

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092120\
 Data File : VN063487.D
 Acq On : 21 Sep 2020 15:45
 Operator : JC/MD
 Sample : L4056-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 229

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.293	161	172	186	rBV	41460	66050	3.73%	0.366%
2	2.422	198	212	224	rBV5	20397	46113	2.61%	0.256%
3	3.232	445	464	498	rBV	361105	893519	50.50%	4.955%
4	3.772	611	632	667	rBV2	206095	560240	31.66%	3.107%
5	4.016	693	708	727	rBV	17474	48253	2.73%	0.268%
6	4.727	910	929	941	rBV4	9528	32939	1.86%	0.183%
7	4.994	986	1012	1031	rBV3	16779	60544	3.42%	0.336%
8	6.502	1465	1481	1487	rBV6	7716	18304	1.03%	0.102%
9	7.550	1790	1807	1818	rBV2	213380	534141	30.19%	2.962%
10	7.627	1818	1831	1855	rVB	368864	885364	50.03%	4.910%
11	7.987	1921	1943	1965	rBV	257278	638188	36.07%	3.539%
12	8.090	1965	1975	1990	rVB6	11300	26762	1.51%	0.148%
13	8.405	2063	2073	2089	rVB3	12385	29635	1.67%	0.164%
14	8.550	2101	2118	2139	rBV	600304	1256029	70.98%	6.966%
15	8.756	2170	2182	2195	rBV4	13998	31646	1.79%	0.175%
16	9.007	2248	2260	2279	rBV	23536	48592	2.75%	0.269%
17	9.952	2545	2554	2566	rBV3	13741	26532	1.50%	0.147%
18	10.061	2572	2588	2600	rBV	960104	1769513	100.00%	9.813%
19	10.125	2600	2608	2618	rVB2	70218	120318	6.80%	0.667%
20	10.691	2772	2784	2792	rBV4	13332	26014	1.47%	0.144%
21	11.379	2984	2998	3017	rBV	963175	1669749	94.36%	9.260%
22	11.595	3050	3065	3075	rBV4	26314	57588	3.25%	0.319%
23	11.920	3148	3166	3178	rBV	90373	159614	9.02%	0.885%
24	12.006	3183	3193	3202	rBV3	15869	26849	1.52%	0.149%
25	12.376	3295	3308	3329	rVB	780630	1313732	74.24%	7.286%
26	12.633	3380	3388	3392	rBV2	21476	34433	1.95%	0.191%
27	12.662	3393	3397	3404	rVV	29109	42025	2.37%	0.233%
28	12.707	3404	3411	3433	rVB	56064	98641	5.57%	0.547%
29	12.865	3451	3460	3470	rVB2	37640	62307	3.52%	0.346%
30	13.013	3497	3506	3515	rVB2	24373	37293	2.11%	0.207%
31	13.318	3588	3601	3625	rBV2	948470	1736841	98.15%	9.632%
32	13.424	3626	3634	3645	rVV3	36878	65020	3.67%	0.361%
33	13.489	3648	3654	3656	rVV6	17689	22562	1.28%	0.125%
34	13.518	3657	3663	3669	rVV3	50690	80340	4.54%	0.446%

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35	13.556	3669	3675	3694	rVB2	61043	113509	6.41%	0.629%
36	13.649	3694	3704	3710	rBV9	17801	28530	1.61%	0.158%
37	13.714	3715	3724	3732	rBV2	20622	33762	1.91%	0.187%
38	13.778	3737	3744	3746	rBV2	20256	23191	1.31%	0.129%
39	13.865	3762	3771	3780	rVB	88390	134466	7.60%	0.746%
40	13.920	3781	3788	3794	rBV5	17726	25052	1.42%	0.139%
41	13.968	3795	3803	3813	rVV3	52153	83280	4.71%	0.462%
42	14.103	3836	3845	3852	rBV3	38881	61797	3.49%	0.343%
43	14.183	3861	3870	3876	rBV	77785	124247	7.02%	0.689%
44	14.222	3876	3882	3891	rVB	103078	155973	8.81%	0.865%
45	14.270	3891	3897	3903	rBV3	25849	32355	1.83%	0.179%
46	14.309	3903	3909	3922	rVB4	15790	28149	1.59%	0.156%
47	14.405	3924	3939	3945	rBV7	25087	50988	2.88%	0.283%
48	14.450	3945	3953	3961	rBV	66987	101985	5.76%	0.566%
49	14.498	3961	3968	3976	rVB4	59283	87354	4.94%	0.484%
50	14.566	3976	3989	3999	rBV3	234381	507073	28.66%	2.812%
51	14.627	4000	4008	4025	rVB4	106585	197689	11.17%	1.096%
52	14.714	4025	4035	4044	rBV2	114725	184750	10.44%	1.025%
53	14.784	4046	4057	4060	rBV7	22284	42376	2.39%	0.235%
54	14.823	4061	4069	4075	rBV4	104814	173374	9.80%	0.961%
55	14.929	4093	4102	4112	rVB3	94613	171809	9.71%	0.953%
56	15.103	4137	4156	4165	rBV3	79591	190512	10.77%	1.057%
57	15.161	4165	4174	4182	rVV	97272	172326	9.74%	0.956%
58	15.222	4182	4193	4211	rVV7	96350	352790	19.94%	1.956%
59	15.299	4212	4217	4232	rVB7	52340	91600	5.18%	0.508%
60	15.383	4232	4243	4257	rBV3	97282	194863	11.01%	1.081%
61	15.543	4277	4293	4296	rBV6	95287	200473	11.33%	1.112%
62	15.579	4297	4304	4322	rVV2	241462	486664	27.50%	2.699%
63	15.665	4322	4331	4340	rVV5	50576	97414	5.51%	0.540%
64	15.736	4343	4353	4373	rVB5	87299	240074	13.57%	1.331%
65	15.878	4384	4397	4410	rBV3	201836	419238	23.69%	2.325%
66	16.064	4445	4455	4465	rVB4	102676	196516	11.11%	1.090%
67	16.122	4465	4473	4485	rBV3	80599	151517	8.56%	0.840%
68	16.308	4520	4531	4544	rBV8	76923	211581	11.96%	1.173%
69	16.447	4565	4574	4590	rVB6	56751	138963	7.85%	0.771%

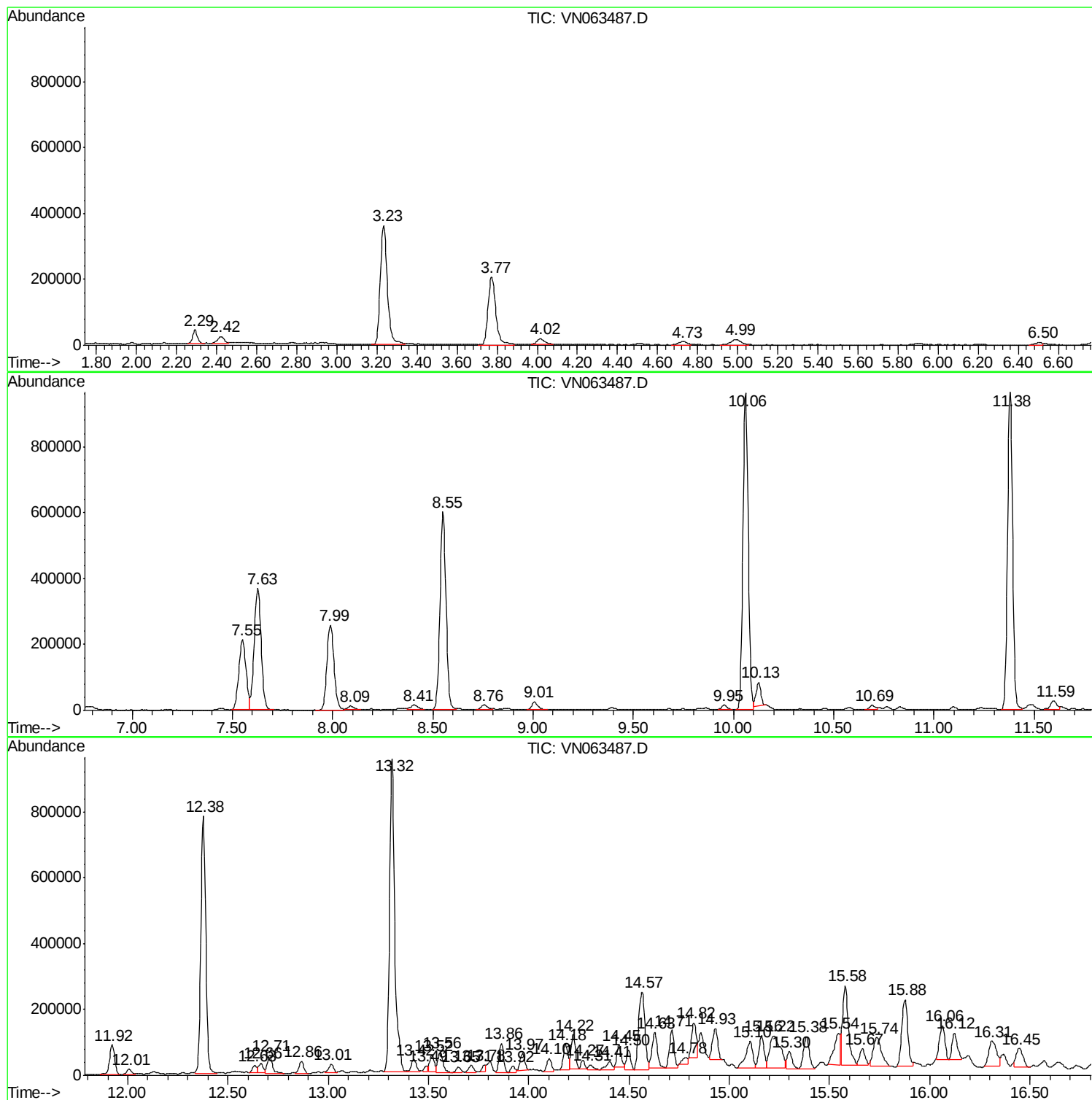
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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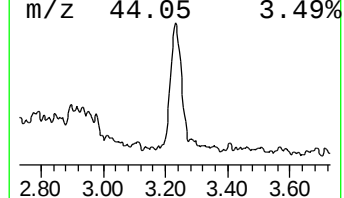
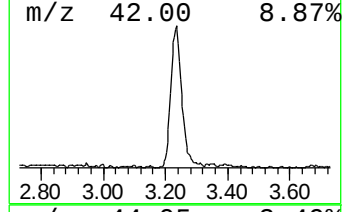
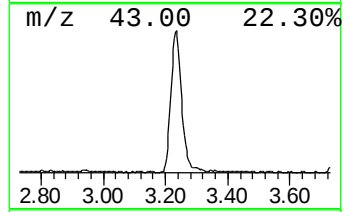
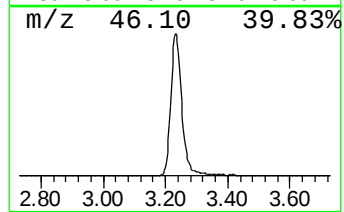
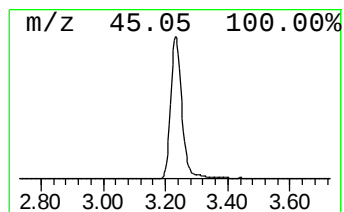
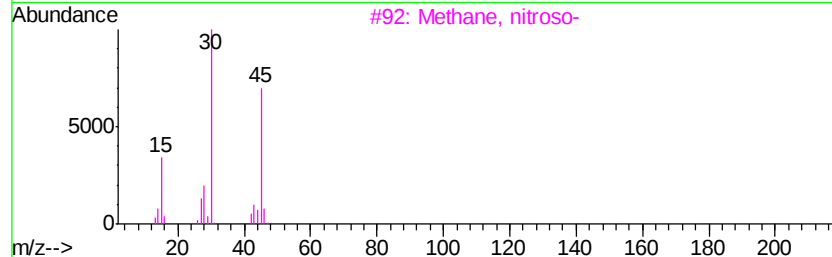
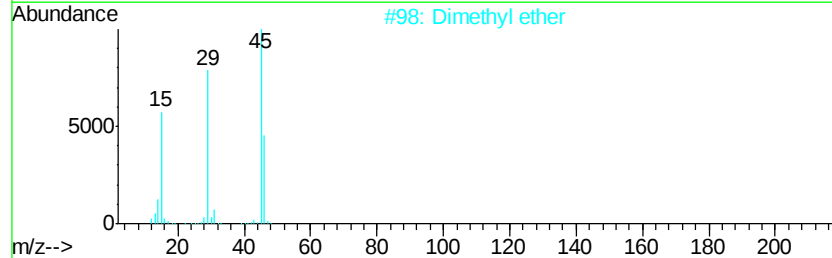
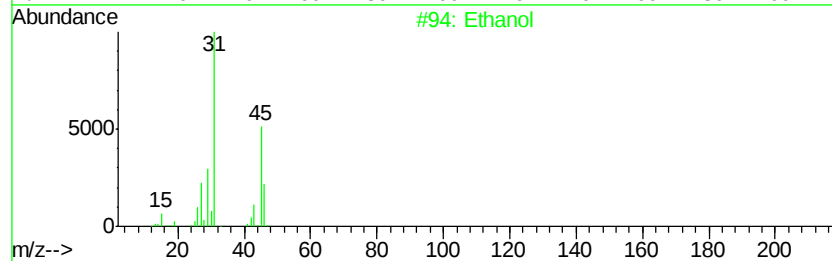
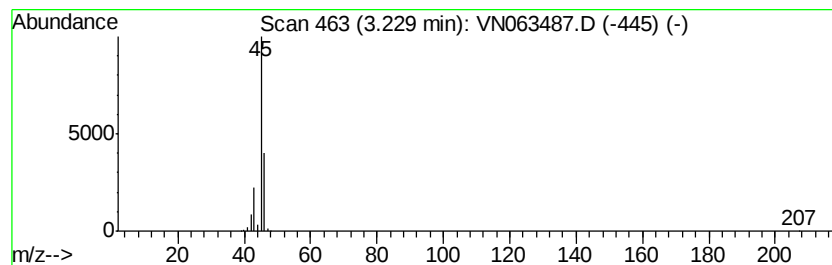
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.23	50.46 ug/l	893519	Pentafluorobenzene	7.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol		46	C2H6O	000064-17-5	74
2	Dimethyl ether		46	C2H6O	000115-10-6	9
3	Methane, nitroso-		45	CH3NO	000865-40-7	4
4	Oxalic acid		90	C2H2O4	000144-62-7	4
5	Formic acid		46	CH2O2	000064-18-6	4



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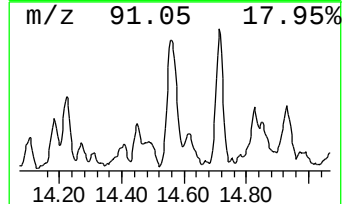
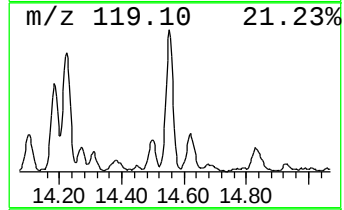
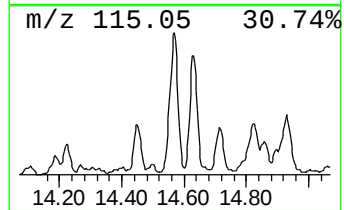
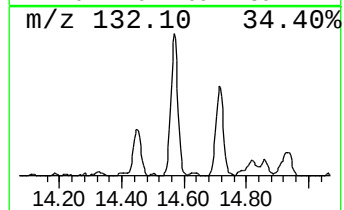
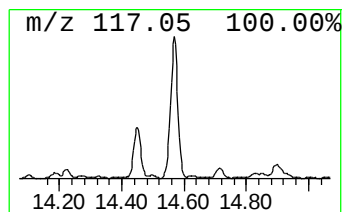
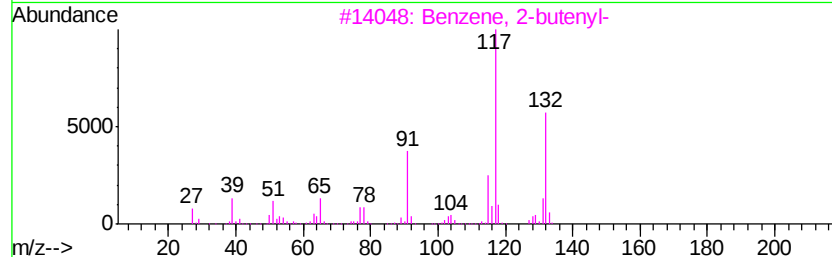
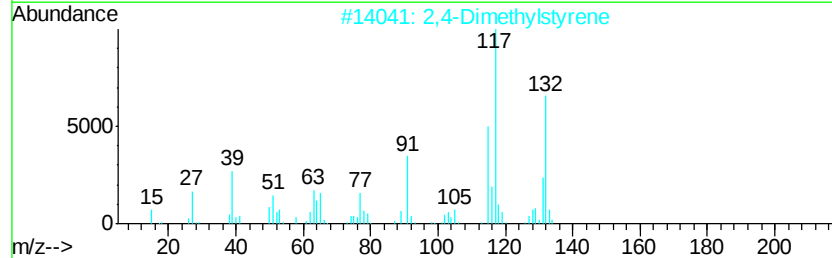
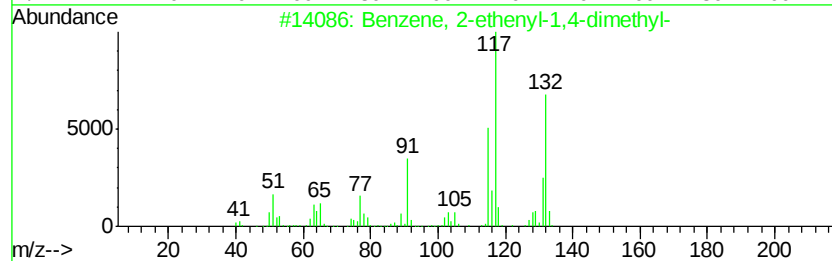
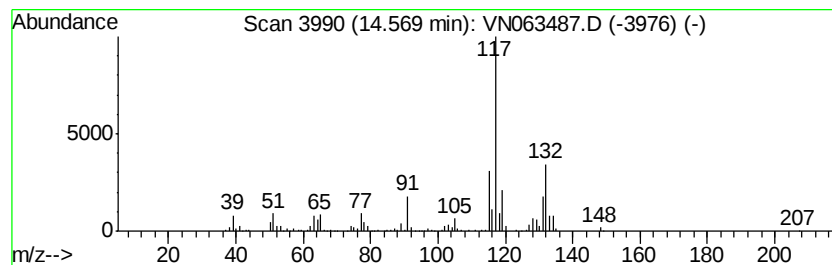
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	14.60 ug/l	507073	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	92
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	81
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81



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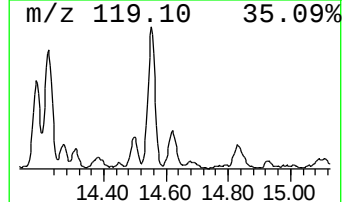
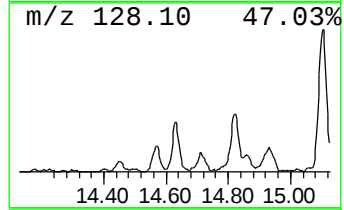
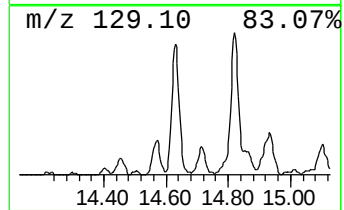
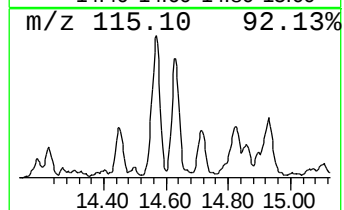
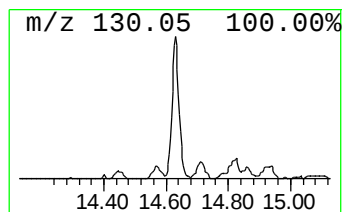
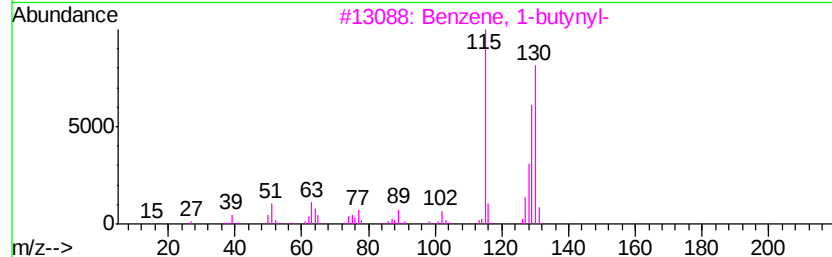
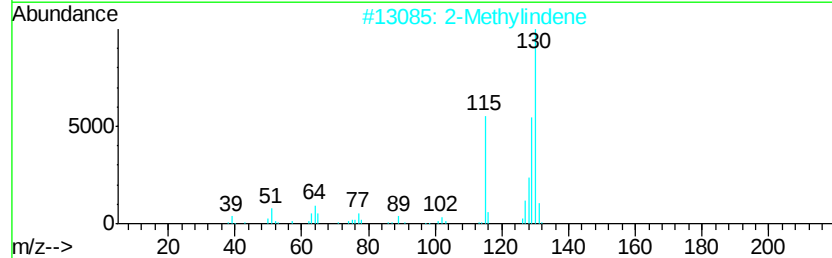
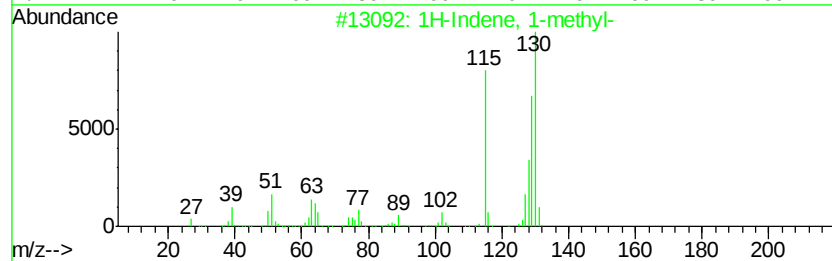
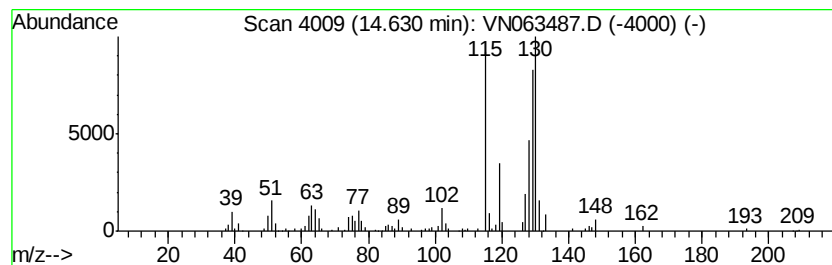
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Peak Number 3 1H-Indene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	5.69 ug/l	197689	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 1-methyl-	130 C10H10	000767-59-9	96
2	2-Methylindene	130 C10H10	002177-47-1	95
3	Benzene, 1-butynyl-	130 C10H10	000622-76-4	94
4	1,4-Dihydronaphthalene	130 C10H10	000612-17-9	92
5	Cycloprop[a]indene, 1,1a,6,6a-te...	130 C10H10	015677-15-3	91



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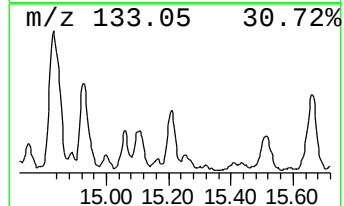
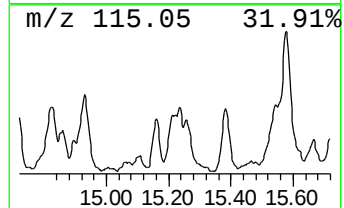
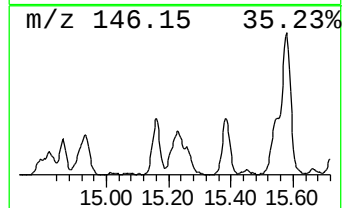
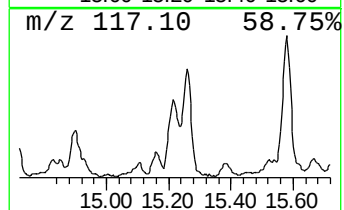
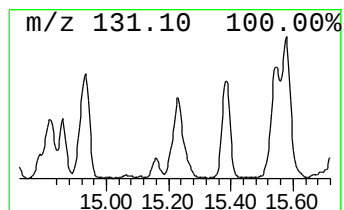
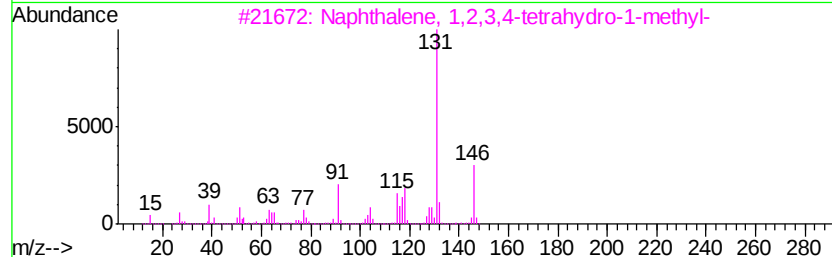
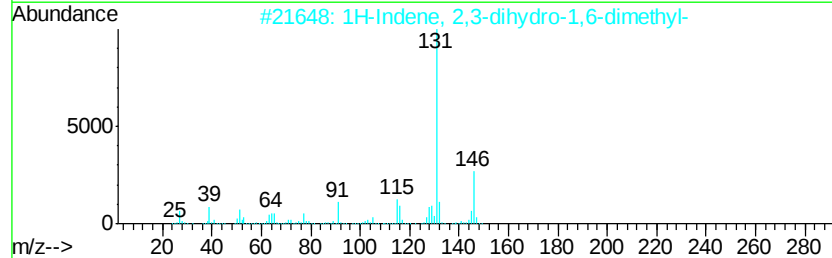
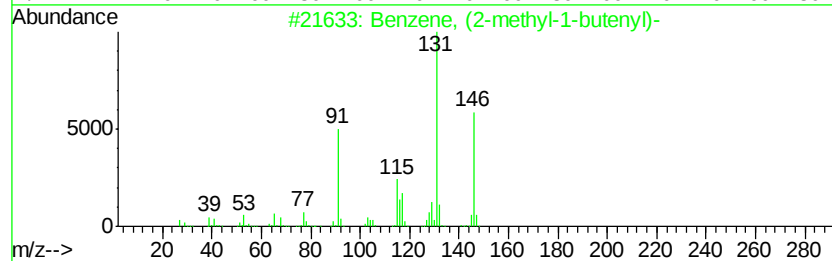
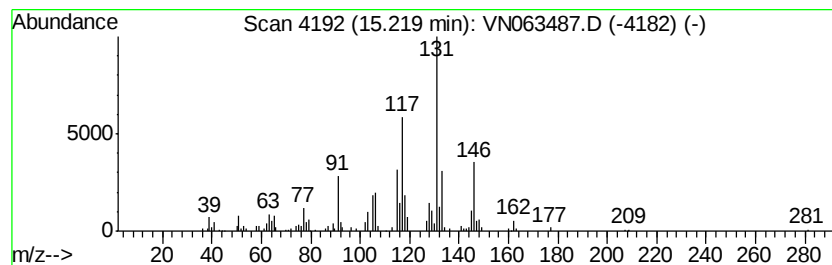
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Peak Number 4 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	10.16 ug/l	352790	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	92
3		Naphthalene, 1,2,3,4-tetrahydro...	146	C11H14	001559-81-5	76
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	76
5		1-Methylindan-2-one	146	C10H10O	035587-60-1	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
229

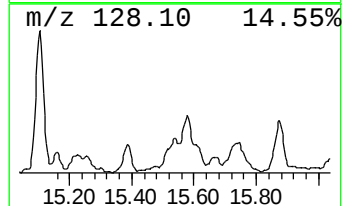
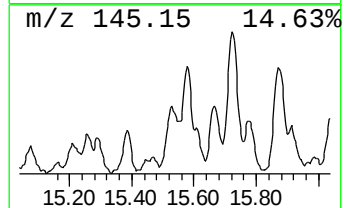
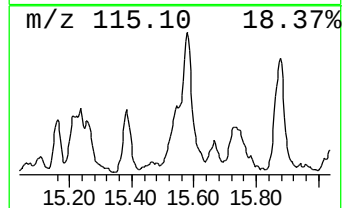
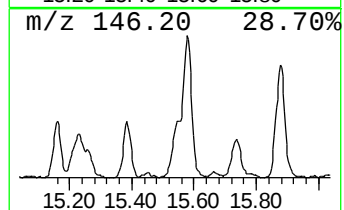
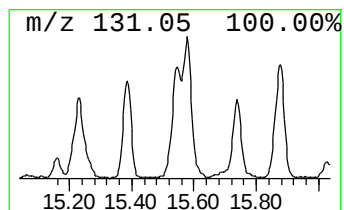
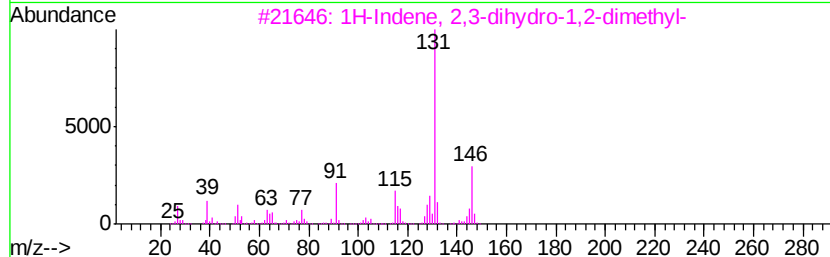
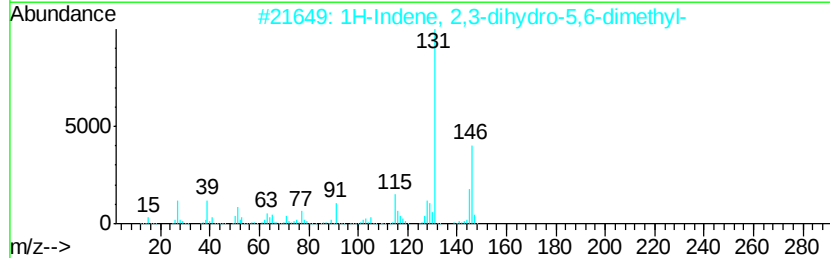
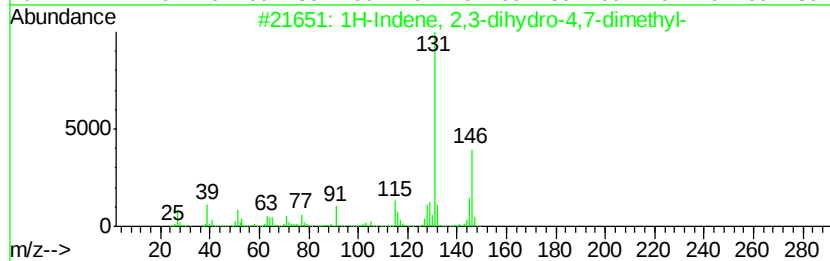
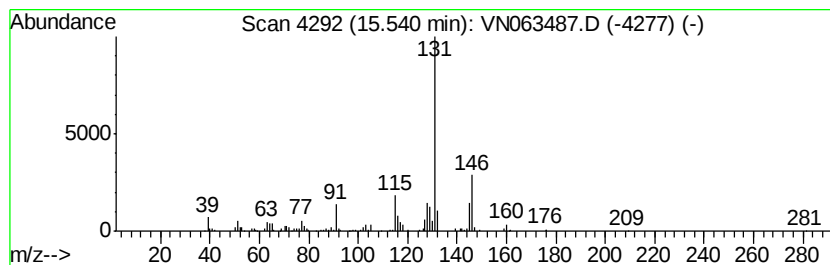
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	5.77 ug/l	200473	1,4-Dichlorobenzene-d4	13.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
5			1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

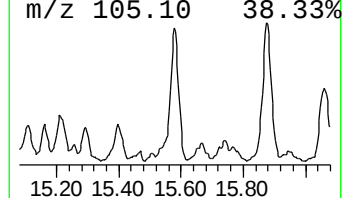
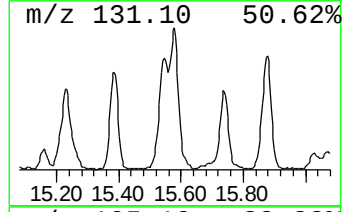
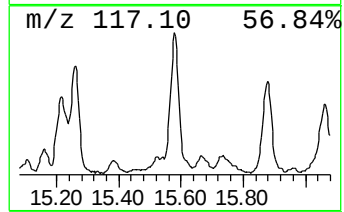
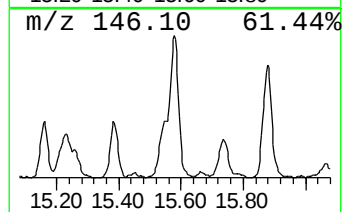
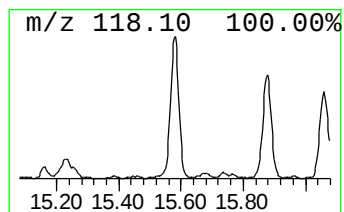
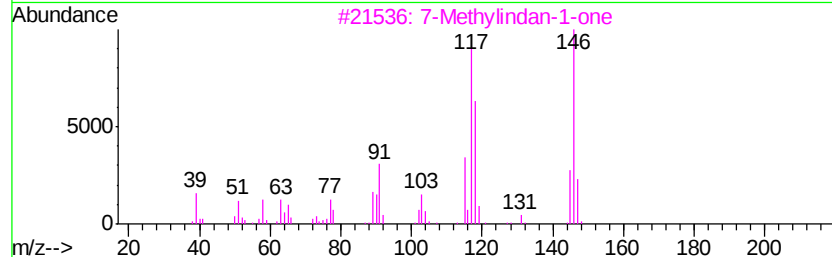
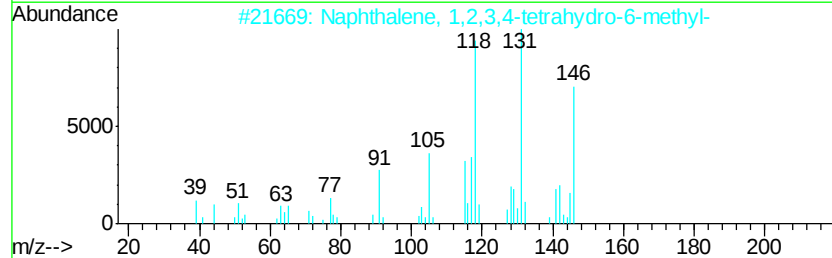
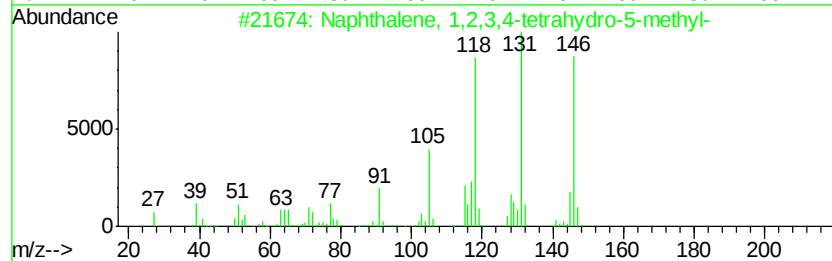
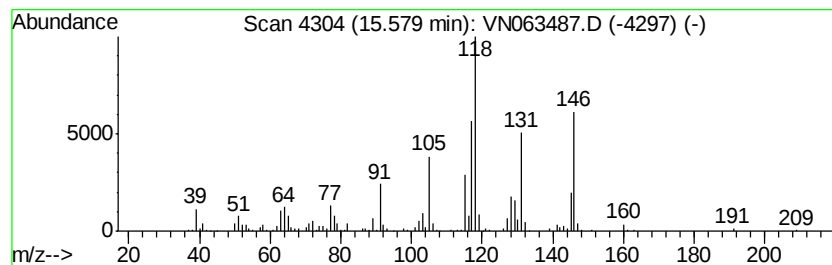
Instrument :
MSVOA_N
ClientSampled :
229

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	14.01 ug/l	486664	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 72
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 62
3		7-Methylindan-1-one	146 C10H10O	039627-61-7 55
4		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 53
5		5H-Benzocycloheptene,6,7,8,9-tet...	146 C11H14	001075-16-7 49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
229

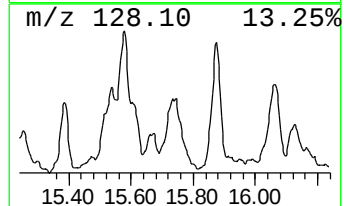
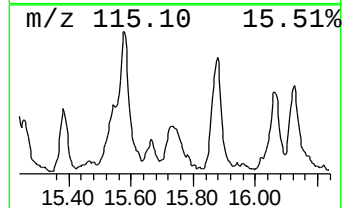
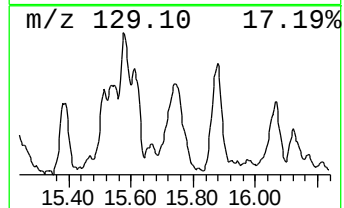
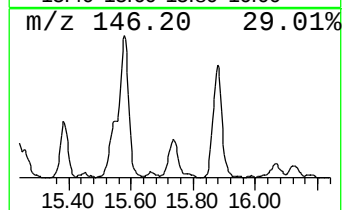
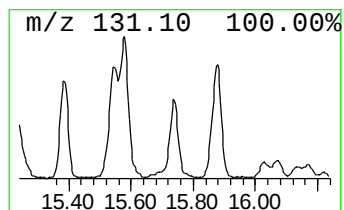
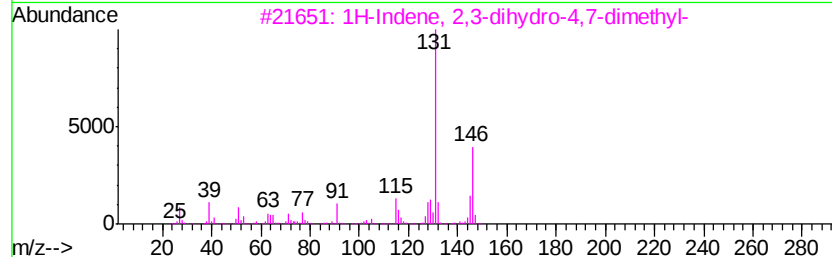
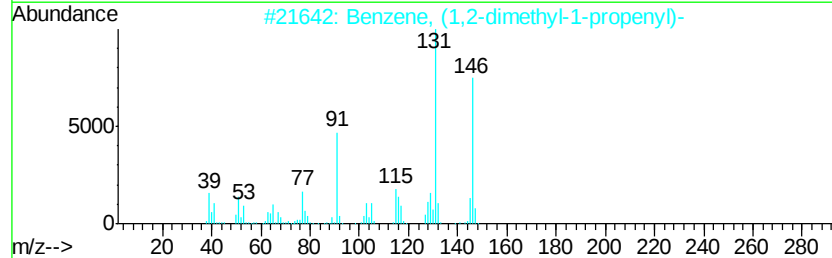
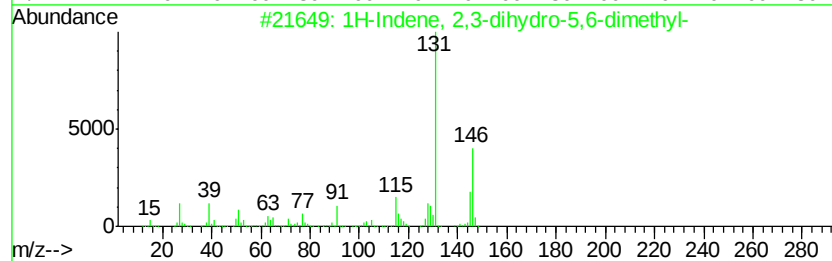
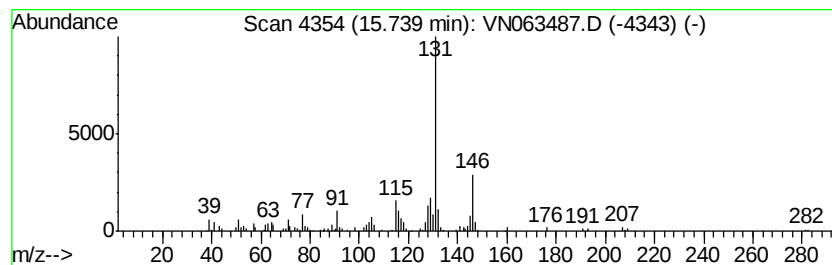
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1H-Indene, 2,3-dihydro-5,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	6.91 ug/l	240074	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	95
2		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	95
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
4		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	94
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

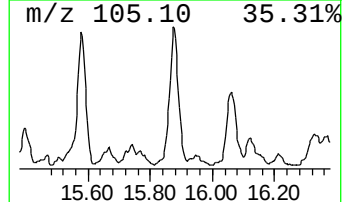
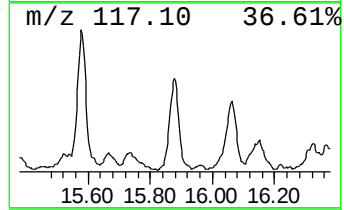
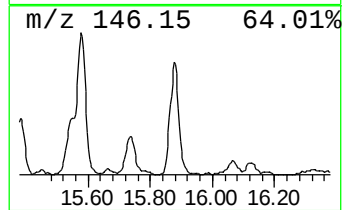
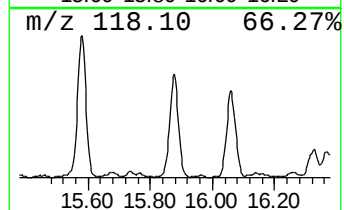
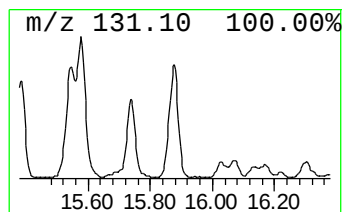
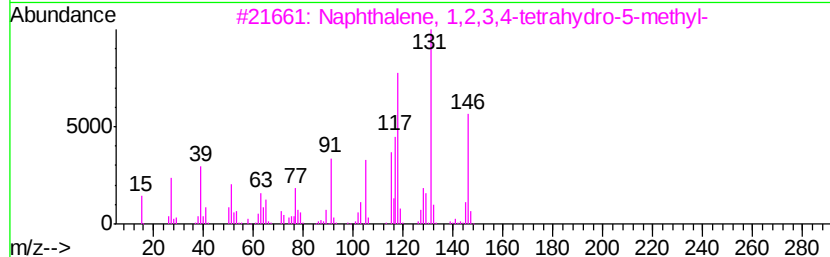
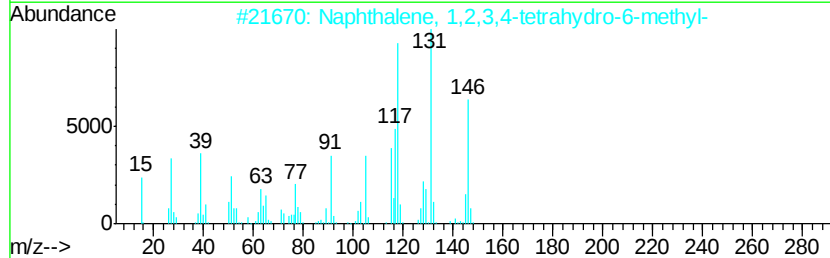
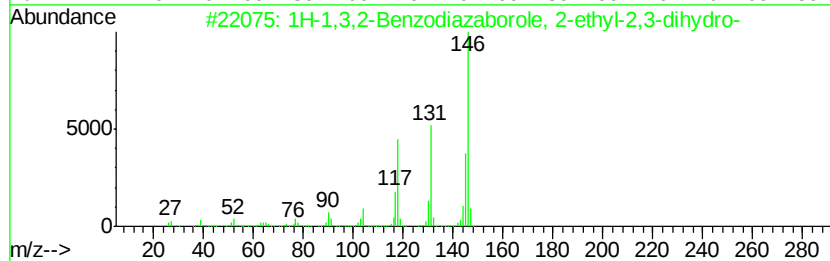
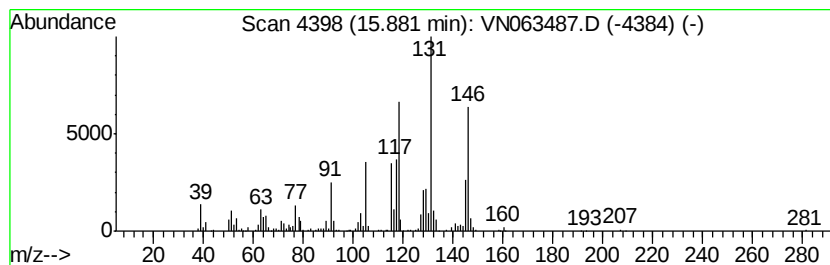
Instrument :
MSVOA_N
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-1,3,2-Benzodiazaborole, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	12.07 ug/l	419238	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 90
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 90
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 87
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 70
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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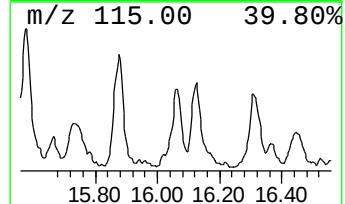
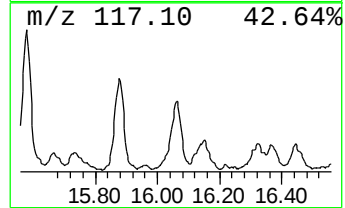
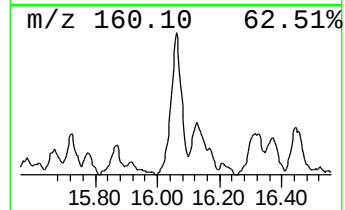
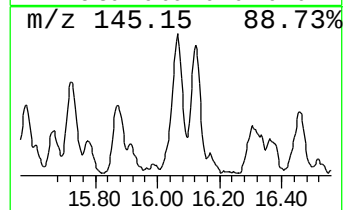
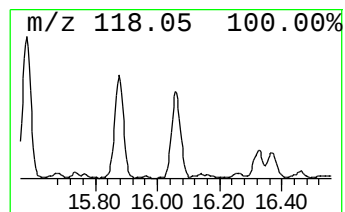
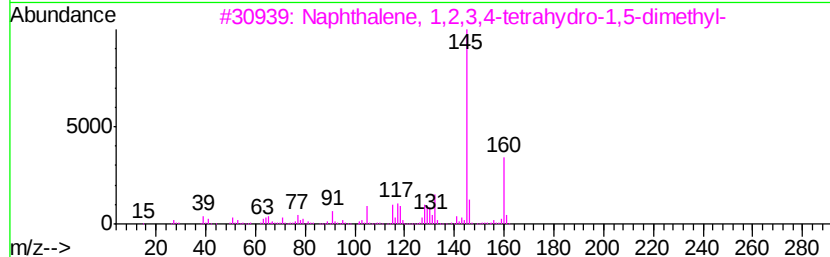
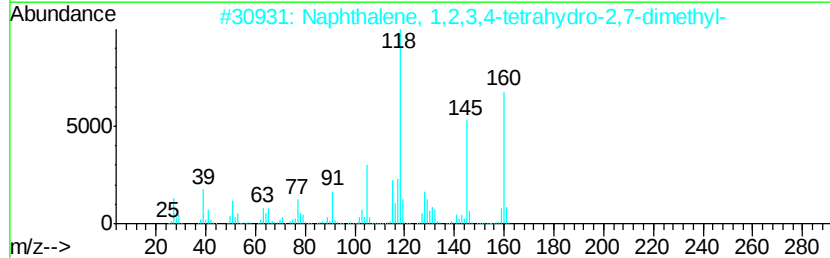
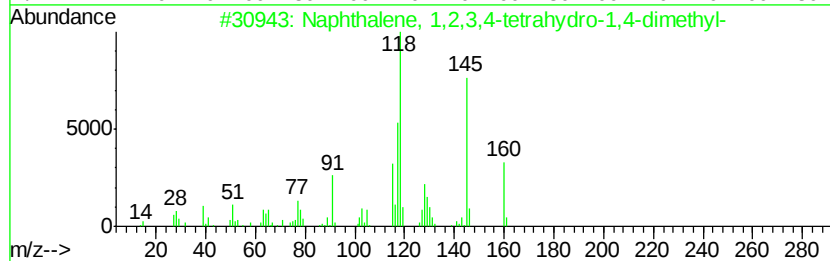
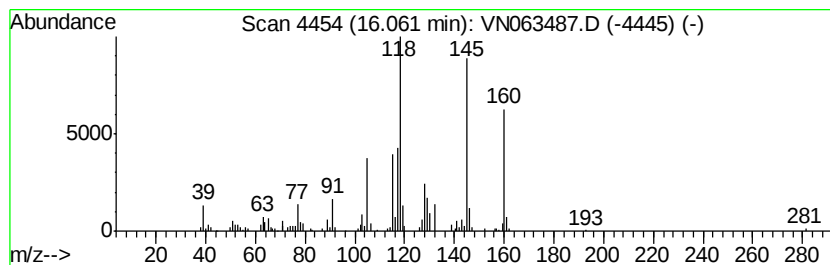
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	5.66 ug/l	196516	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	91
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	64
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	64
4		Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	60
5		Benzene, (1,3-dimethyl-2-butenyl)-	160	C12H16	050704-01-3	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

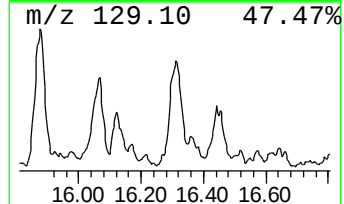
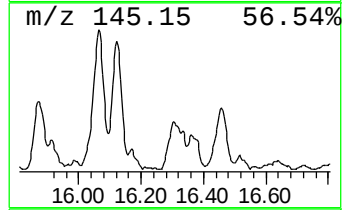
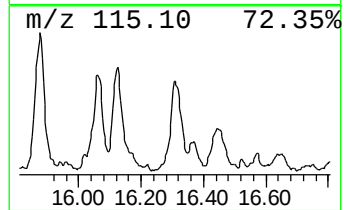
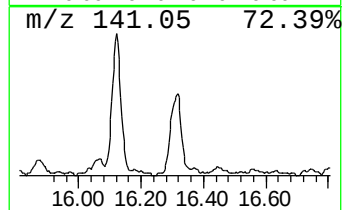
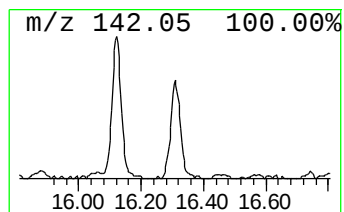
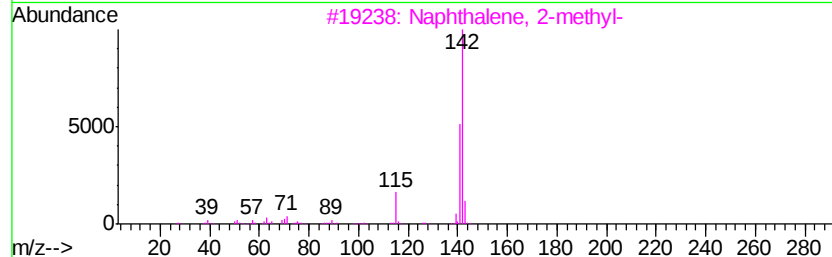
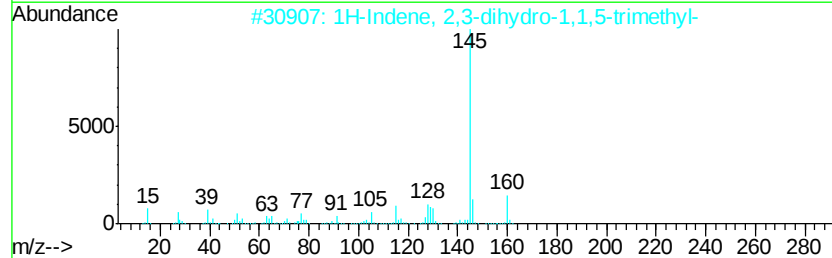
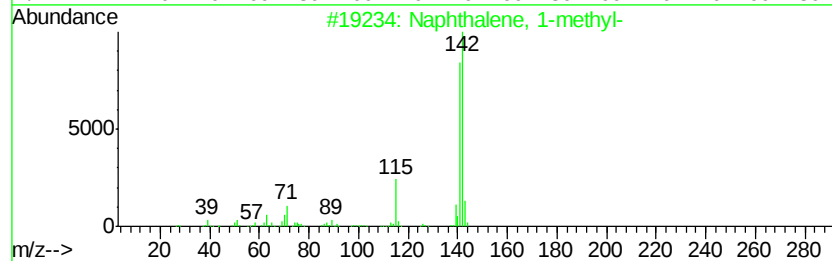
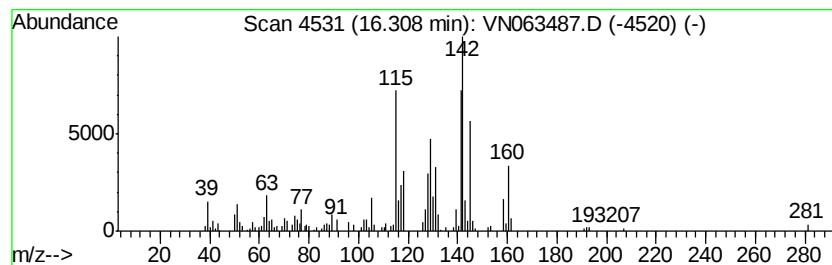
Instrument :
MSVOA_N
ClientSampled :
229

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.31	6.09 ug/l	211581	1,4-Dichlorobenzene-d4	13.32		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	59
2		1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	42
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	41
4		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	38
5		1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	30



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN092120\
Data File : VN063487.D
Acq On : 21 Sep 2020 15:45
Operator : JC/MD
Sample : L4056-03
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
229

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 2-etheny...	14.57	14.6	ug/l	507073	4	13.32	1736840	50.0
1H-Indene, 1-methyl-	14.63	5.7	ug/l	197689	4	13.32	1736840	50.0
Benzene, (2-methy...	15.22	10.2	ug/l	352790	4	13.32	1736840	50.0
1H-Indene, 2,3-di...	15.54	5.8	ug/l	200473	4	13.32	1736840	50.0
Naphthalene, 1,2,...	15.58	14.0	ug/l	486664	4	13.32	1736840	50.0
1H-Indene, 2,3-di...	15.74	6.9	ug/l	240074	4	13.32	1736840	50.0
1H-1,3,2-Benzodia...	15.88	12.1	ug/l	419238	4	13.32	1736840	50.0
Naphthalene, 1,2,...	16.06	5.7	ug/l	196516	4	13.32	1736840	50.0
Naphthalene, 1-me...	16.31	6.1	ug/l	211581	4	13.32	1736840	50.0