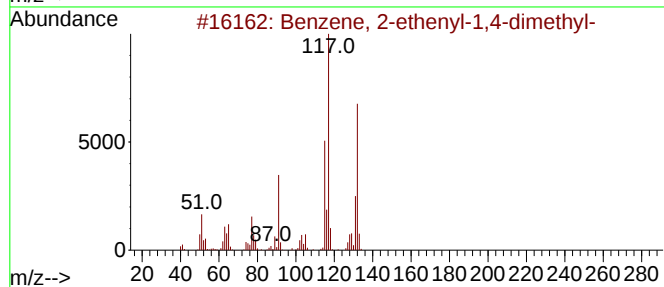
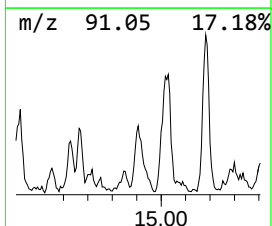
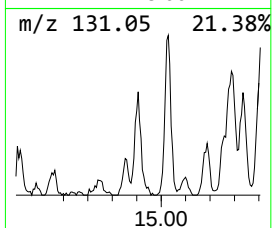
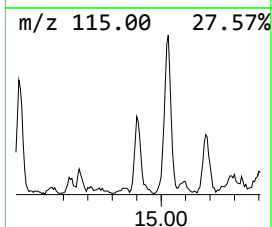
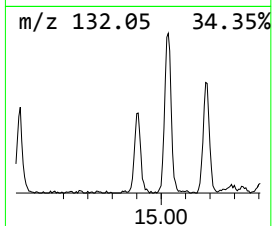
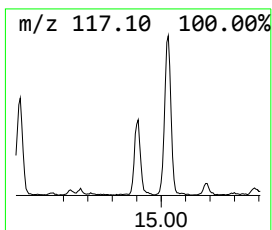
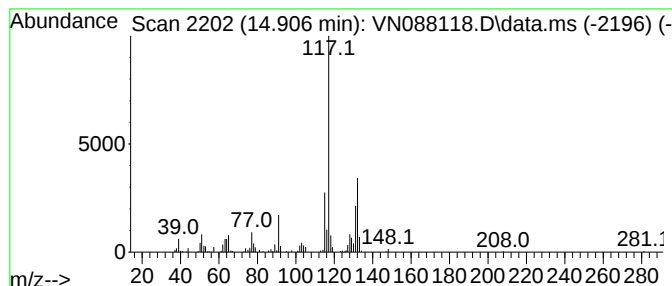
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.906	7.48 ug/l	335575	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	92
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
Data File : VN088118.D
Acq On : 27 Oct 2025 16:37
Operator : JC\MD
Sample : Q3426-02 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TWP3

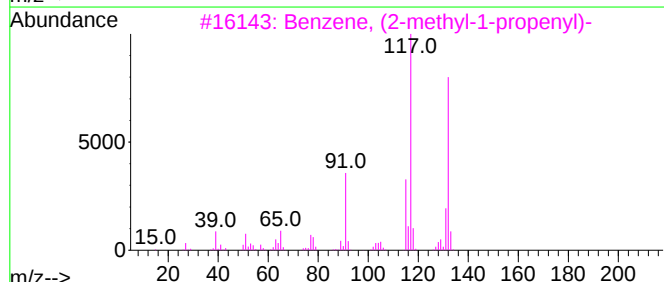
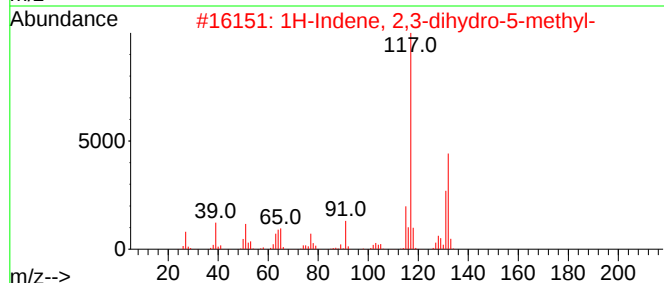
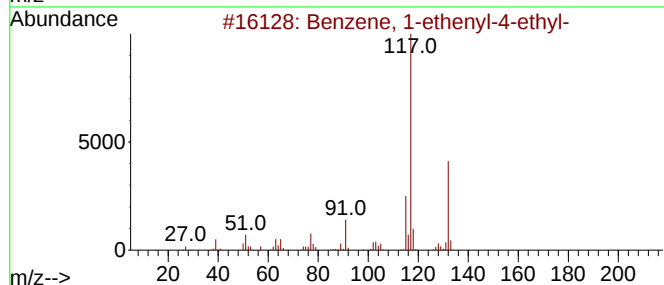
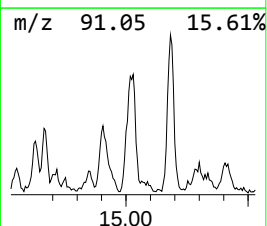
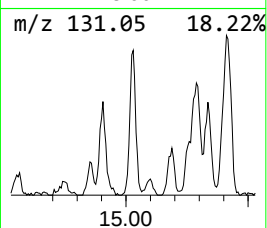
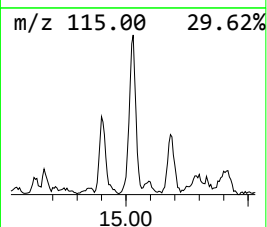
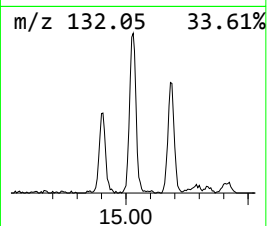
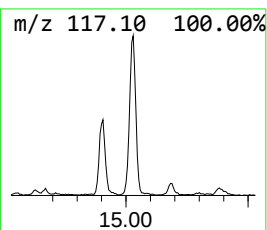
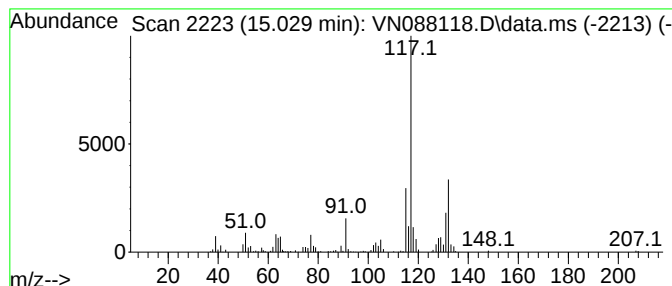
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.029	17.93 ug/l	804965	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
4			Indan, 1-methyl-	132	C10H12	000767-58-8	90
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
Data File : VN088118.D
Acq On : 27 Oct 2025 16:37
Operator : JC\MD
Sample : Q3426-02 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TWP3

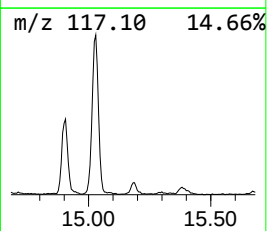
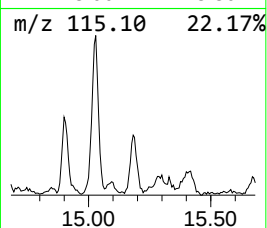
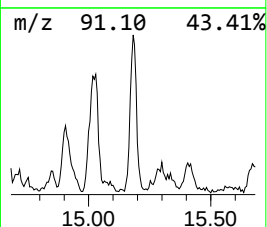
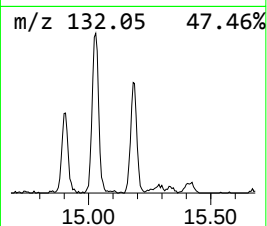
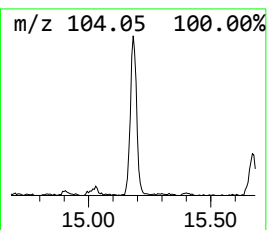
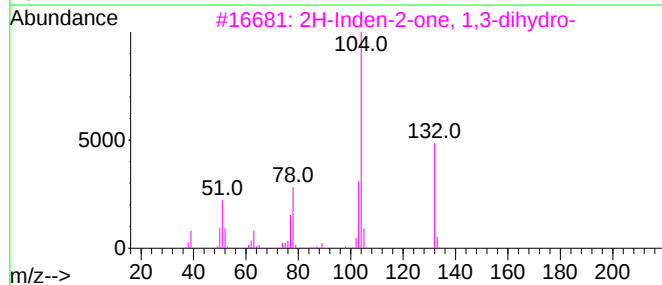
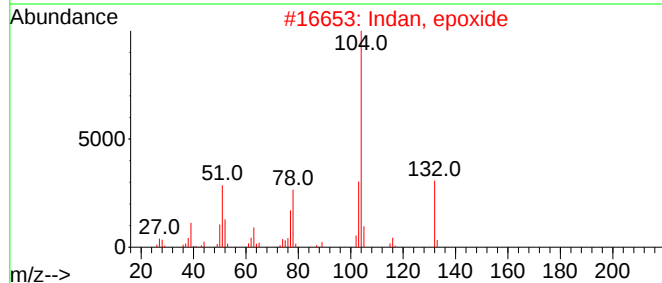
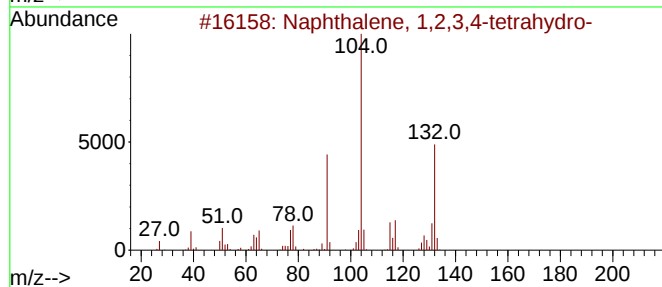
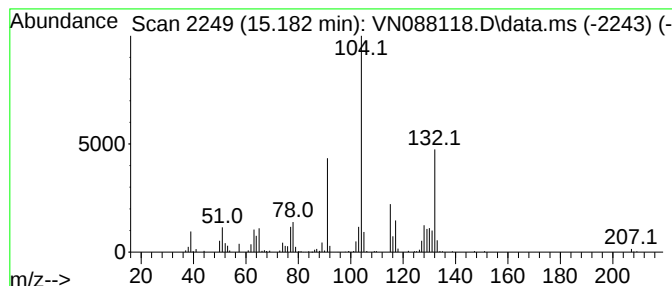
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.182	8.48 ug/l	380766	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Indan, epoxide	132	C9H8O	000768-22-9	49
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	49
4			5,5-Dimethyl-4-phenyldihydrofura...	190	C12H14O2	013133-96-5	38
5			(3-Phenylpropionyl)glycine	207	C11H13NO3	056613-60-6	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
Data File : VN088118.D
Acq On : 27 Oct 2025 16:37
Operator : JC\MD
Sample : Q3426-02 10X
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TWP3

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
Cyclopropane, 1...	4.148	5.1	ug/l	119158	1	8.206	1171120	50.0
Cyclopentene, 3...	7.988	6.8	ug/l	158650	1	8.206	1171120	50.0
Benzene, 1-ethy...	13.312	12.3	ug/l	549770	4	13.771	2244380	50.0
Indane	13.970	30.7	ug/l	1376950	4	13.771	2244380	50.0
Benzene, 1-ethy...	14.306	5.2	ug/l	233873	4	13.771	2244380	50.0
Benzene, 2-ethe...	14.906	7.5	ug/l	335575	4	13.771	2244380	50.0
Benzene, 1-ethe...	15.029	17.9	ug/l	804965	4	13.771	2244380	50.0
Naphthalene, 1,...	15.182	8.5	ug/l	380766	4	13.771	2244380	50.0