

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD061820\
 Data File : VD065802.D
 Acq On : 18 Jun 2020 12:31
 Operator : VA/SY
 Sample : L2811-01
 Misc : 6.52G/5.00ml/MSVOA D/SOIL
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 NB-08-061620

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_D\METHOD\82D060520S.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	2.221	44	51	53	rBV4	4299	7941	1.16%	0.128%
2	2.644	118	123	129	rVB5	5893	12195	1.78%	0.196%
3	4.967	503	518	531	rBV3	32021	109428	15.95%	1.761%
4	5.997	682	693	704	rVB4	12028	38487	5.61%	0.619%
5	7.914	1007	1019	1025	rBV2	108512	266719	38.87%	4.292%
6	7.979	1025	1030	1041	rVB	200559	448430	65.35%	7.217%
7	8.332	1078	1090	1101	rBV	96393	226689	33.04%	3.648%
8	8.867	1172	1181	1194	rBV	300283	613179	89.36%	9.868%
9	10.344	1423	1432	1439	rBV	379649	686188	100.00%	11.043%
10	10.402	1439	1442	1447	rVB3	8237	15009	2.19%	0.242%
11	10.726	1492	1497	1503	rVB3	4067	7966	1.16%	0.128%
12	11.649	1646	1654	1664	rBV	385665	669823	97.62%	10.780%
13	11.749	1666	1671	1676	rVB3	4655	7251	1.06%	0.117%
14	11.849	1680	1688	1694	rBV4	7105	14090	2.05%	0.227%
15	12.191	1741	1746	1752	rBV6	4635	8888	1.30%	0.143%
16	12.638	1815	1822	1829	rVB	219704	372800	54.33%	6.000%
17	12.920	1861	1870	1873	rBV5	2960	7742	1.13%	0.125%
18	13.279	1926	1931	1939	rVB3	6691	11305	1.65%	0.182%
19	13.514	1967	1971	1976	rBV5	5221	9070	1.32%	0.146%
20	13.579	1976	1982	1992	rVB	281744	479588	69.89%	7.718%
21	13.779	2011	2016	2019	rVV2	25986	41802	6.09%	0.673%
22	13.820	2019	2023	2032	rVB4	27366	55242	8.05%	0.889%
23	13.902	2032	2037	2042	rBV5	14221	27939	4.07%	0.450%
24	13.973	2044	2049	2056	rBV	14214	22966	3.35%	0.370%
25	14.038	2056	2060	2061	rBV2	7319	7569	1.10%	0.122%
26	14.067	2061	2065	2069	rVB2	12923	19491	2.84%	0.314%
27	14.120	2069	2074	2080	rVB3	30907	52971	7.72%	0.852%
28	14.185	2080	2085	2088	rBV3	9593	16489	2.40%	0.265%
29	14.232	2088	2093	2099	rVB2	48411	80227	11.69%	1.291%
30	14.285	2099	2102	2104	rBV3	4914	7048	1.03%	0.113%
31	14.361	2109	2115	2120	rBV4	16200	29507	4.30%	0.475%
32	14.414	2120	2124	2127	rBV5	13778	23802	3.47%	0.383%
33	14.449	2127	2130	2133	rVB2	9437	9979	1.45%	0.161%
34	14.485	2133	2136	2140	rVB	23750	30474	4.44%	0.490%

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ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
NB-08-061620

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

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35	14.532	2140	2144	2147	rBV2	25239	32740	4.77%	0.527%
36	14.573	2147	2151	2158	rVB5	23219	40770	5.94%	0.656%
37	14.632	2158	2161	2163	rBV4	10430	15015	2.19%	0.242%
38	14.667	2163	2167	2172	rVB3	27632	40280	5.87%	0.648%
39	14.720	2172	2176	2181	rBV4	28484	61546	8.97%	0.990%
40	14.761	2181	2183	2188	rVB	23499	28448	4.15%	0.458%
41	14.838	2188	2196	2202	rBV3	123276	282084	41.11%	4.540%
42	14.890	2202	2205	2215	rVB6	25122	44073	6.42%	0.709%
43	14.990	2216	2222	2230	rVB	103285	173646	25.31%	2.795%
44	15.067	2230	2235	2236	rBV4	12426	19487	2.84%	0.314%
45	15.102	2236	2241	2244	rBV3	34707	57891	8.44%	0.932%
46	15.220	2253	2261	2267	rVB4	56302	138485	20.18%	2.229%
47	15.385	2280	2289	2290	rBV7	12315	26930	3.92%	0.433%
48	15.408	2290	2293	2297	rVB5	11069	16546	2.41%	0.266%
49	15.461	2297	2302	2307	rBV3	49853	95212	13.88%	1.532%
50	15.514	2307	2311	2314	rVV6	30739	59022	8.60%	0.950%
51	15.543	2314	2316	2335	rVB8	44330	137769	20.08%	2.217%
52	15.702	2335	2343	2350	rBV5	20413	53705	7.83%	0.864%
53	15.779	2350	2356	2362	rBV9	5462	11214	1.63%	0.180%
54	15.873	2364	2372	2374	rBV6	24073	48179	7.02%	0.775%
55	15.914	2374	2379	2389	rVB3	66395	135920	19.81%	2.187%
56	16.008	2389	2395	2400	rBV8	10122	20128	2.93%	0.324%
57	16.073	2400	2406	2407	rBV2	16003	20918	3.05%	0.337%
58	16.232	2426	2433	2440	rBV5	38561	88870	12.95%	1.430%
59	16.432	2456	2467	2474	rBV4	35363	94306	13.74%	1.518%
60	16.508	2474	2480	2486	rVV8	8980	21498	3.13%	0.346%
61	16.696	2507	2512	2513	rBV4	7078	10758	1.57%	0.173%

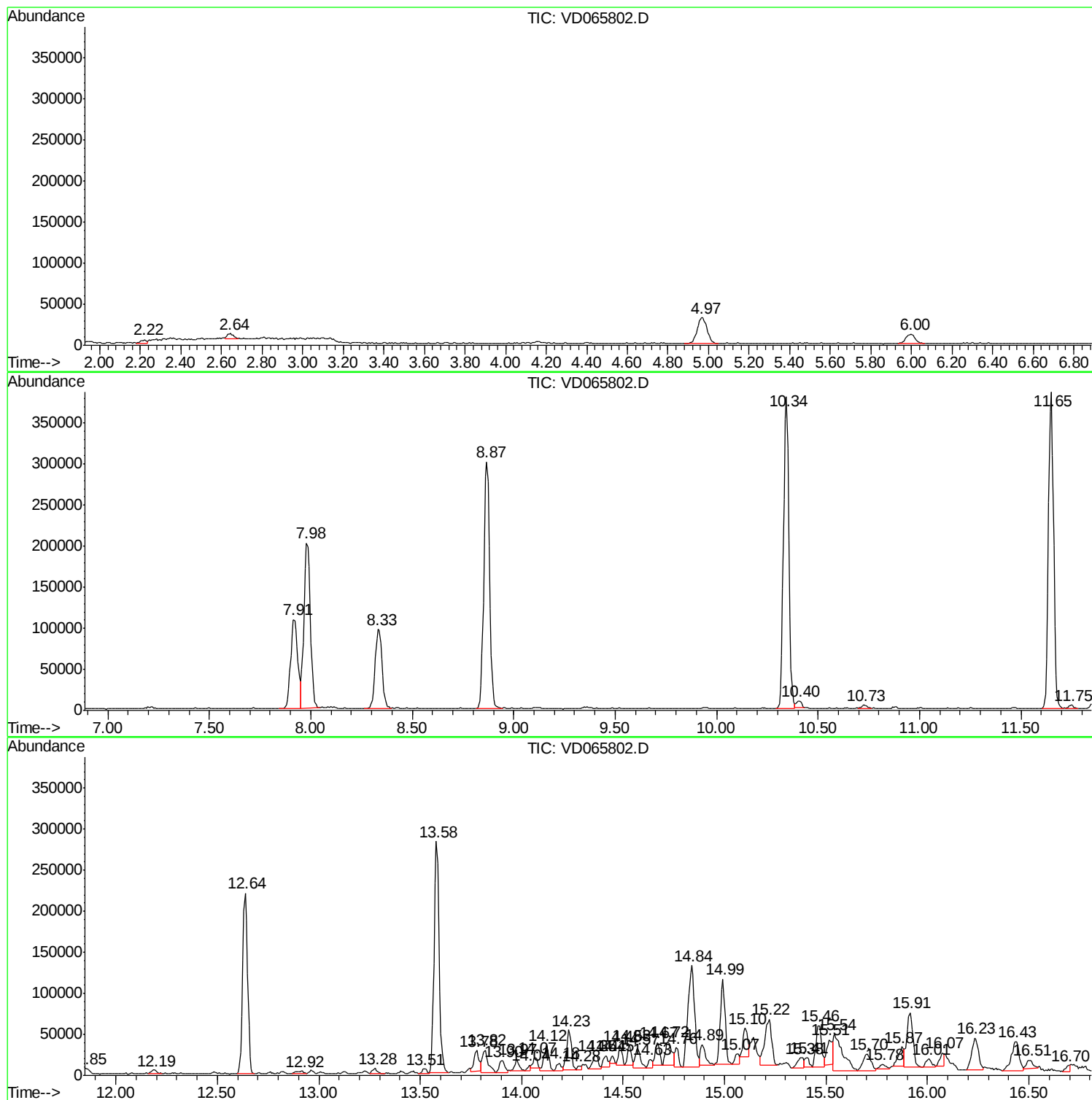
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Quant Title : SW846 8260

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



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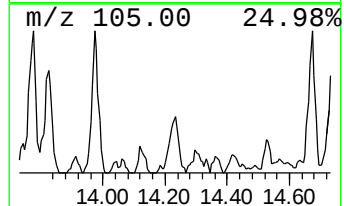
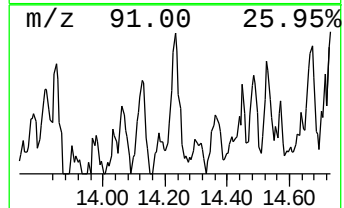
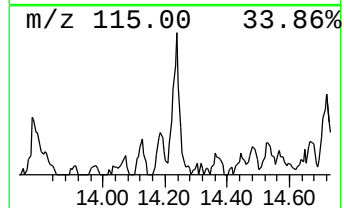
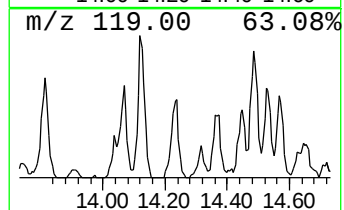
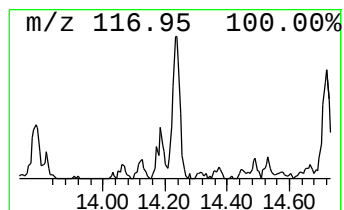
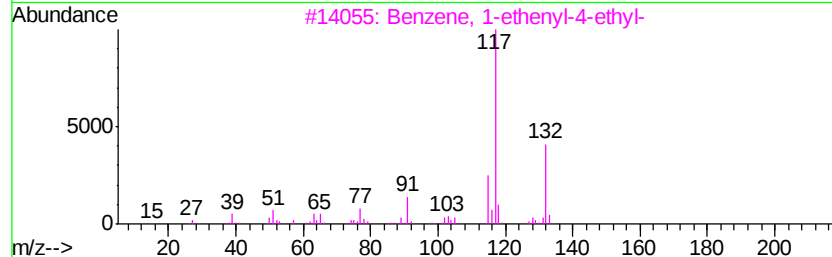
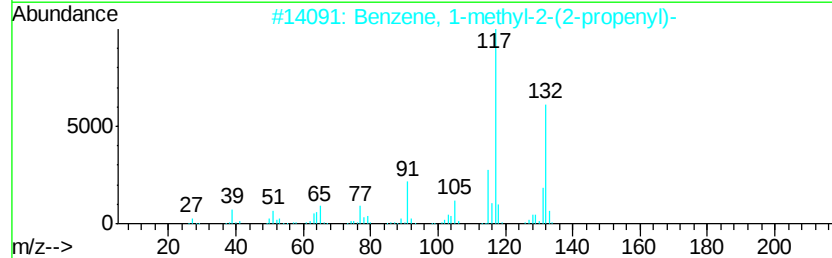
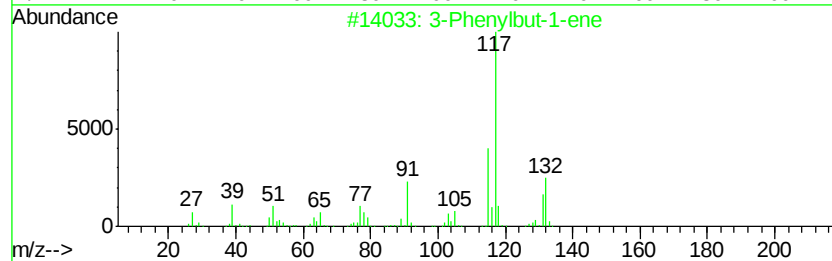
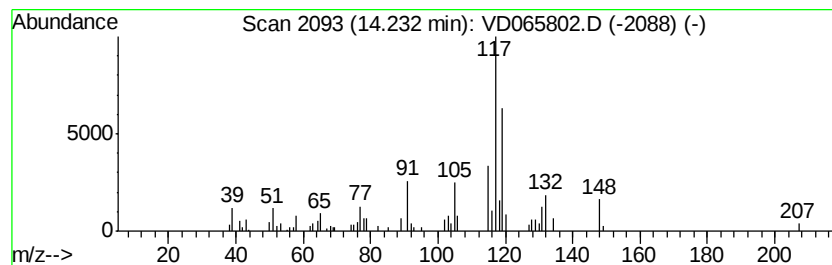
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

Peak Number 1 3-Phenylbut-1-ene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	8.36 ug/l	80227	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	60
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	58
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	58



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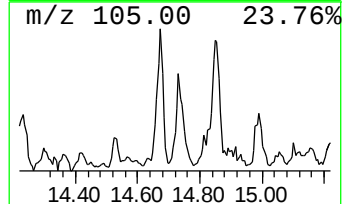
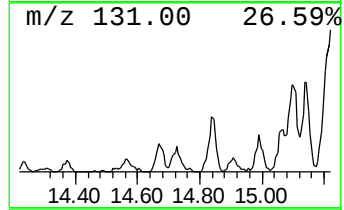
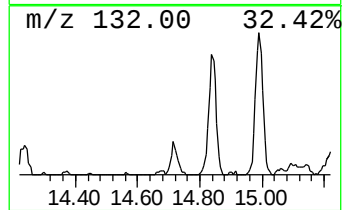
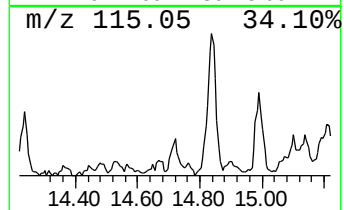
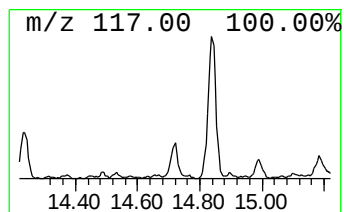
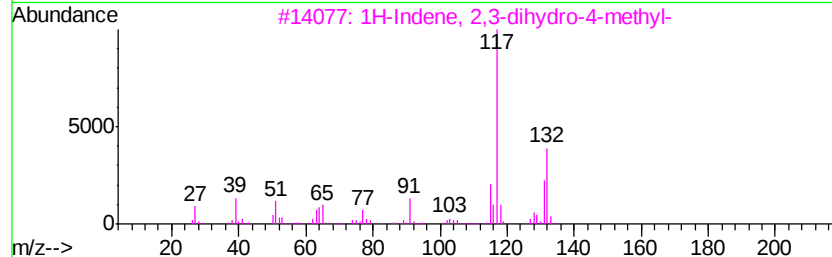
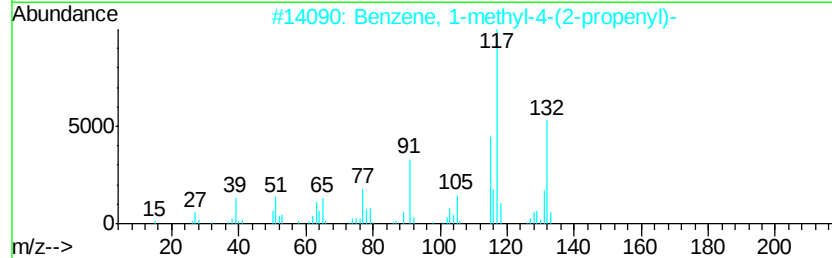
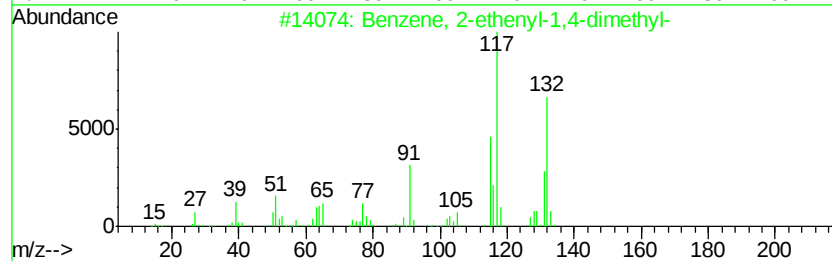
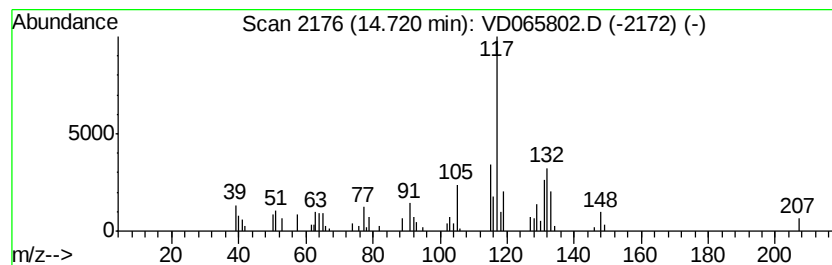
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	6.42 ug/l	61546	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	62
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	62
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	62



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ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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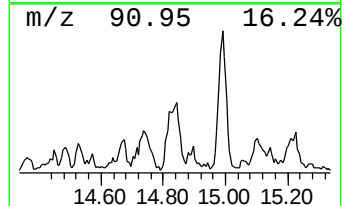
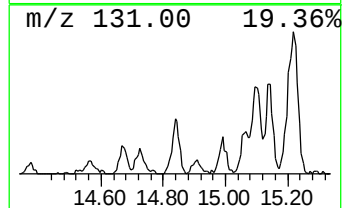
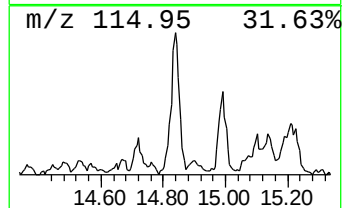
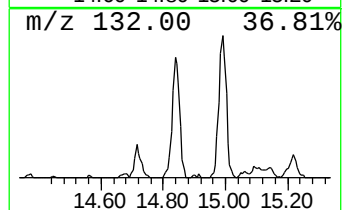
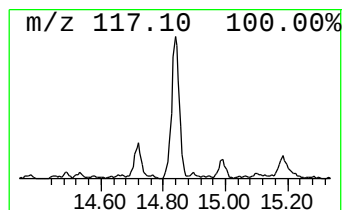
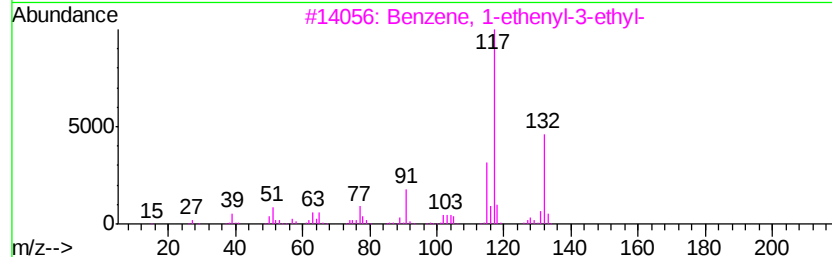
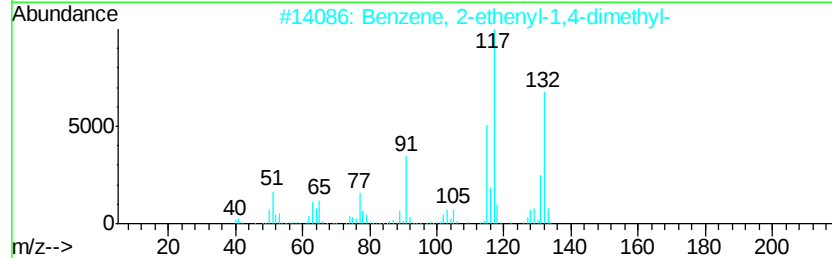
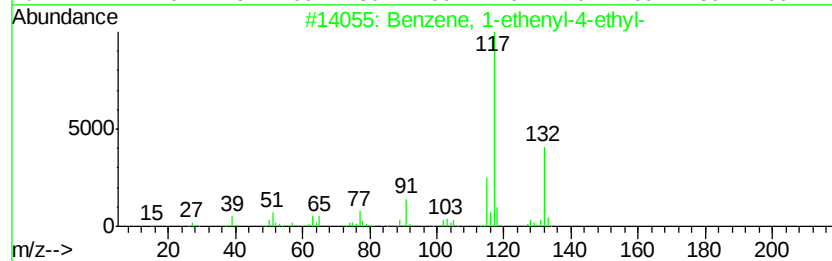
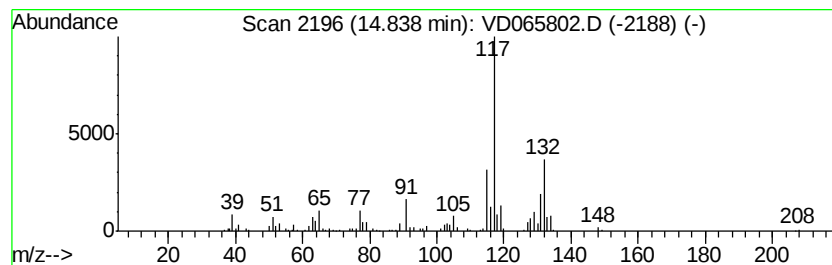
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	29.41 ug/l	282084	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87



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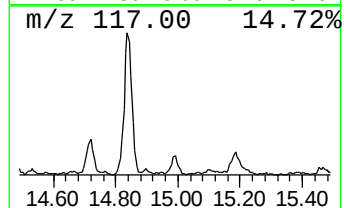
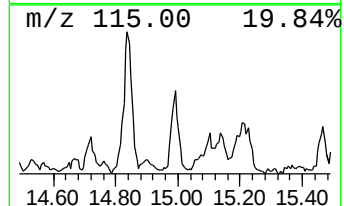
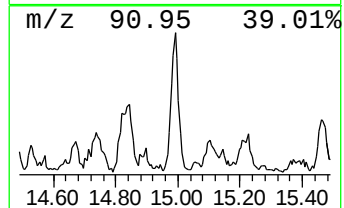
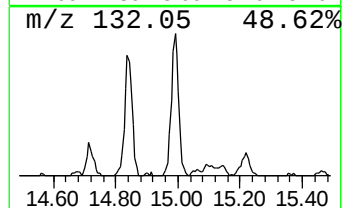
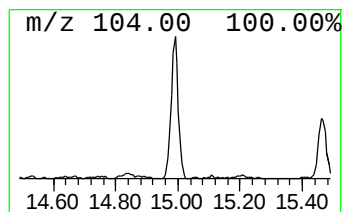
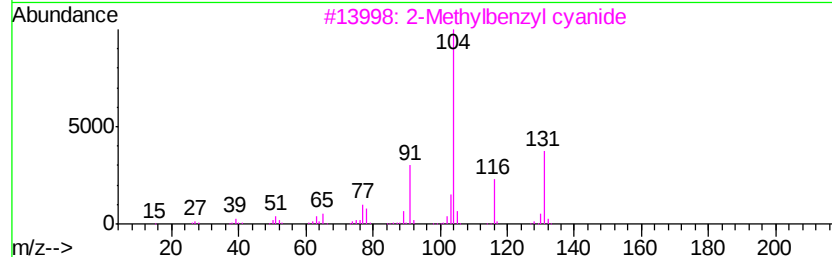
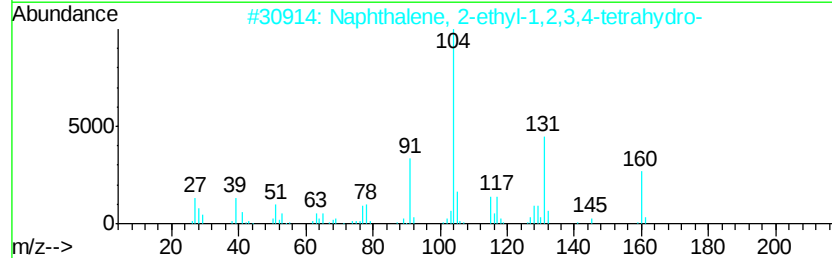
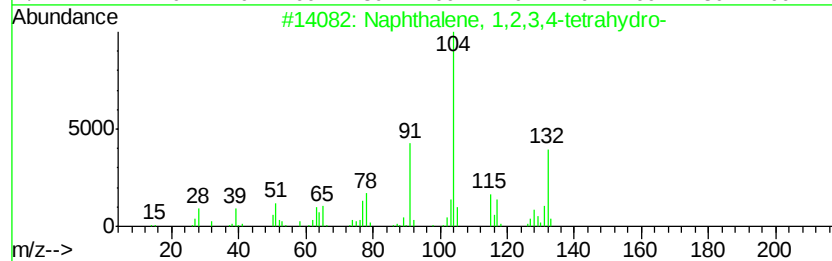
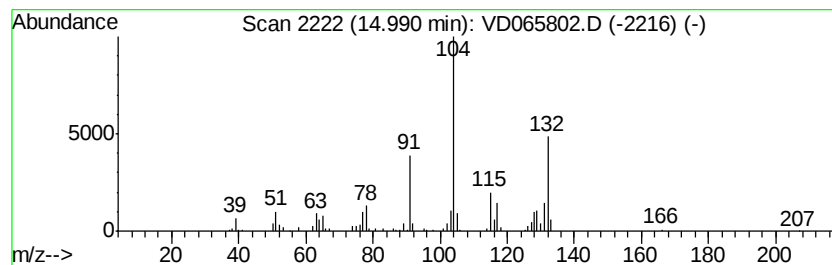
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	18.10 ug/l	173646	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	59
3		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	35
4		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	35
5		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	35



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Misc : 6.52G/5.00ml/MSVOA D/SOIL
ALS Vial : 5 Sample Multiplier: 1

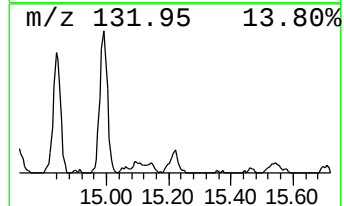
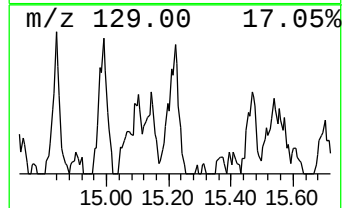
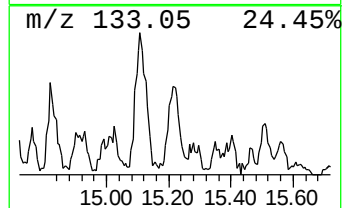
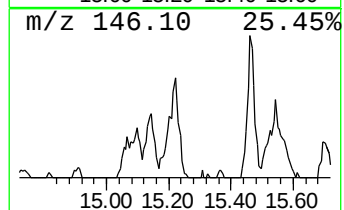
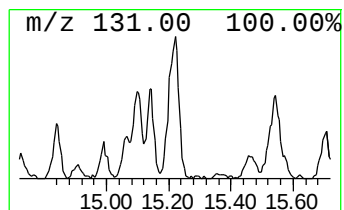
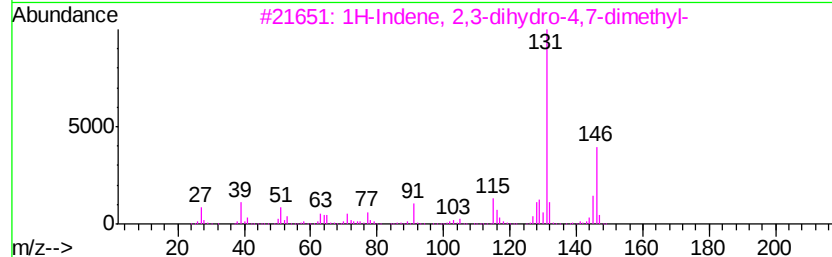
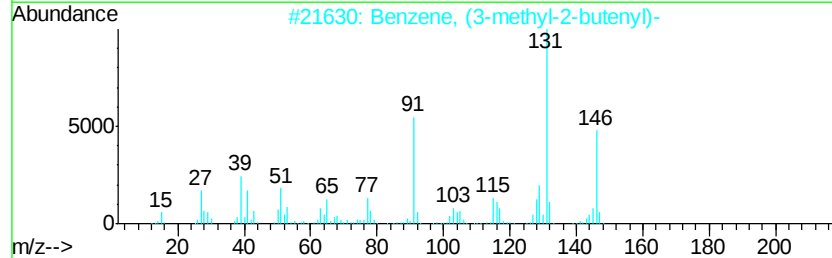
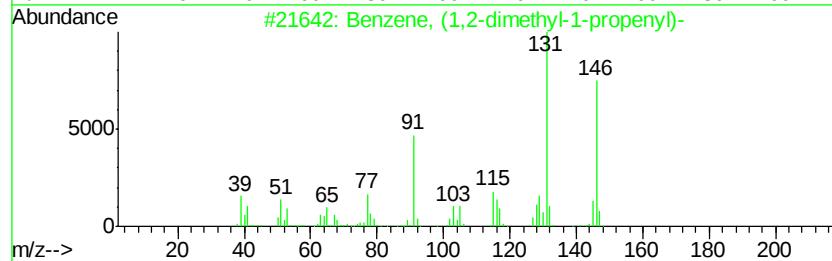
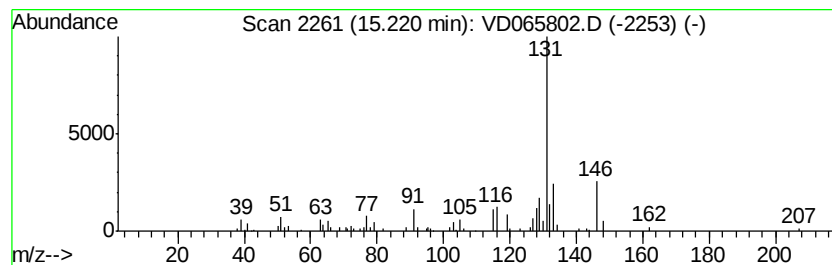
Instrument :
MSVOA_D
ClientSampleId :
NB-08-061620

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	14.44 ug/l	138485	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 87
2		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 87
3		1H-Indene, 2,3-dihydro-4,7-dimethyl-	146 C11H14	006682-71-9 87
4		1H-Indene, 2,3-dihydro-1,6-dimethyl-	146 C11H14	017059-48-2 83
5		Bicyclo[4.2.0]octa-1,3,5-triene,	146 C11H14	027087-54-3 81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD061820\
Data File : VD065802.D
Acq On : 18 Jun 2020 12:31
Operator : VA/SY
Sample : L2811-01
Misc : 6.52G/5.00ml/MSVOA D/SOIL
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
NB-08-061620

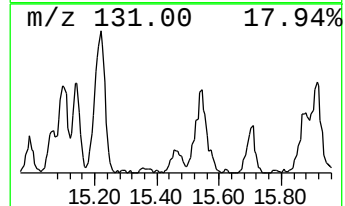
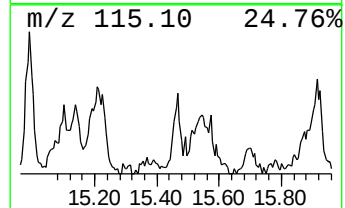
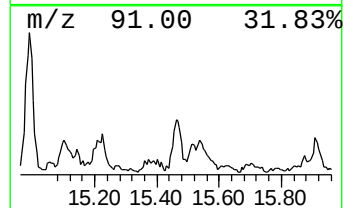
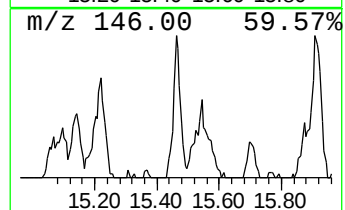
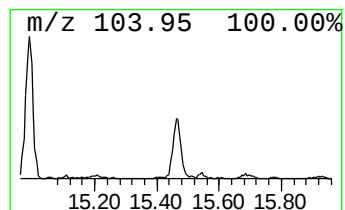
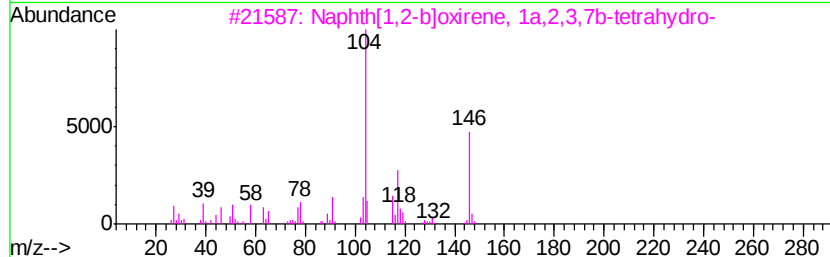
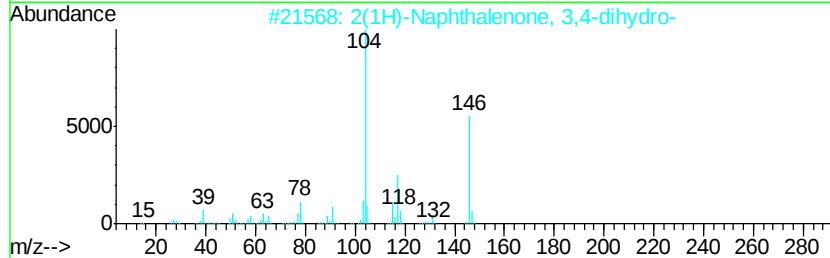
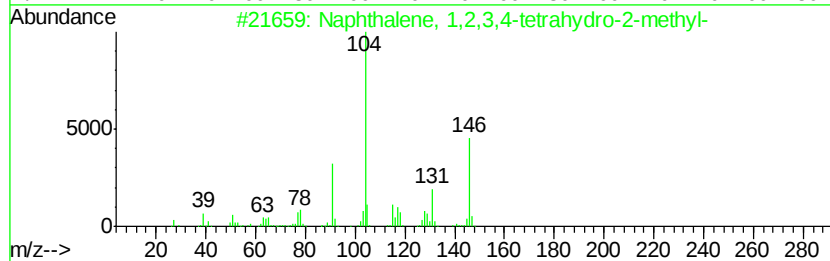
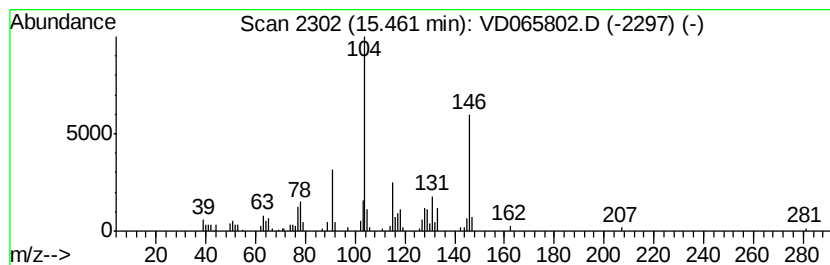
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	9.93 ug/l	95212	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	76
3		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	68
4		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	45
5		2-Azetidinone, 4-phenyl-	147	C9H9NO	005661-55-2	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD061820\
Data File : VD065802.D
Acq On : 18 Jun 2020 12:31
Operator : VA/SY
Sample : L2811-01
Misc : 6.52G/5.00ml/MSVOA D/SOIL
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
NB-08-061620

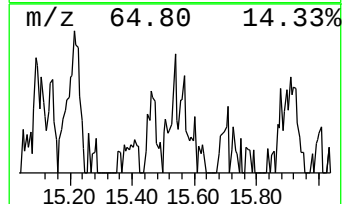
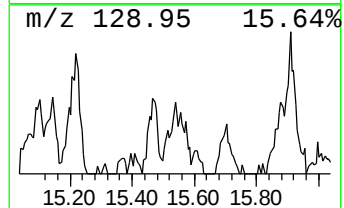
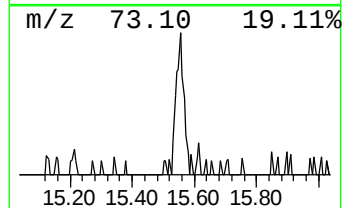
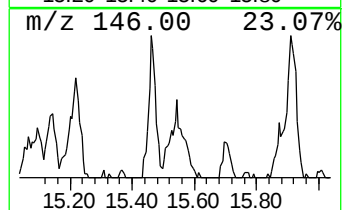
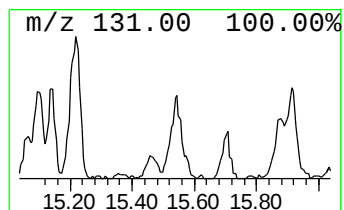
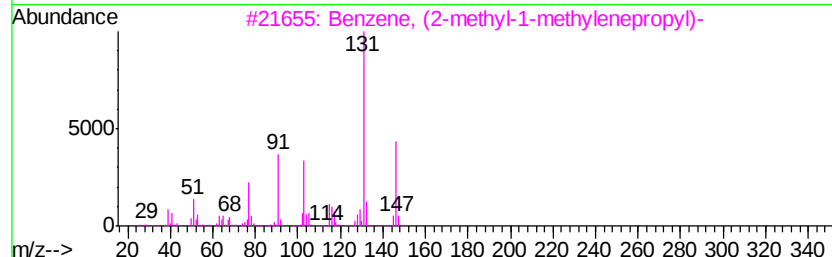
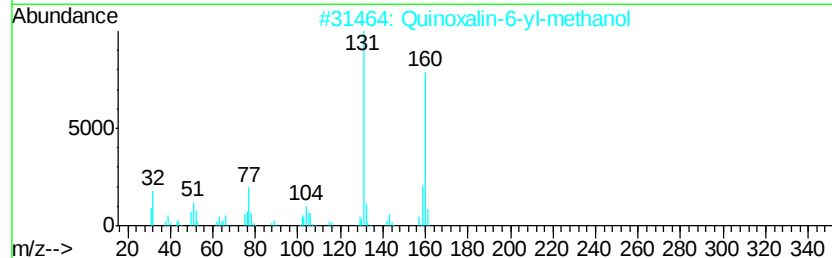
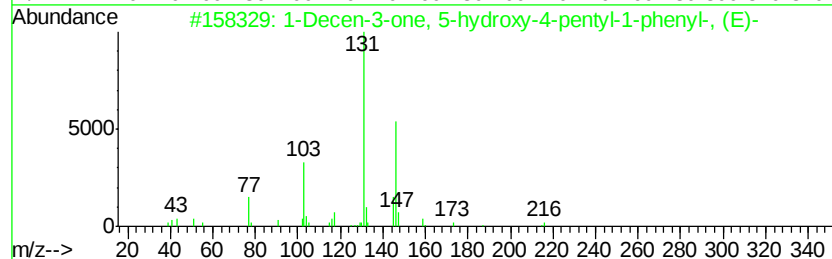
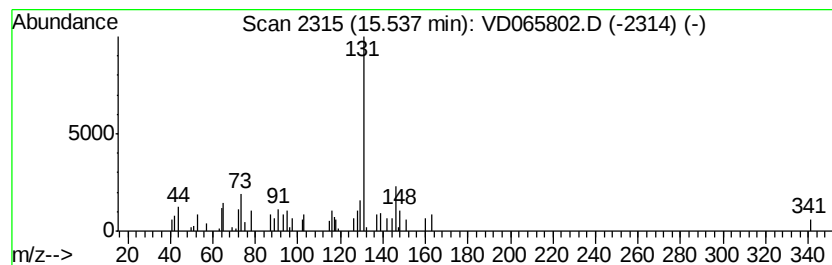
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Quant Title : SW846 8260

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

Peak Number 7 unknown15.54 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	14.36 ug/l	137769	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Decen-3-one, 5-hydroxy-4-pentyl...	316	C21H32O2	1000115-33-2	40
2		Quinoxalin-6-yl-methanol	160	C9H8N2O	1000318-21-1	35
3		Benzene, (2-methyl-1-methylenepropyl)-	146	C11H14	017498-71-4	33
4		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	28
5		Aziridine, 1,2-dimethyl-3-phenyl-	147	C10H13N	000936-43-6	14



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD061820\
Data File : VD065802.D
Acq On : 18 Jun 2020 12:31
Operator : VA/SY
Sample : L2811-01
Misc : 6.52G/5.00ml/MSVOA D/SOIL
ALS Vial : 5 Sample Multiplier: 1

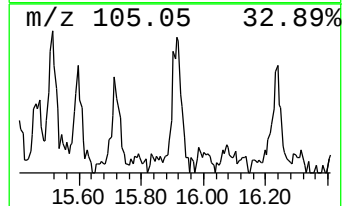
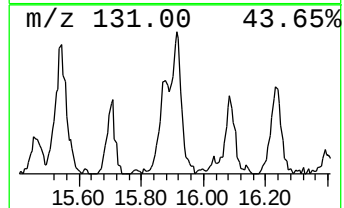
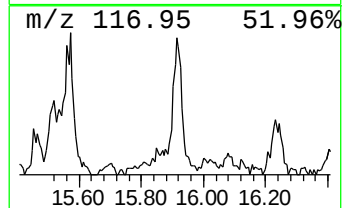
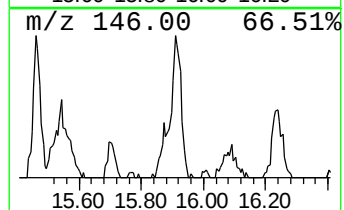
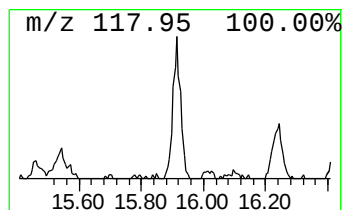
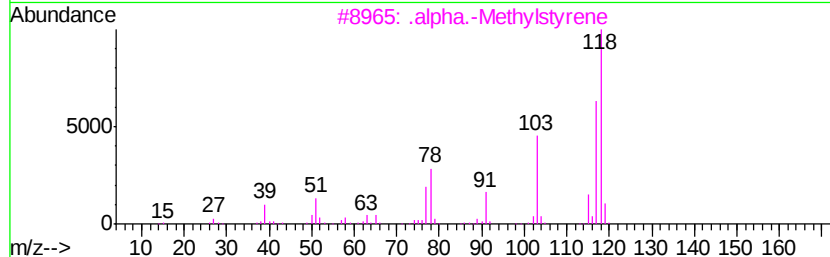
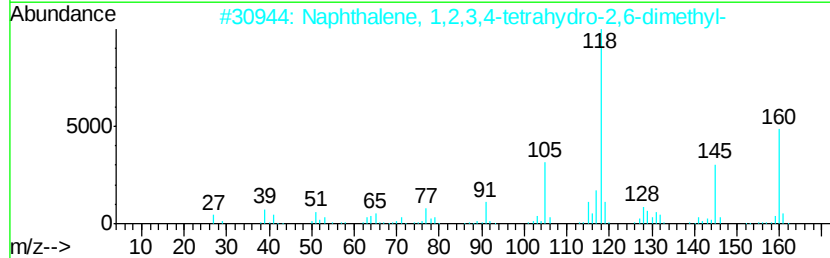
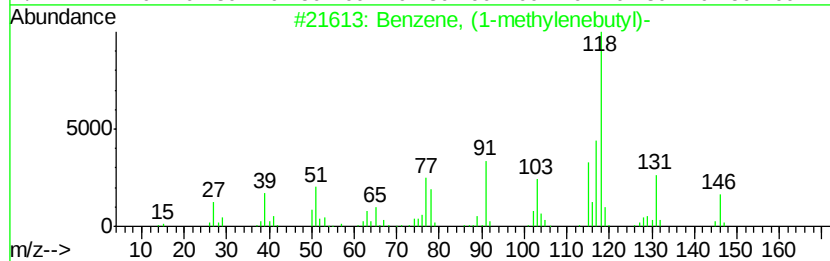
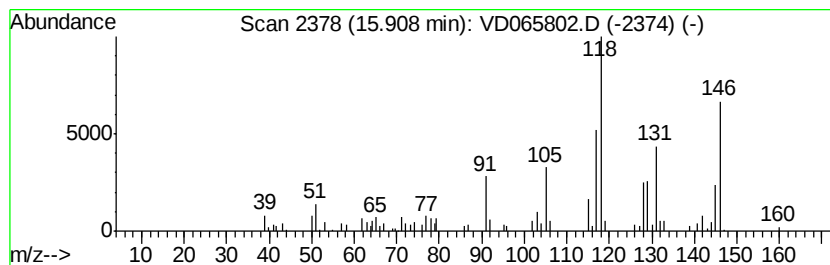
Instrument :
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ClientSampleId :
NB-08-061620

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methylenebutyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	14.17 ug/l	135920	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methylenebutyl)-	146 C11H14	005676-32-4 74
2		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	007524-63-2 43
3		.alpha.-Methylstyrene	118 C9H10	000098-83-9 38
4		Phthalic acid, methyl 3-phenylpr...	298 C18H18O4	1000308-99-2 38
5		3-Chloropropionic acid, 3-phenyl...	226 C12H15ClO2	078987-69-6 38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD061820\
Data File : VD065802.D
Acq On : 18 Jun 2020 12:31
Operator : VA/SY
Sample : L2811-01
Misc : 6.52G/5.00ml/MSVOA D/SOIL
ALS Vial : 5 Sample Multiplier: 1

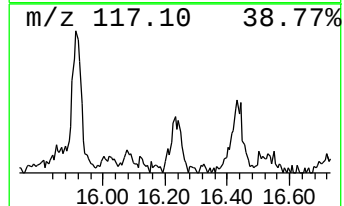
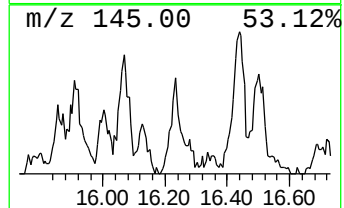
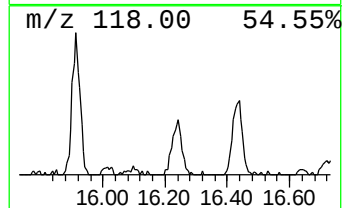
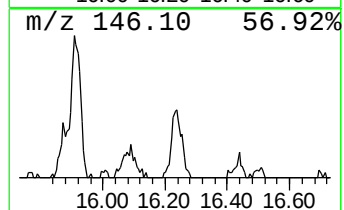
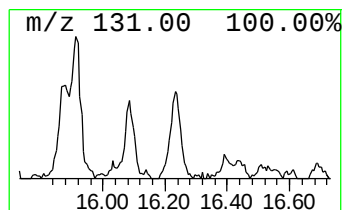
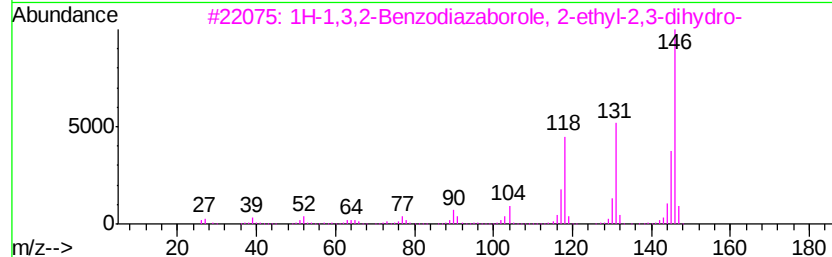
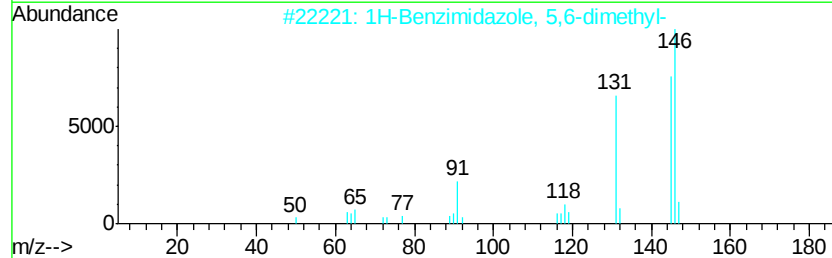
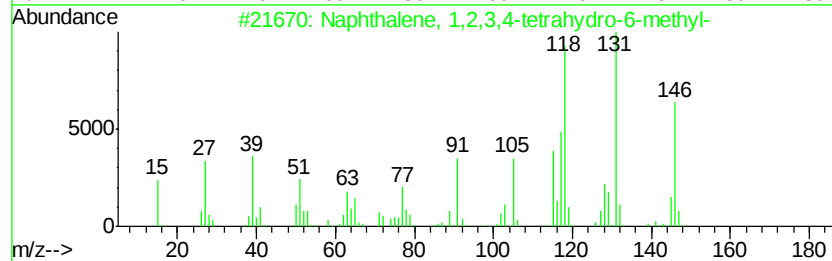
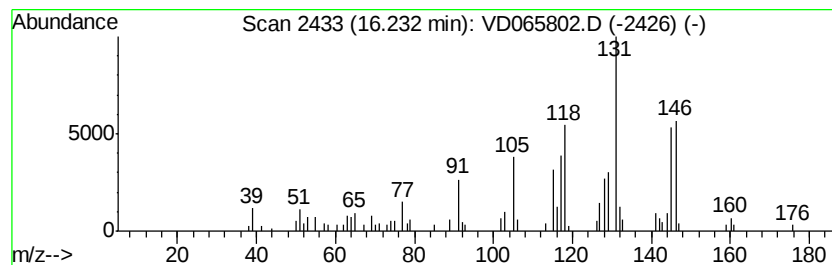
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ClientSampleId :
NB-08-061620

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
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 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	9.27 ug/l	88870	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 78
2		1H-Benzimidazole, 5,6-dimethyl-	146 C9H10N2	000582-60-5 60
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 58
4		2-Propyn-1-ol, 3-(4-methylphenyl)-	146 C10H10O	016017-24-6 53
5		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD061820\
Data File : VD065802.D
Acq On : 18 Jun 2020 12:31
Operator : VA/SY
Sample : L2811-01
Misc : 6.52G/5.00ml/MSVOA D/SOIL
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
NB-08-061620

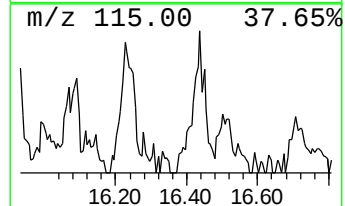
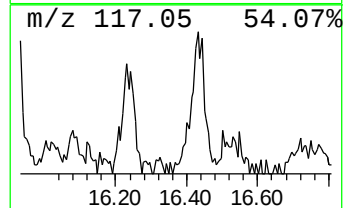
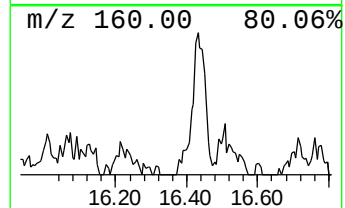
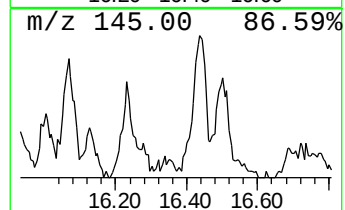
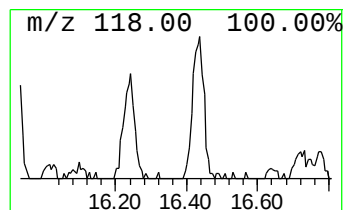
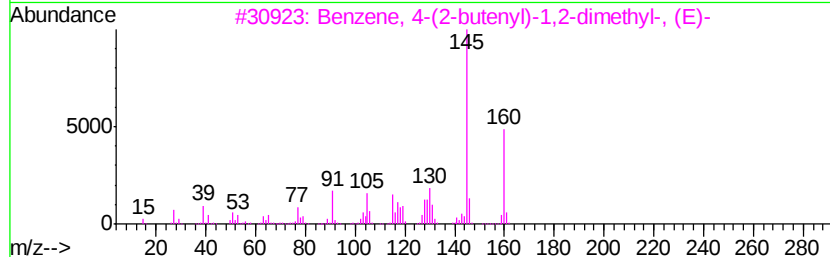
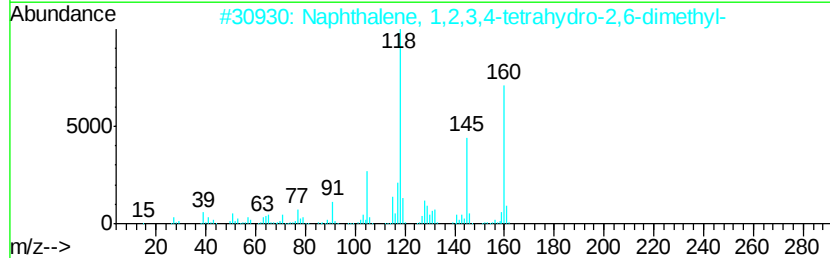
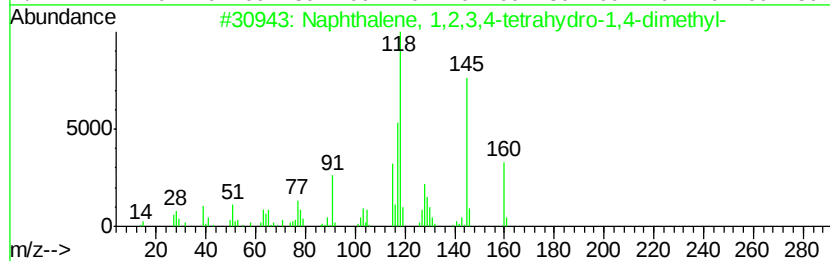
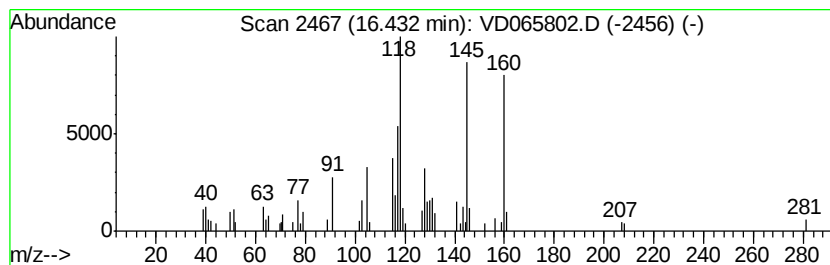
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	9.83 ug/l	94306	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	94
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	83
3		Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	76
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	60
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	49



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD061820\
Data File : VD065802.D
Acq On : 18 Jun 2020 12:31
Operator : VA/SY
Sample : L2811-01
Misc : 6.52G/5.00ml/MSVOA_D/SOIL
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
NB-08-061620

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D060520S.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
3-Phenylbut-1-ene	14.23	8.4	ug/l	80227	4	13.58	479588	50.0
Benzene, 2-etheny...	14.72	6.4	ug/l	61546	4	13.58	479588	50.0
Benzene, 1-etheny...	14.84	29.4	ug/l	282084	4	13.58	479588	50.0
Naphthalene, 1,2,...	14.99	18.1	ug/l	173646	4	13.58	479588	50.0
Benzene, (1,2-dim...	15.22	14.4	ug/l	138485	4	13.58	479588	50.0
Naphthalene, 1,2,...	15.46	9.9	ug/l	95212	4	13.58	479588	50.0
unknown15.54	15.54	14.4	ug/l	137769	4	13.58	479588	50.0
Benzene, (1-methy...	15.91	14.2	ug/l	135920	4	13.58	479588	50.0
Naphthalene, 1,2,...	16.23	9.3	ug/l	88870	4	13.58	479588	50.0
Naphthalene, 1,2,...	16.43	9.8	ug/l	94306	4	13.58	479588	50.0