

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
 Data File : VN068900.D
 Acq On : 13 Oct 2021 16:23
 Operator : JC/MD
 Sample : M4180-01
 Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NYPD-69TH-WC-1

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\82N101221W.M
 Title : SW846 8260

Signal : TIC: VN068900.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.029	1853	1879	1898	rBV3	31481	95777	1.19%	0.109%
2	7.131	1900	1917	1940	rVB3	29240	85919	1.07%	0.098%
3	8.019	2222	2248	2257	rBV	268847	653618	8.13%	0.742%
4	8.080	2259	2271	2289	rVV	561475	1372993	17.08%	1.559%
5	8.166	2290	2303	2328	rVB3	113277	283634	3.53%	0.322%
6	8.289	2328	2349	2367	rBV3	129312	303212	3.77%	0.344%
7	8.434	2384	2403	2427	rBV	285052	651061	8.10%	0.739%
8	8.609	2433	2468	2492	rBV	571448	1622790	20.19%	1.843%
9	8.702	2494	2503	2526	rVV4	45836	114299	1.42%	0.130%
10	8.820	2529	2547	2572	rVB3	73649	169053	2.10%	0.192%
11	8.965	2580	2601	2628	rBV	753332	1658022	20.63%	1.883%
12	9.456	2753	2784	2794	rBV4	242820	638675	7.95%	0.725%
13	9.507	2795	2803	2819	rVV2	155365	305010	3.80%	0.346%
14	9.590	2820	2834	2843	rVV	58373	126160	1.57%	0.143%
15	9.730	2864	2886	2902	rVB6	58547	147542	1.84%	0.168%
16	9.880	2925	2942	2964	rVB	686207	1402588	17.45%	1.593%
17	10.014	2966	2992	3006	rBV	876108	1889024	23.51%	2.145%
18	10.078	3007	3016	3024	rBV	120982	201899	2.51%	0.229%
19	10.215	3046	3067	3095	rBV2	269390	707182	8.80%	0.803%
20	10.371	3111	3125	3137	rVB2	248835	454968	5.66%	0.517%
21	10.443	3137	3152	3179	rBV	1248073	2456301	30.56%	2.789%
22	10.623	3205	3219	3234	rVB8	213896	482684	6.01%	0.548%
23	10.784	3269	3279	3295	rBV5	57223	125092	1.56%	0.142%
24	10.872	3296	3312	3335	rBV10	172892	532179	6.62%	0.604%
25	10.963	3337	3346	3358	rVB2	106196	186479	2.32%	0.212%
26	11.055	3360	3380	3392	rBV3	105210	229090	2.85%	0.260%
27	11.154	3408	3417	3434	rVB	155967	322880	4.02%	0.367%
28	11.344	3481	3488	3500	rVB3	103273	164369	2.05%	0.187%
29	11.401	3503	3509	3519	rVB6	65805	96577	1.20%	0.110%
30	11.454	3521	3529	3537	rBV4	97135	139917	1.74%	0.159%
31	11.527	3538	3556	3582	rVB6	455105	1242860	15.46%	1.411%
32	11.628	3582	3594	3617	rBV	414036	781953	9.73%	0.888%
33	11.746	3622	3638	3654	rBV	1108694	2100358	26.13%	2.385%
34	11.846	3666	3675	3695	rVB	630300	1109801	13.81%	1.260%
35	12.184	3793	3801	3814	rVB5	118944	208391	2.59%	0.237%
36	12.256	3816	3828	3840	rBV3	287345	573176	7.13%	0.651%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
 Data File : VN068900.D
 Acq On : 13 Oct 2021 16:23
 Operator : JC/MD
 Sample : M4180-01
 Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NYPD-69TH-WC-1

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\82N101221W.M
 Title : SW846 8260

37	12.369	3863	3870	3880	rVB2	252854	385544	4.80%	0.438%
38	12.583	3941	3950	3959	rBV	295636	533270	6.64%	0.606%
39	12.876	4051	4059	4064	rBV2	167183	250388	3.12%	0.284%
40	12.921	4066	4076	4089	rVB	1424329	2351826	29.26%	2.670%
41	13.050	4117	4124	4143	rVB3	449544	812247	10.11%	0.922%
42	13.417	4254	4261	4272	rVB3	267301	438950	5.46%	0.498%
43	13.500	4277	4292	4297	rBV3	369789	860085	10.70%	0.977%
44	13.608	4325	4332	4338	rBV2	253614	328785	4.09%	0.373%
45	13.677	4349	4358	4367	rBV	1156356	1759558	21.89%	1.998%
46	13.728	4369	4377	4401	rVB4	846491	2214986	27.56%	2.515%
47	13.841	4407	4419	4426	rBV	1777936	2830286	35.22%	3.214%
48	13.879	4427	4433	4442	rVV	1199115	2063916	25.68%	2.344%
49	13.922	4443	4449	4468	rVB3	1017607	2445920	30.43%	2.777%
50	13.999	4469	4478	4490	rBV3	589318	949964	11.82%	1.079%
51	14.139	4519	4530	4547	rVB3	544768	1487368	18.51%	1.689%
52	14.219	4547	4560	4573	rBV	5057151	8036630	100.00%	9.126%
53	14.329	4590	4601	4614	rVB3	1689669	2956066	36.78%	3.357%
54	14.396	4617	4626	4639	rVB6	232517	447636	5.57%	0.508%
55	14.466	4645	4652	4662	rBV6	201907	299705	3.73%	0.340%
56	14.541	4664	4680	4690	rBV	2846157	4840315	60.23%	5.496%
57	14.622	4702	4710	4720	rVB	1046472	1387709	17.27%	1.576%
58	14.812	4771	4781	4791	rBV2	1691201	2935339	36.52%	3.333%
59	14.930	4808	4825	4837	rBV3	2914251	6940330	86.36%	7.881%
60	14.984	4837	4845	4862	rVB	989962	1820571	22.65%	2.067%
61	15.198	4901	4925	4935	rBV4	2016219	4714138	58.66%	5.353%
62	15.311	4955	4967	4984	rVB4	1381495	3151680	39.22%	3.579%
63	15.472	5015	5027	5031	rBV5	304507	530973	6.61%	0.603%
64	15.617	5069	5081	5091	rBV4	613119	1113847	13.86%	1.265%
65	15.670	5093	5101	5136	rVB5	552247	1596955	19.87%	1.813%
66	15.807	5137	5152	5168	rBV	831780	1745627	21.72%	1.982%
67	16.110	5256	5265	5277	rVB	288548	492616	6.13%	0.559%
68	16.338	5340	5350	5362	rBV8	104164	200392	2.49%	0.228%
69	16.620	5441	5455	5498	rVB	603473	1508058	18.76%	1.712%

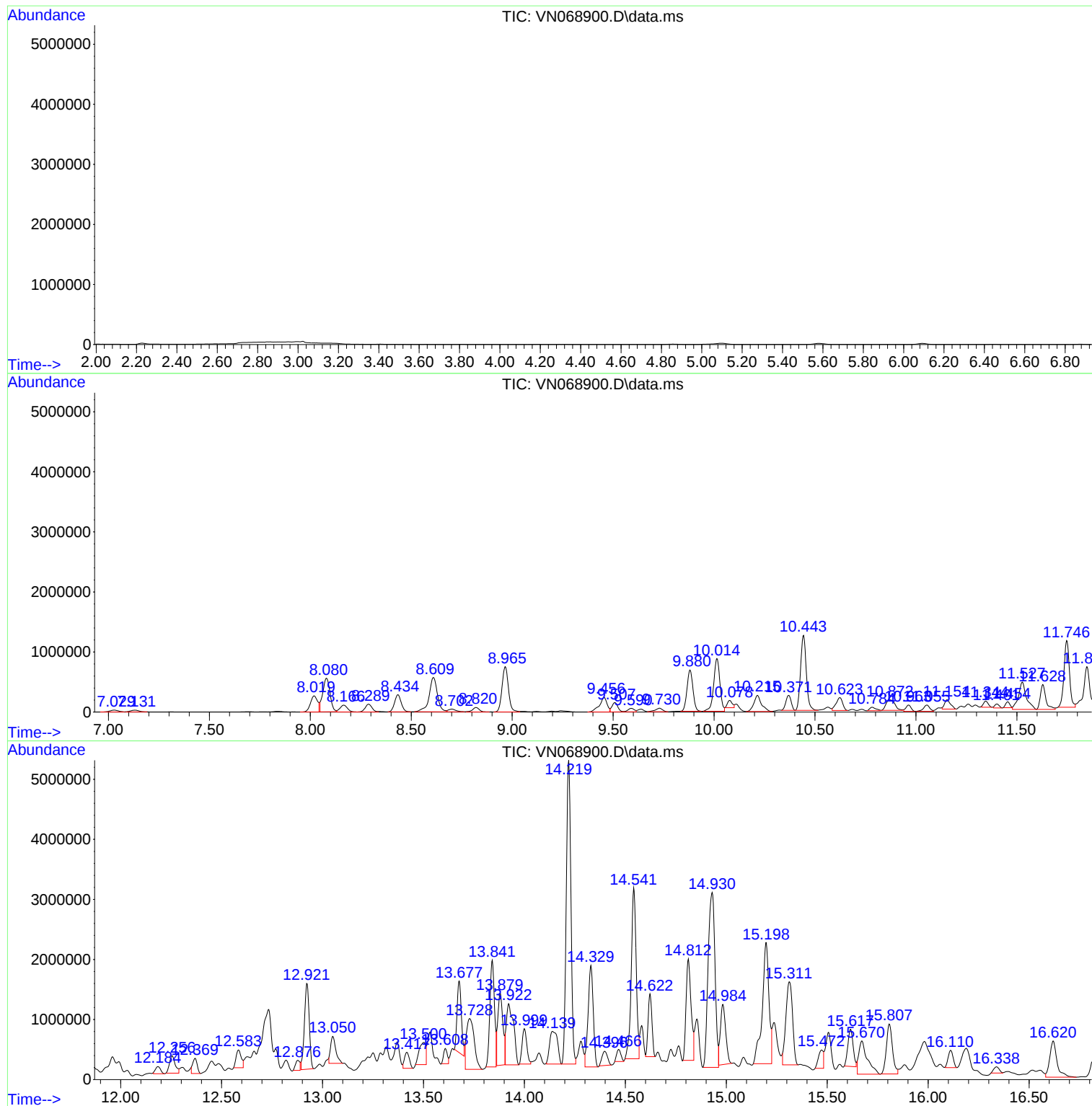
Sum of corrected areas: 88067213

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

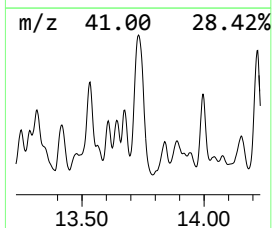
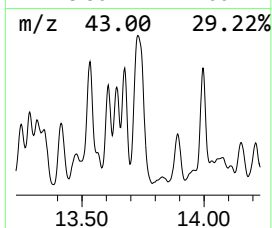
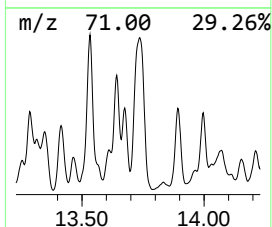
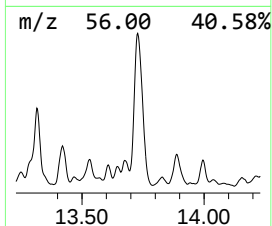
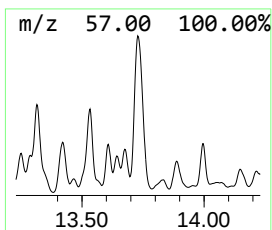
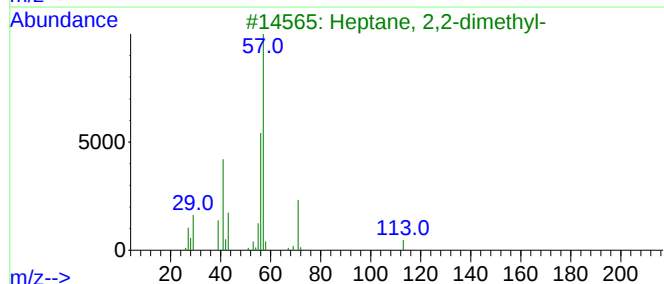
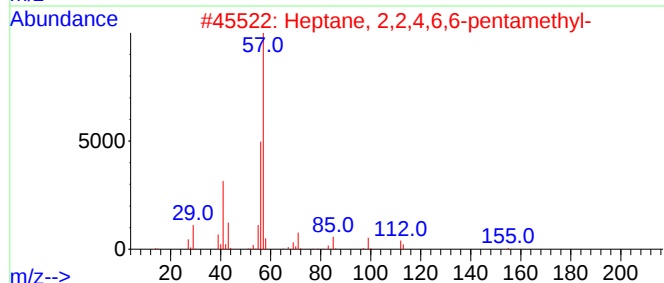
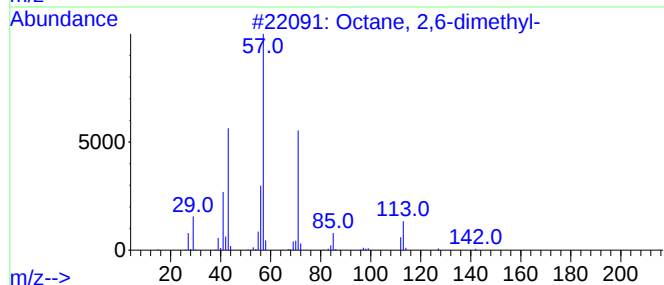
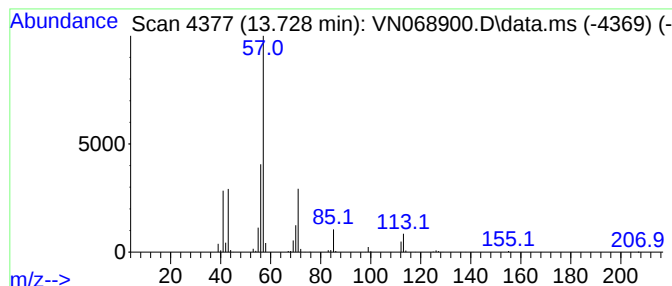
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.729	62.94 ug/l	2214990	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
2			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	59
3			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	59
4			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	53
5			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

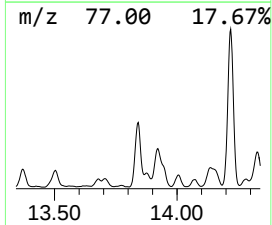
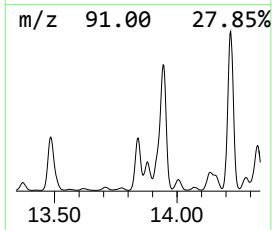
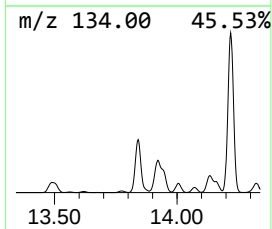
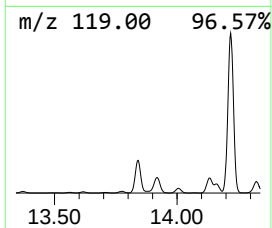
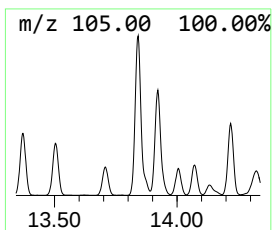
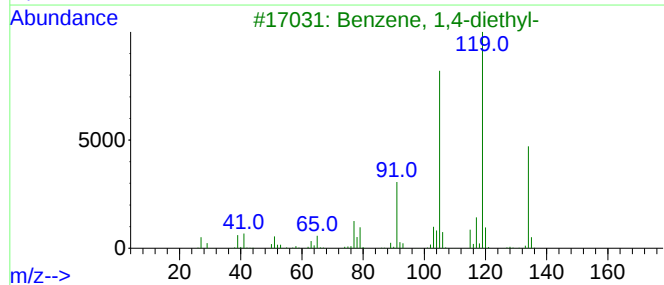
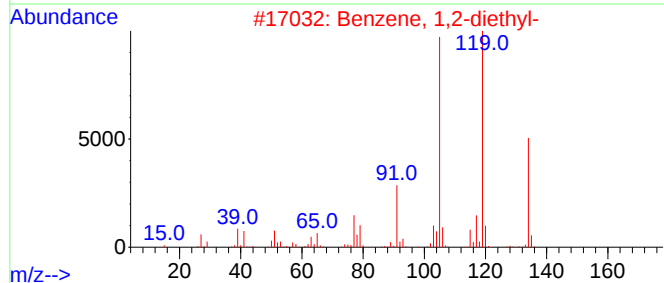
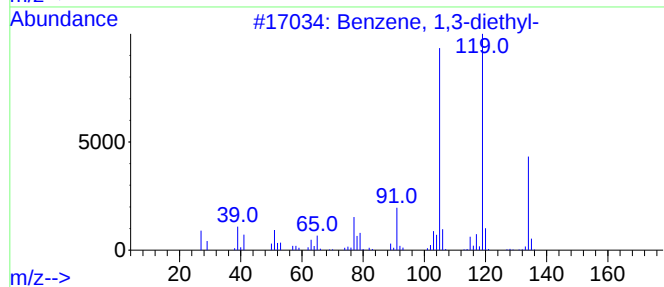
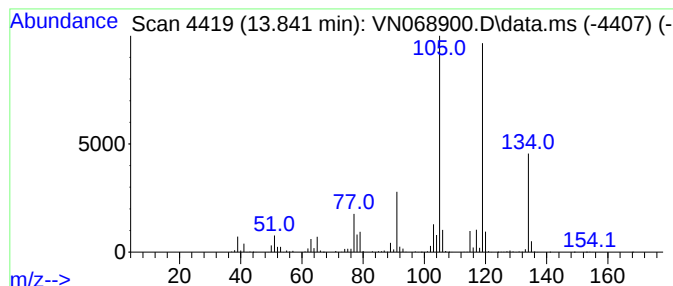
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.841	80.43 ug/l	2830290	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

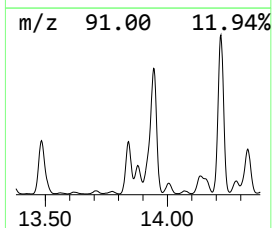
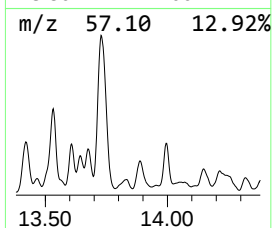
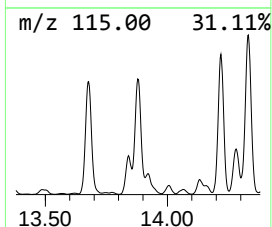
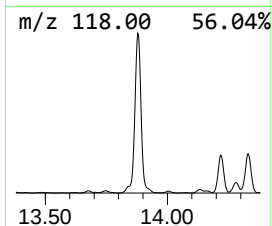
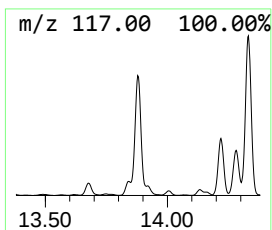
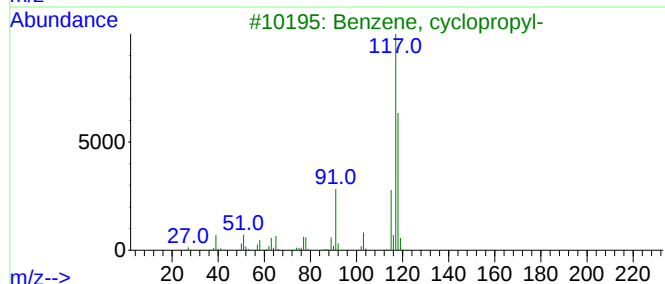
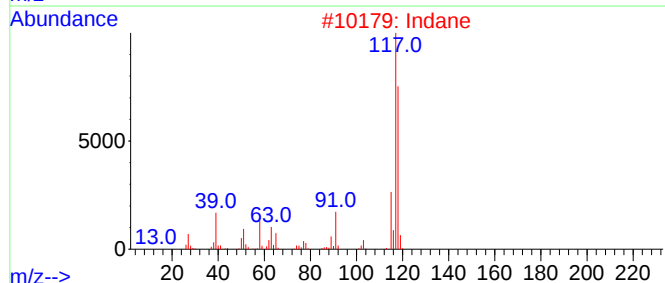
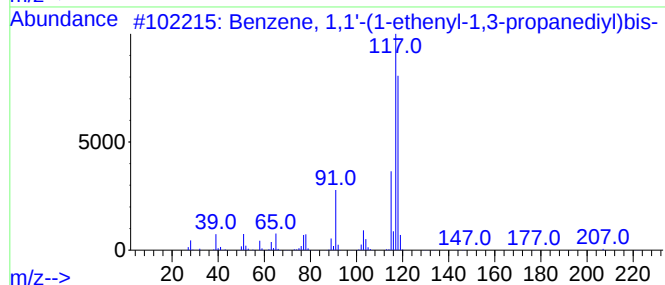
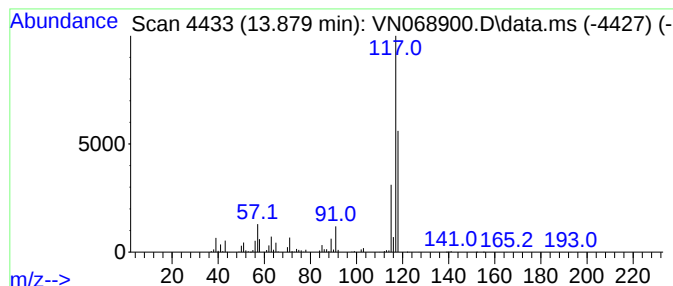
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,1'-(1-ethenyl-1,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.879	58.65 ug/l	2063920	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	78
2			Indane	118	C9H10	000496-11-7	74
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
5			Indole	117	C8H7N	000120-72-9	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

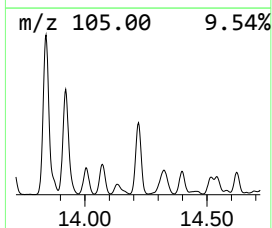
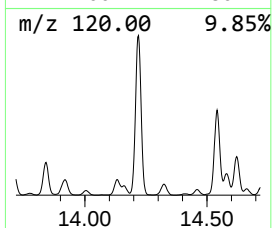
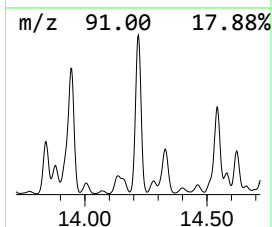
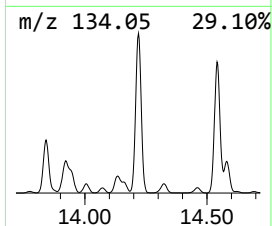
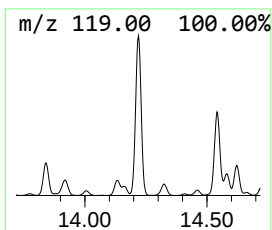
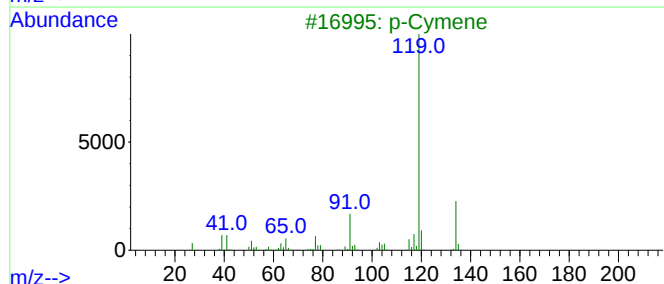
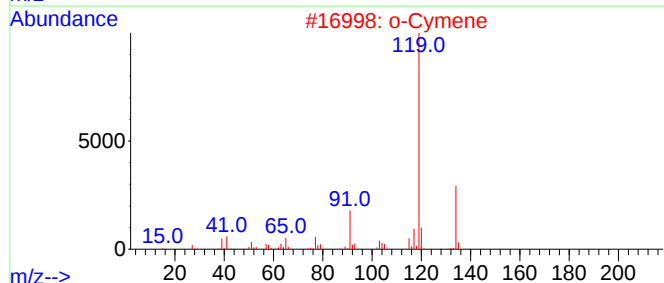
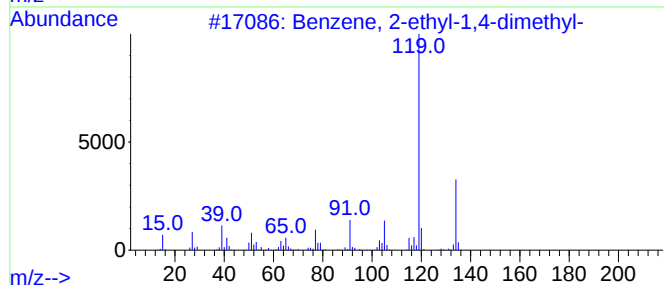
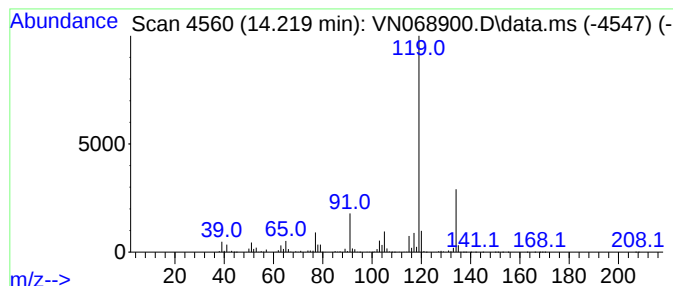
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.219	228.37 ug/l	8036630	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

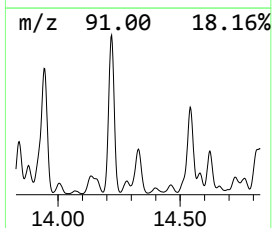
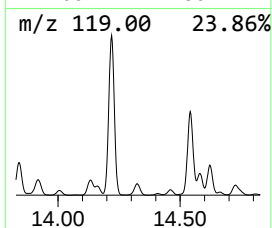
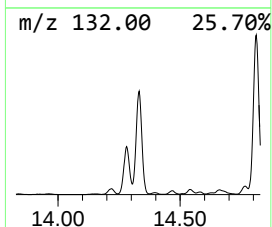
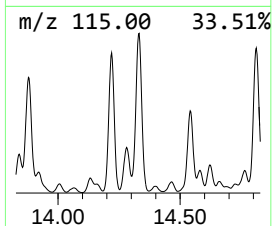
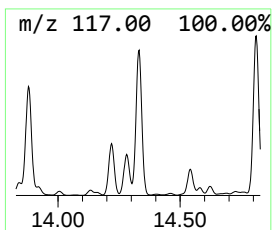
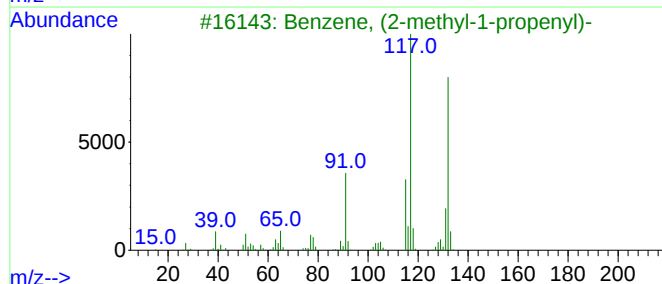
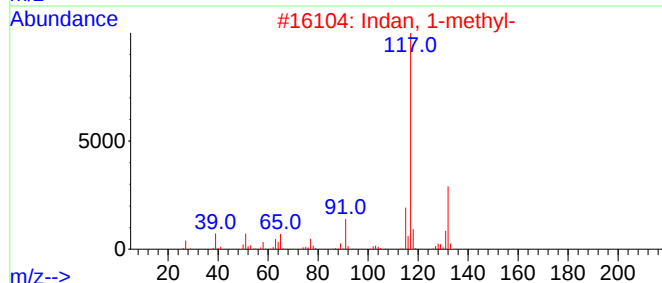
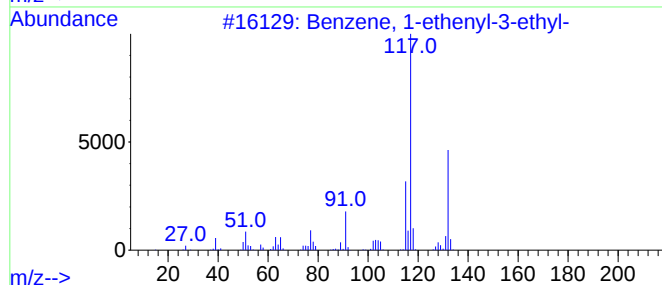
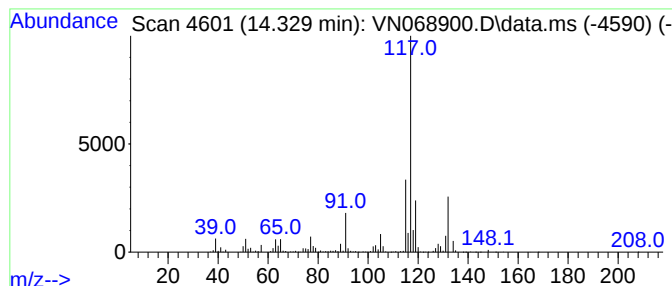
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.329	84.00 ug/l	2956070	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
2			Indan, 1-methyl-	132	C10H12	000767-58-8	81
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	76
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

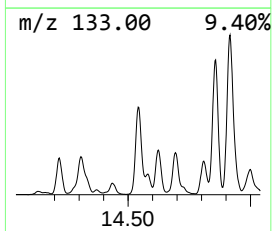
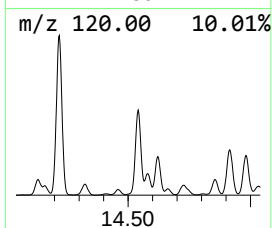
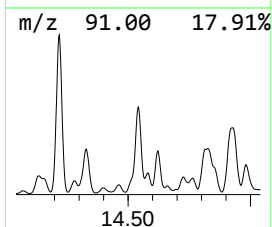
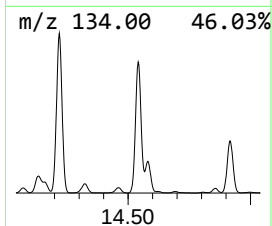
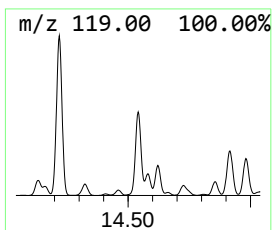
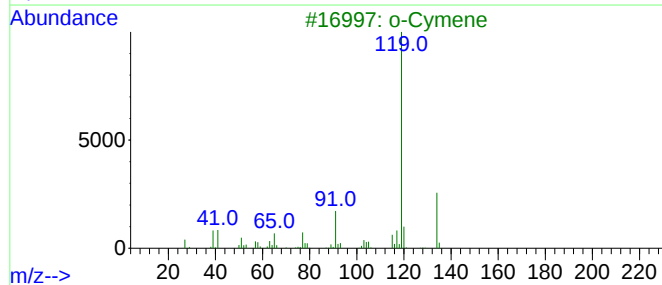
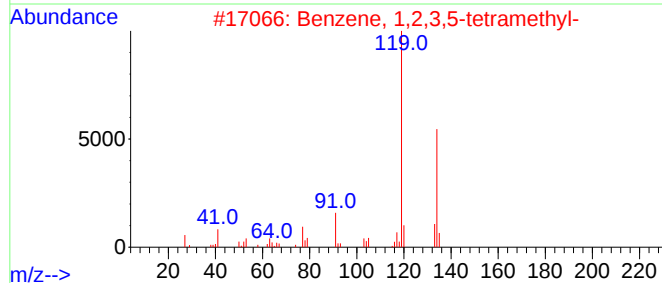
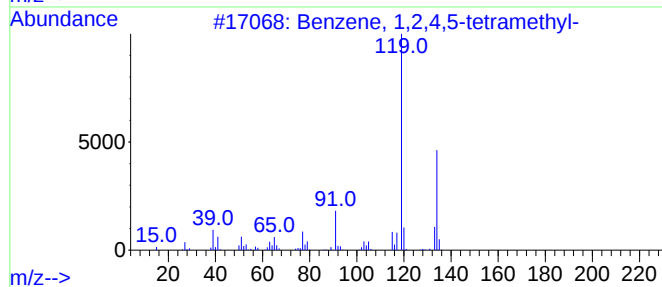
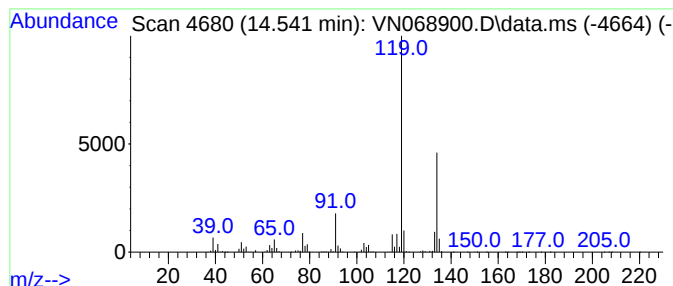
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.541	137.54 ug/l	4840320	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

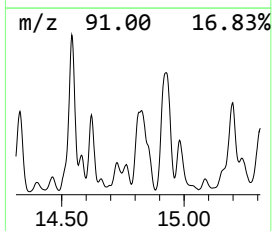
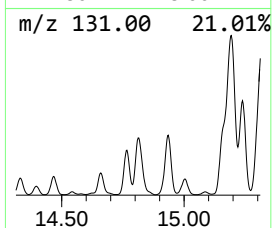
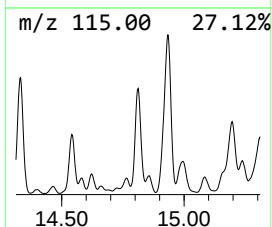
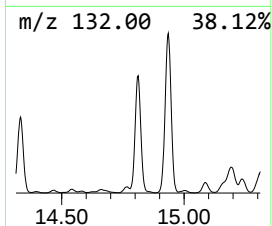
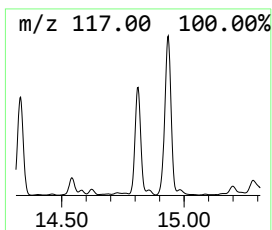
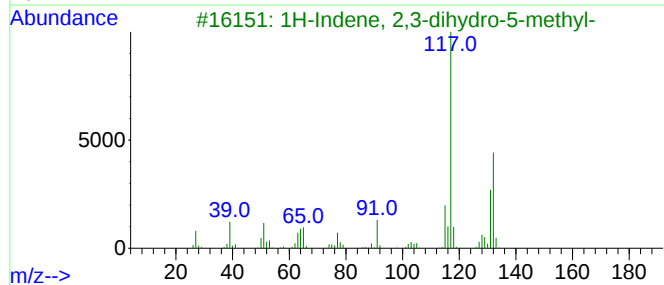
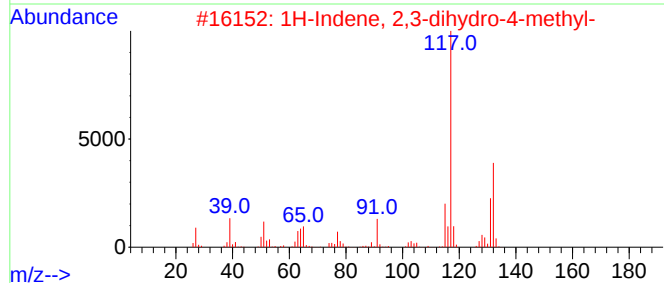
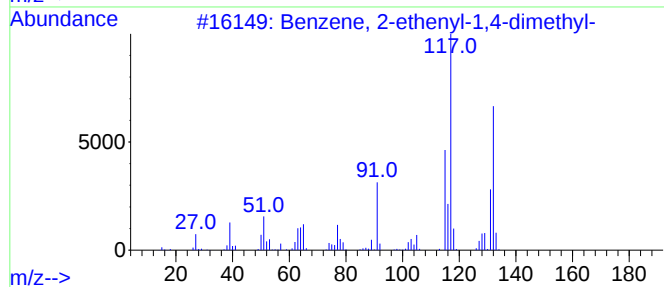
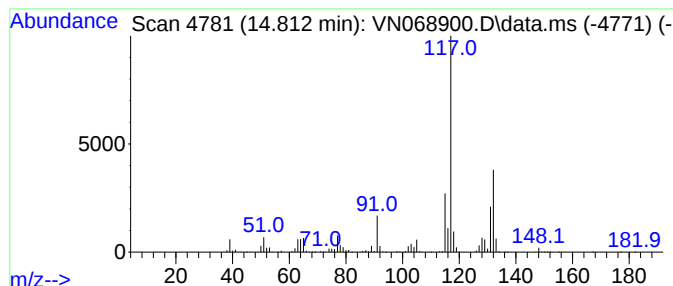
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.812	83.41 ug/l	2935340	1,4-Dichlorobenzene-d4	13.678

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-4-methyl-	132	C10H12	000824-22-6	94
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

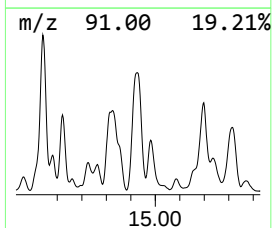
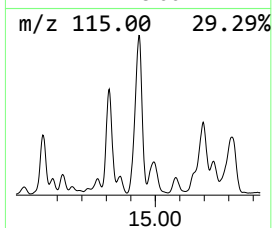
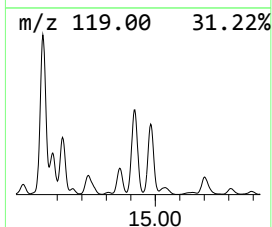
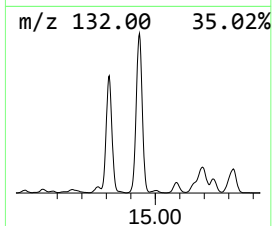
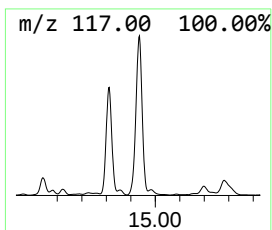
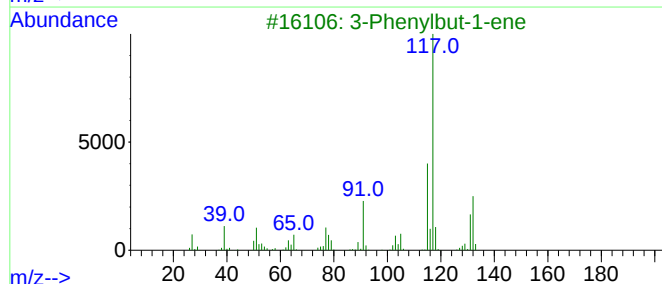
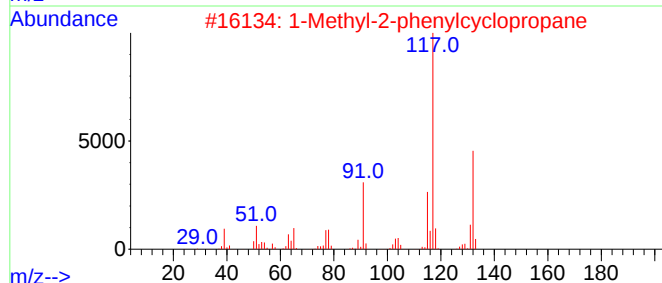
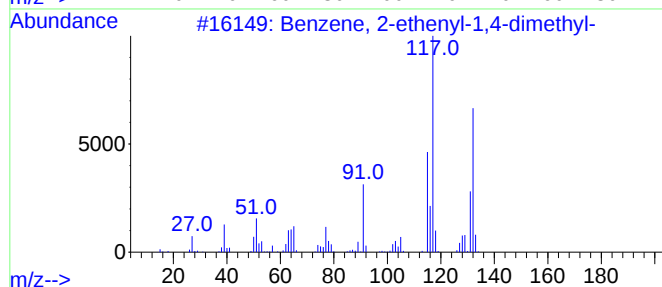
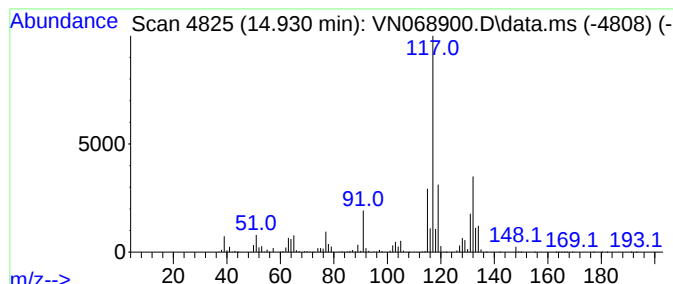
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 8 1-Methyl-2-phenylcyclopropane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.930	197.22 ug/l	6940330	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

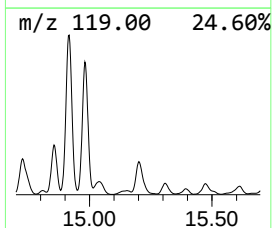
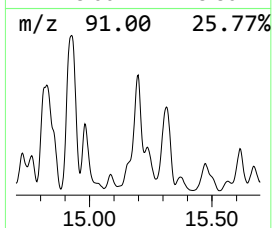
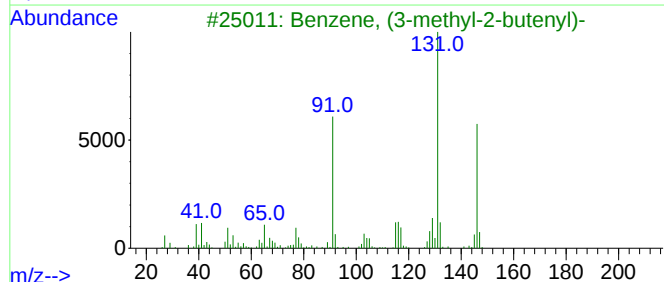
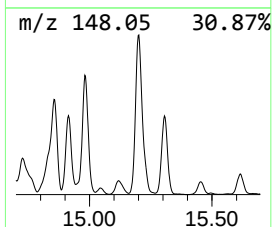
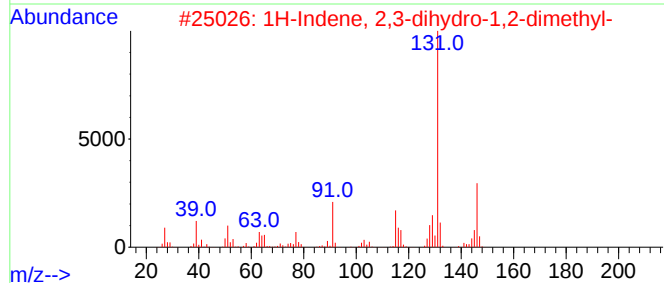
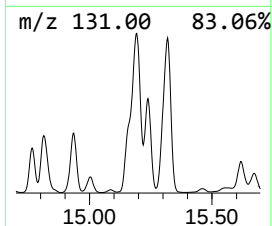
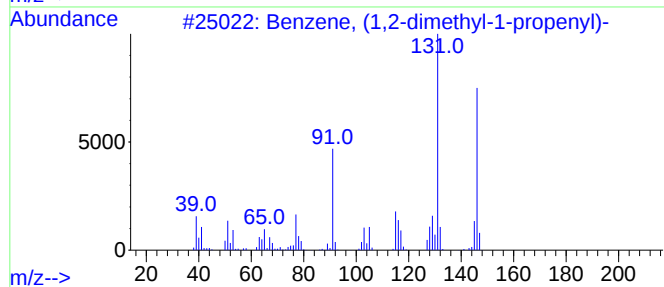
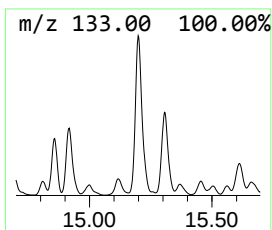
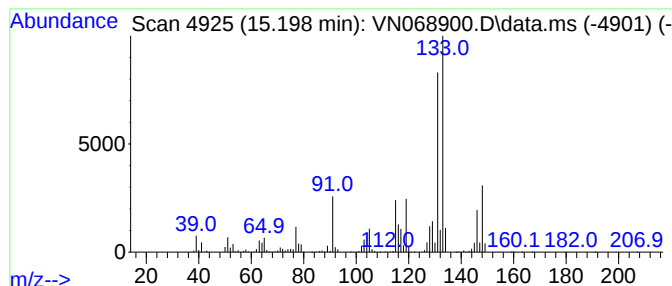
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.198	133.96 ug/l	4714140	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	50
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

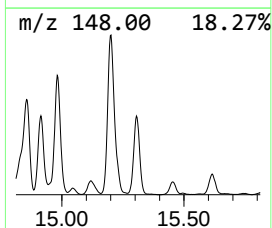
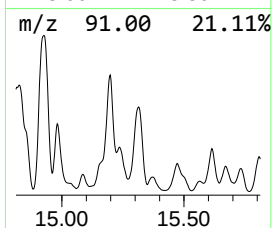
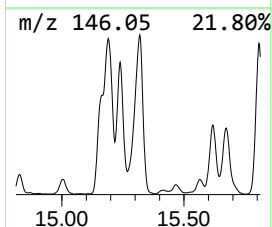
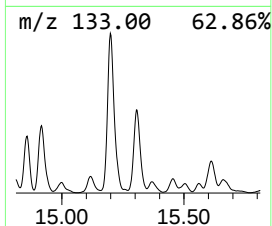
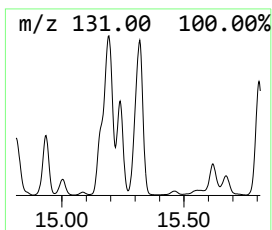
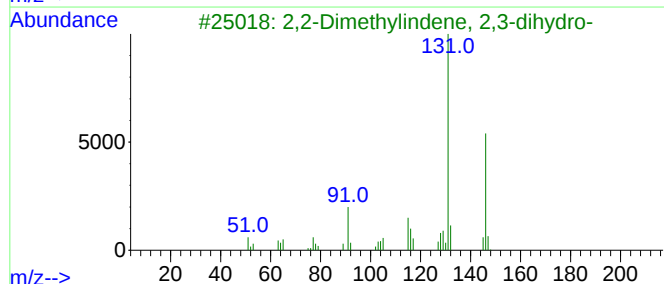
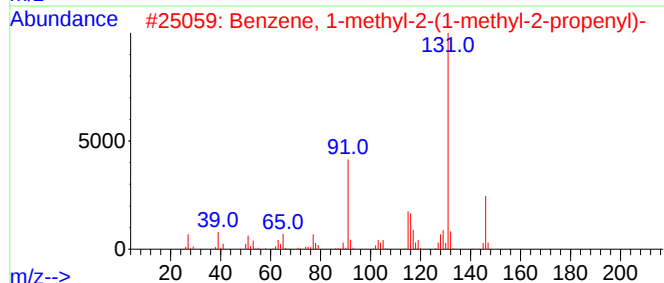
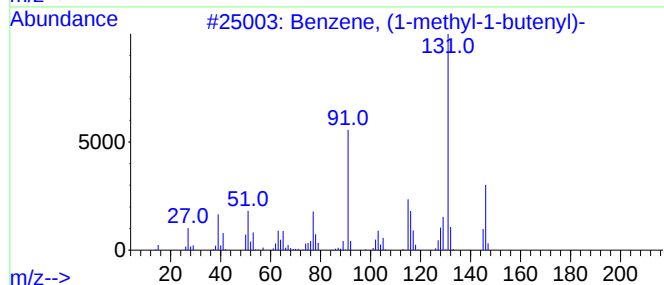
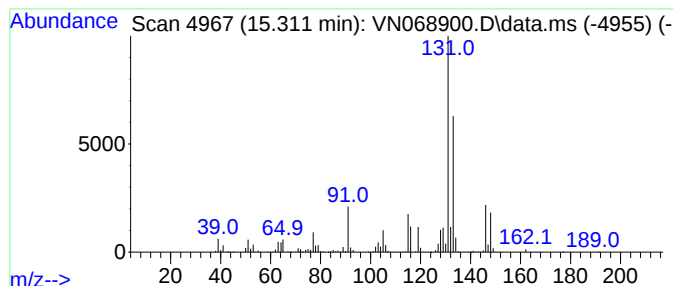
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.311	89.56 ug/l	3151680	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	83
2			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	70
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	70
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101321\
Data File : VN068900.D
Acq On : 13 Oct 2021 16:23
Operator : JC/MD
Sample : M4180-01
Misc : 9.31g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
NYPD-69TH-WC-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.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
Octane, 2,6-dim...	13.729	62.9	ug/l	2214990	4	13.678	1759560	50.0
Benzene, 1,3-di...	13.841	80.4	ug/l	2830290	4	13.678	1759560	50.0
Benzene, 1,1'-(...	13.879	58.6	ug/l	2063920	4	13.678	1759560	50.0
Benzene, 2-ethy...	14.219	228.4	ug/l	8036630	4	13.678	1759560	50.0
Benzene, 1-ethe...	14.329	84.0	ug/l	2956070	4	13.678	1759560	50.0
Benzene, 1,2,4,...	14.541	137.5	ug/l	4840320	4	13.678	1759560	50.0
Benzene, 2-ethe...	14.812	83.4	ug/l	2935340	4	13.678	1759560	50.0
1-Methyl-2-phen...	14.930	197.2	ug/l	6940330	4	13.678	1759560	50.0
Benzene, (1,2-d...	15.198	134.0	ug/l	4714140	4	13.678	1759560	50.0
Benzene, (1-met...	15.311	89.6	ug/l	3151680	4	13.678	1759560	50.0