

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081822\
 Data File : VN074000.D
 Acq On : 18 Aug 2022 13:20
 Operator : JC\MD
 Sample : N4204-24DL2 100X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1DL2

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 1 % 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\82N081522W.M
 Title : SW846 8260

Signal : TIC: VN074000.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.087	197	204	211	rVB	37567	77535	2.30%	0.260%
2	3.481	260	271	283	rVB3	37981	114740	3.40%	0.385%
3	3.657	290	301	312	rBV3	38637	99115	2.94%	0.333%
4	3.828	320	330	338	rBV3	21800	58120	1.72%	0.195%
5	3.928	338	347	360	rVB	99891	271528	8.05%	0.911%
6	4.881	498	509	521	rBV2	32585	101390	3.01%	0.340%
7	5.110	536	548	570	rVB3	23884	106369	3.15%	0.357%
8	5.534	609	620	632	rBV6	12807	46533	1.38%	0.156%
9	6.387	758	765	775	rVB2	20801	55909	1.66%	0.188%
10	6.534	778	790	796	rBV4	11428	35644	1.06%	0.120%
11	6.816	830	838	850	rVB4	16254	48636	1.44%	0.163%
12	7.069	867	881	889	rBV4	25148	75466	2.24%	0.253%
13	7.240	900	910	918	rBV2	32657	89256	2.65%	0.300%
14	7.769	992	1000	1010	rVB2	48312	119097	3.53%	0.400%
15	7.951	1021	1031	1036	rBV	415077	986678	29.25%	3.312%
16	8.016	1036	1042	1054	rVB	792799	1800392	53.38%	6.043%
17	8.375	1092	1103	1115	rBV2	440458	1176449	34.88%	3.949%
18	8.522	1120	1128	1136	rVV4	21518	53770	1.59%	0.180%
19	8.904	1184	1193	1208	rBV	1016800	2146160	63.63%	7.204%
20	10.380	1436	1444	1449	rBV	1434423	2566169	76.08%	8.613%
21	10.445	1449	1455	1470	rVB	1901454	3373009	100.00%	11.322%
22	11.686	1658	1666	1675	rBV	1516030	2622194	77.74%	8.801%
23	11.786	1675	1683	1692	rVV	1033886	1719198	50.97%	5.771%
24	11.892	1693	1701	1715	rVB	1261418	2403124	71.25%	8.066%
25	12.222	1746	1757	1766	rBV	390632	648738	19.23%	2.178%
26	12.522	1800	1808	1815	rBV	178967	293986	8.72%	0.987%
27	12.674	1826	1834	1845	rBV	1312932	2217290	65.74%	7.442%
28	12.863	1859	1866	1871	rBV	183504	290831	8.62%	0.976%
29	12.921	1871	1876	1879	rVV	129729	207446	6.15%	0.696%
30	12.957	1879	1882	1885	rVV	136377	201604	5.98%	0.677%
31	12.998	1885	1889	1896	rVB	225234	364786	10.81%	1.224%
32	13.163	1911	1917	1924	rVB	96323	159811	4.74%	0.536%
33	13.310	1933	1942	1952	rBV	628377	1004274	29.77%	3.371%
34	13.616	1986	1994	2007	rBV	1688098	3087071	91.52%	10.362%
35	13.816	2023	2028	2044	rVB	260961	496988	14.73%	1.668%
36	14.004	2054	2060	2066	rVB2	23275	42745	1.27%	0.143%

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37	14.068	2066	2071	2075	rBV3	19177	36072	1.07%	0.121%
38	14.157	2080	2086	2092	rVB	43921	68985	2.05%	0.232%
39	14.268	2099	2105	2112	rVB3	26967	44748	1.33%	0.150%
40	14.480	2136	2141	2144	rBV	27532	42990	1.27%	0.144%
41	14.516	2144	2147	2153	rVB	26018	39784	1.18%	0.134%
42	14.751	2181	2187	2194	rBV	22498	37794	1.12%	0.127%
43	14.868	2199	2207	2214	rBV4	48316	101134	3.00%	0.339%
44	15.433	2293	2303	2311	rBV	109289	215081	6.38%	0.722%
45	16.533	2481	2490	2496	rBV	19506	44127	1.31%	0.148%

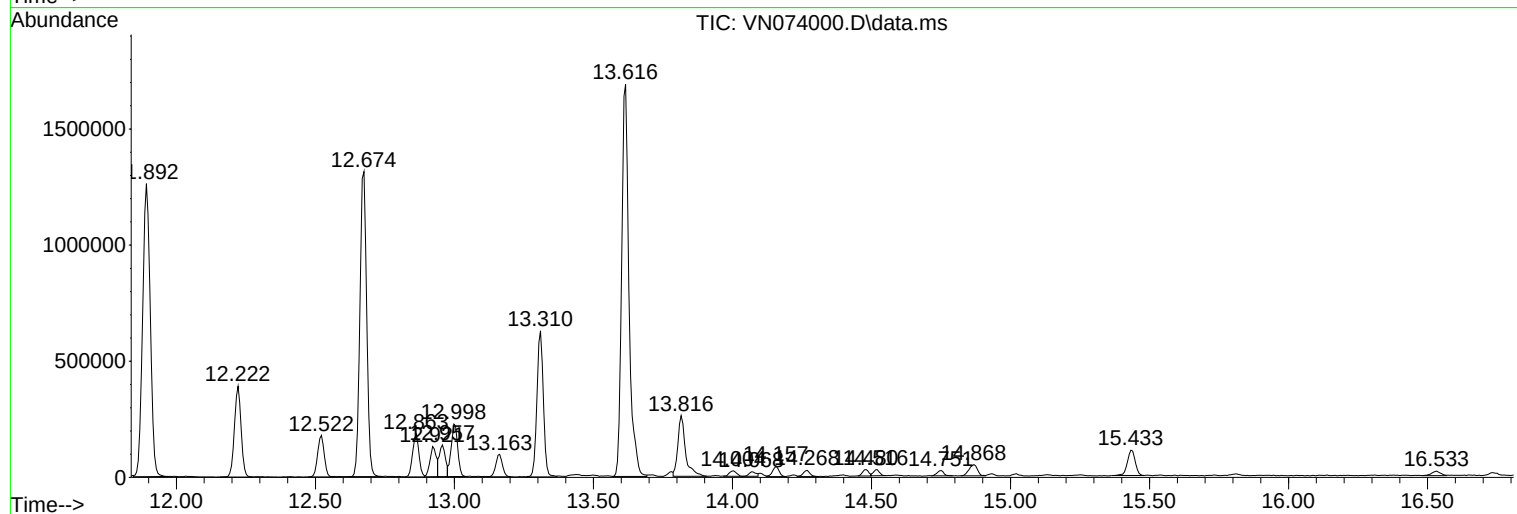
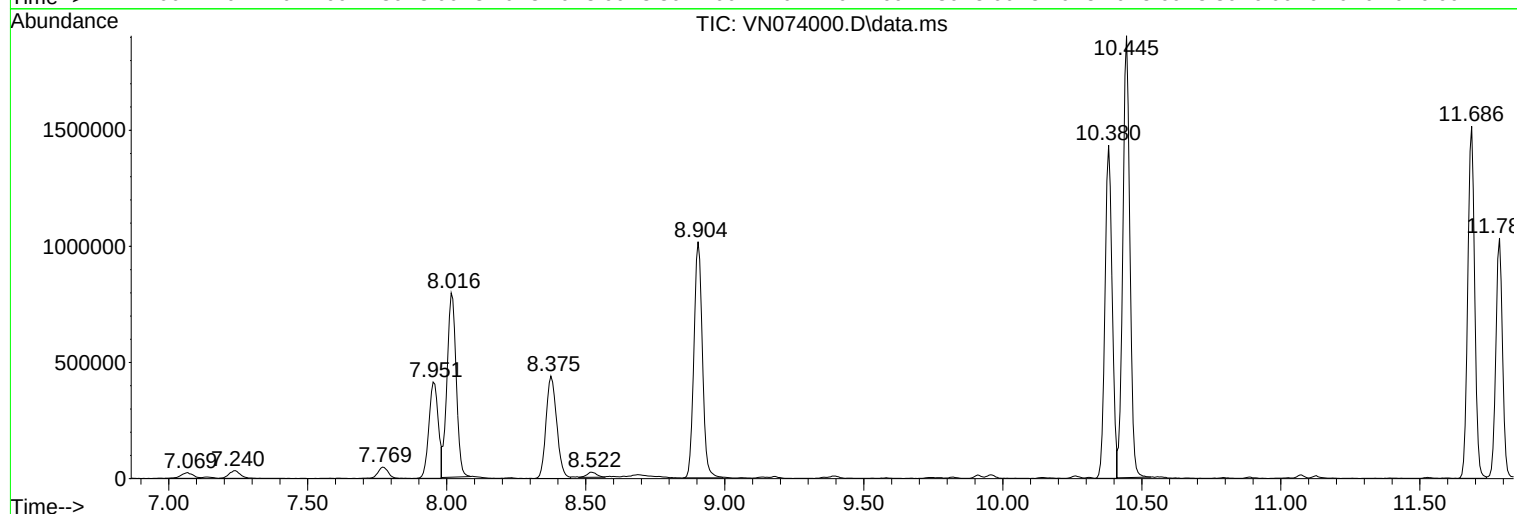
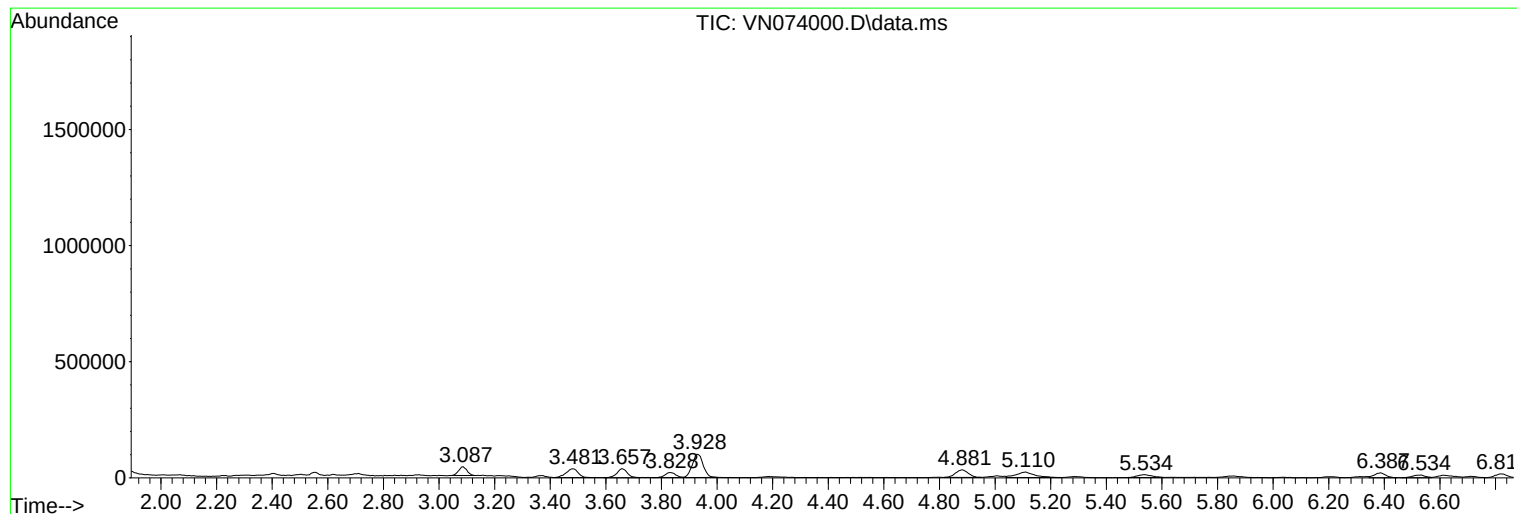
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Quant Title : SW846 8260

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



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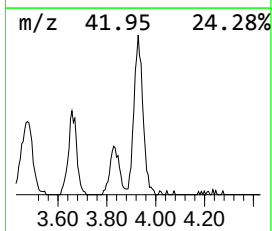
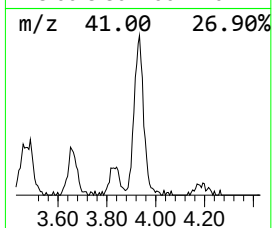
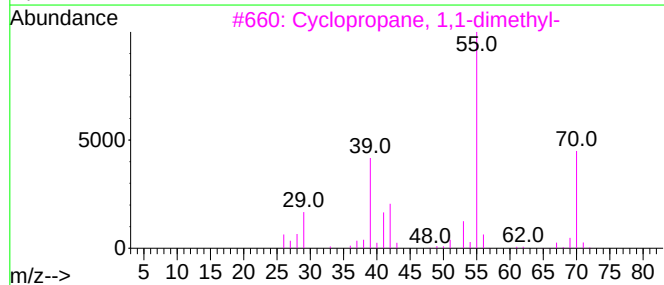
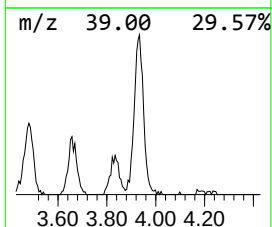
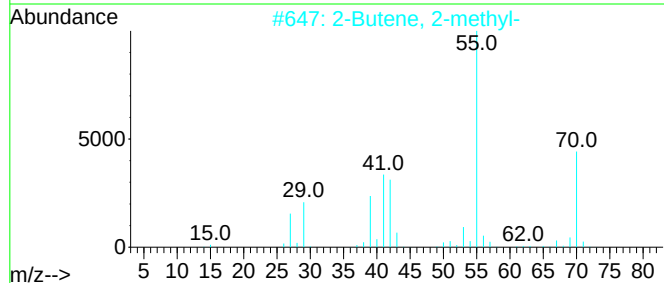
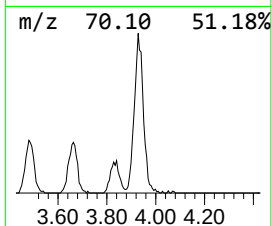
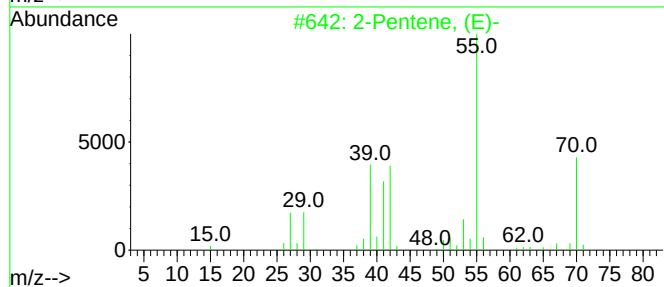
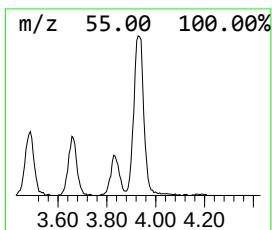
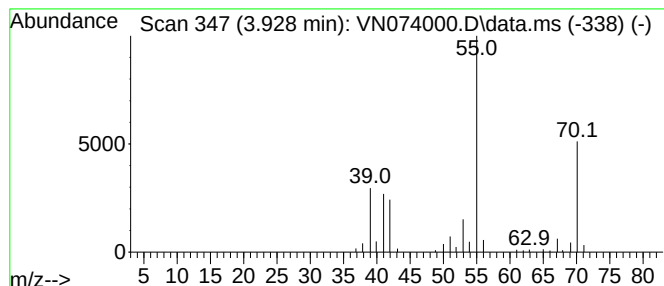
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentene, (E)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.928	7.54 ug/l	271528	Pentafluorobenzene	8.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (E)-			70	C5H10	000646-04-8	91
2	2-Butene, 2-methyl-			70	C5H10	000513-35-9	90
3	Cyclopropane, 1,1-dimethyl-			70	C5H10	001630-94-0	87
4	Cyclopropane, 1,2-dimethyl-, trans-			70	C5H10	002402-06-4	83
5	Cyclopropane, 1,2-dimethyl-, cis-			70	C5H10	000930-18-7	78



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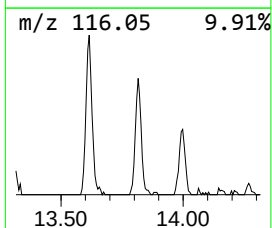
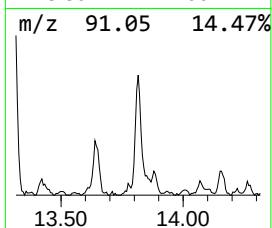
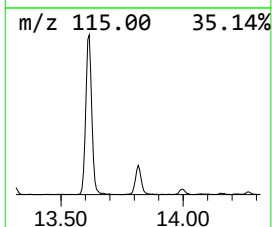
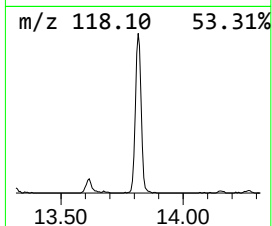
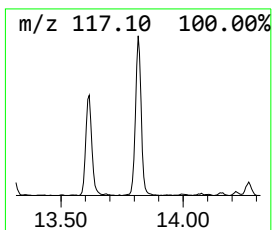
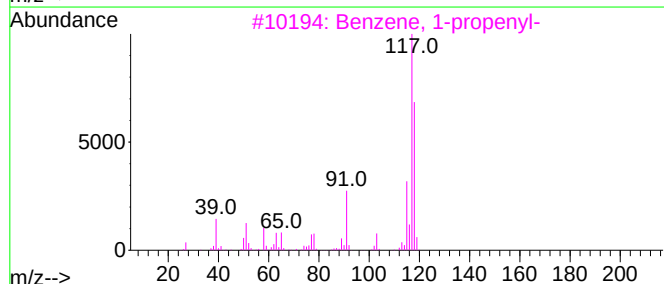
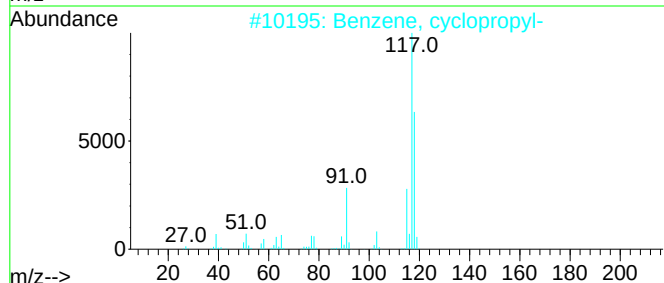
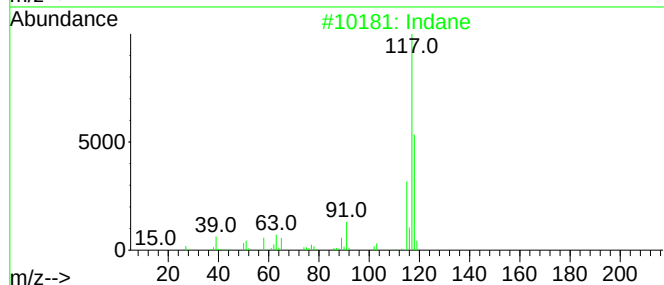
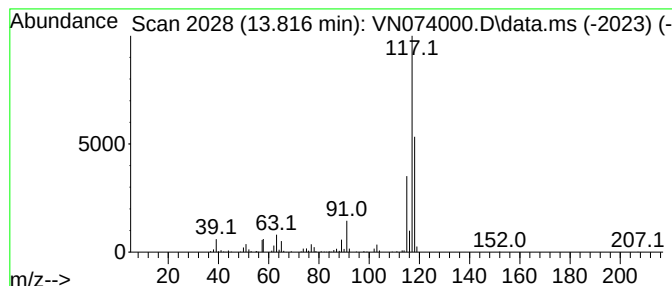
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	8.05 ug/l	496988	1,4-Dichlorobenzene-d4	13.616

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	74
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentene, (E)-	3.928	7.5	ug/l	271528	1	8.016	1800390	50.0
Indane	13.816	8.0	ug/l	496988	4	13.616	3087070	50.0