

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD090420\
 Data File : VD066571.D
 Acq On : 04 Sep 2020 18:57
 Operator : VA/SY
 Sample : L3896-01
 Misc : 5.01G/5.00ml/MSVOA D/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 WC-1

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_D\METHOD\82D090320S.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	4.967	506	518	525	rBV3	27287	88608	1.29%	0.240%
2	7.191	888	896	908	rBV	31157	87915	1.28%	0.238%
3	7.479	936	945	956	rVB3	48720	125284	1.83%	0.340%
4	7.914	1009	1019	1024	rBV	362336	896680	13.10%	2.432%
5	7.979	1024	1030	1040	rVV2	618224	1408076	20.57%	3.820%
6	8.326	1080	1089	1101	rVB	344064	792604	11.58%	2.150%
7	8.720	1149	1156	1162	rBV4	37902	80313	1.17%	0.218%
8	8.861	1169	1180	1189	rBV	885011	1811595	26.46%	4.914%
9	9.350	1251	1263	1269	rBV3	182225	476563	6.96%	1.293%
10	9.620	1300	1309	1317	rBV6	53583	130929	1.91%	0.355%
11	9.767	1329	1334	1345	rVB4	50333	100502	1.47%	0.273%
12	9.973	1363	1369	1373	rBV3	71232	129304	1.89%	0.351%
13	10.102	1383	1391	1406	rVB4	74107	225297	3.29%	0.611%
14	10.255	1406	1417	1423	rBV10	26137	81495	1.19%	0.221%
15	10.338	1423	1431	1445	rBV	1378811	2666141	38.94%	7.233%
16	10.514	1454	1461	1470	rVB6	93417	252380	3.69%	0.685%
17	10.673	1484	1488	1494	rVB	85923	147487	2.15%	0.400%
18	10.773	1499	1505	1509	rVV5	88417	176466	2.58%	0.479%
19	10.861	1515	1520	1525	rVB2	83291	139484	2.04%	0.378%
20	10.950	1525	1535	1541	rBV	263772	468482	6.84%	1.271%
21	11.050	1542	1552	1560	rBV2	313956	736137	10.75%	1.997%
22	11.120	1560	1564	1568	rVV3	109059	165179	2.41%	0.448%
23	11.191	1571	1576	1580	rVV	216994	360288	5.26%	0.977%
24	11.244	1580	1585	1590	rVV	498260	869605	12.70%	2.359%
25	11.297	1590	1594	1598	rVB3	125159	167260	2.44%	0.454%
26	11.355	1598	1604	1607	rBV3	229361	399020	5.83%	1.082%
27	11.426	1607	1616	1628	rVB5	585971	1745379	25.49%	4.735%
28	11.526	1628	1633	1645	rVB	549623	943799	13.78%	2.560%
29	11.644	1646	1653	1662	rBV	1058070	1820174	26.58%	4.938%
30	11.744	1662	1670	1679	rBV	1240298	2247480	32.82%	6.097%
31	11.849	1679	1688	1699	rBV	3261692	6846877	100.00%	18.574%
32	11.991	1707	1712	1715	rBV6	36185	69249	1.01%	0.188%
33	12.179	1732	1744	1752	rBV2	716313	1803201	26.34%	4.892%
34	12.267	1756	1759	1764	rVB	166678	234889	3.43%	0.637%

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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 >
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35	12.349	1764	1773	1777	rBV6	193290	522787	7.64%	1.418%
36	12.479	1790	1795	1800	rBV2	156846	239397	3.50%	0.649%
37	12.632	1815	1821	1827	rBV	756868	1204638	17.59%	3.268%
38	12.720	1831	1836	1840	rBV5	57318	113530	1.66%	0.308%
39	12.820	1845	1853	1857	rVB2	200552	520962	7.61%	1.413%
40	12.885	1858	1864	1866	rBV4	117532	192765	2.82%	0.523%
41	12.914	1866	1869	1872	rVV2	72707	90805	1.33%	0.246%
42	12.961	1873	1877	1882	rVB7	89927	159017	2.32%	0.431%
43	13.191	1912	1916	1921	rBV4	54071	75641	1.10%	0.205%
44	13.273	1924	1930	1935	rBV	206623	322924	4.72%	0.876%
45	13.408	1948	1953	1954	rBV5	58959	85231	1.24%	0.231%
46	13.579	1975	1982	1991	rVB	635066	1168093	17.06%	3.169%
47	13.896	2032	2036	2042	rVB4	98195	153885	2.25%	0.417%
48	14.114	2071	2073	2078	rVB2	62628	91515	1.34%	0.248%
49	14.226	2088	2092	2098	rVB3	60007	100912	1.47%	0.274%
50	14.291	2098	2103	2109	rVB8	40086	72032	1.05%	0.195%
51	14.408	2119	2123	2126	rBV5	69814	111788	1.63%	0.303%
52	14.567	2146	2150	2155	rVV7	38549	73153	1.07%	0.198%
53	14.832	2189	2195	2200	rVV5	82007	200744	2.93%	0.545%
54	14.885	2200	2204	2209	rVB4	67041	132178	1.93%	0.359%
55	15.096	2236	2240	2245	rBV5	55823	88157	1.29%	0.239%
56	15.214	2255	2260	2266	rVB3	54662	112450	1.64%	0.305%
57	15.408	2282	2293	2304	rBV	502307	1156779	16.89%	3.138%
58	15.714	2336	2345	2350	rBV7	52877	138620	2.02%	0.376%
59	15.779	2350	2356	2363	rVB7	34883	79117	1.16%	0.215%
60	16.502	2473	2479	2491	rVB	205296	438269	6.40%	1.189%
61	16.708	2506	2514	2527	rVB	200091	523163	7.64%	1.419%

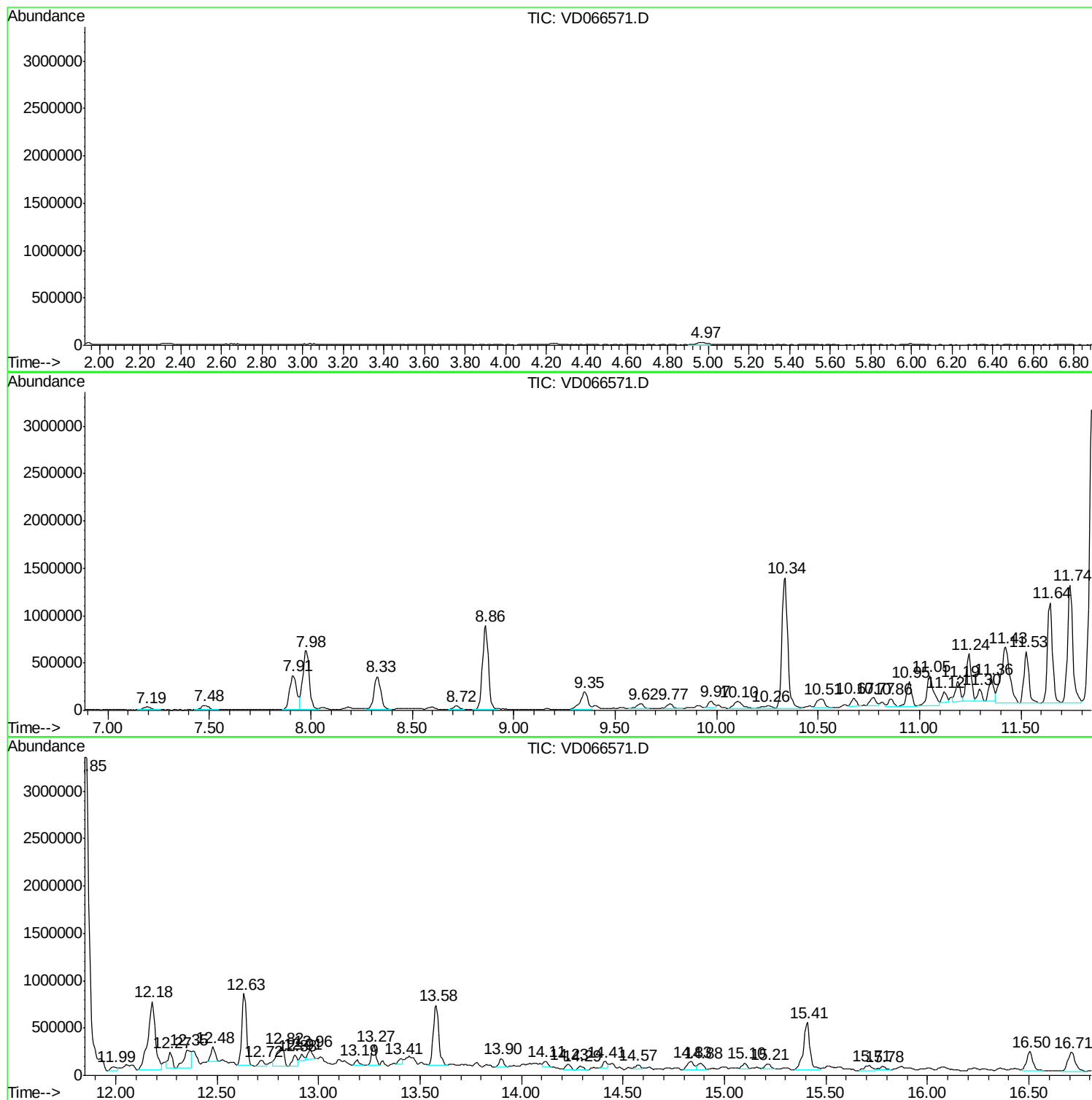
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Quant Title : SW846 8260

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



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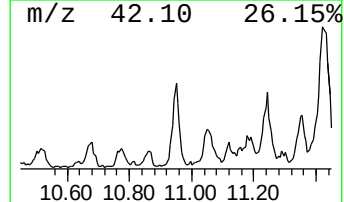
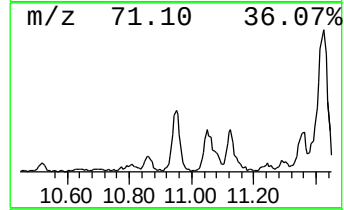
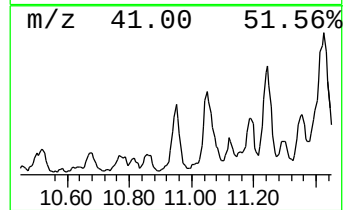
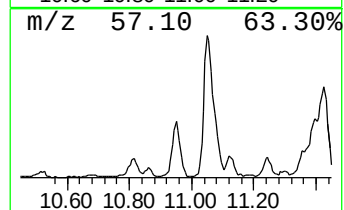
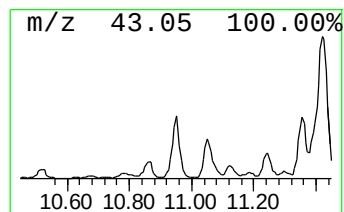
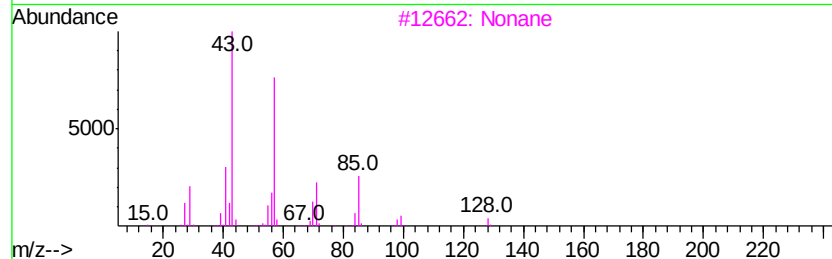
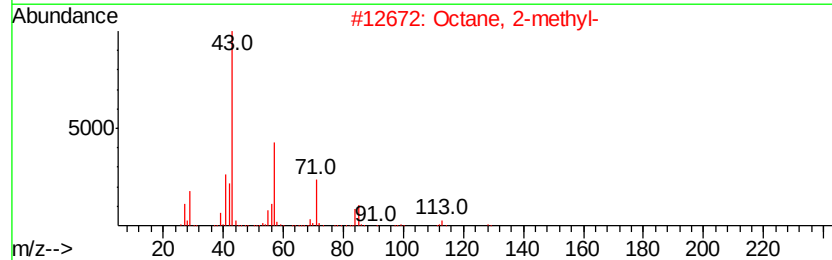
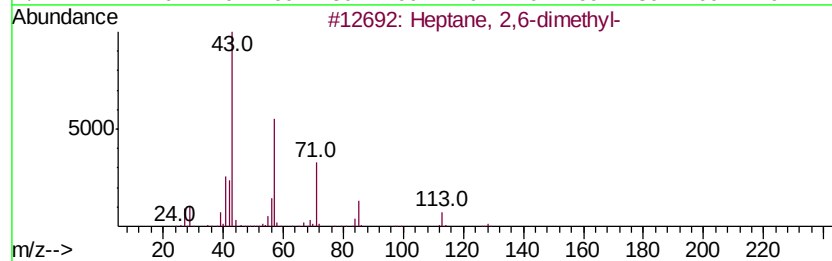
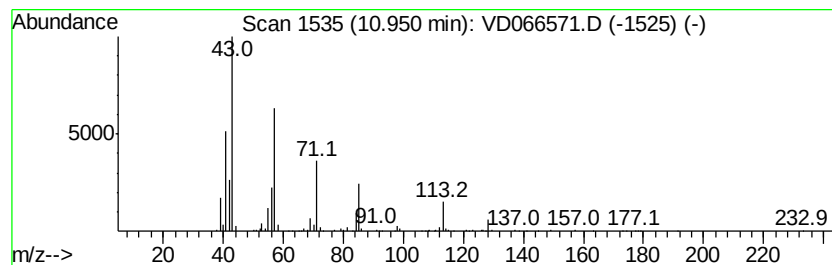
Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.95	12.87 ug/l	468482	Chlorobenzene-d5	11.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	87
2	Octane, 2-methyl-	128	C9H20	003221-61-2	59
3	Nonane	128	C9H20	000111-84-2	58
4	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	58
5	Dodecane, 5-methyl-	184	C13H28	017453-93-9	53



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ALS Vial : 19 Sample Multiplier: 1

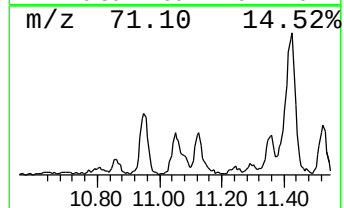
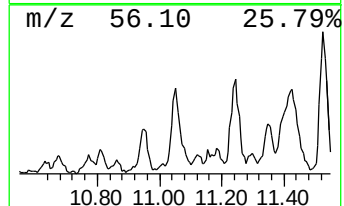
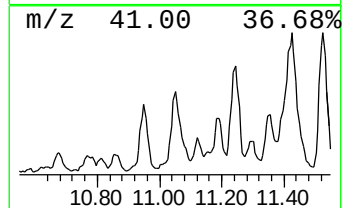
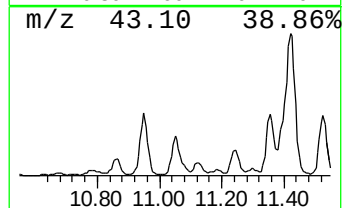
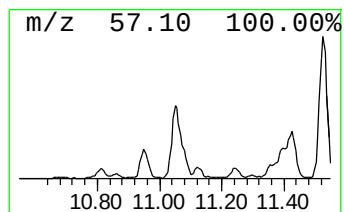
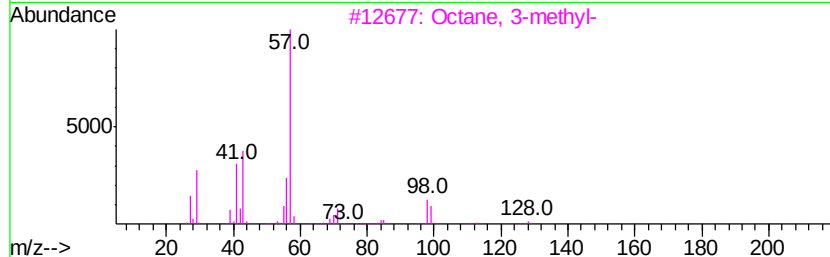
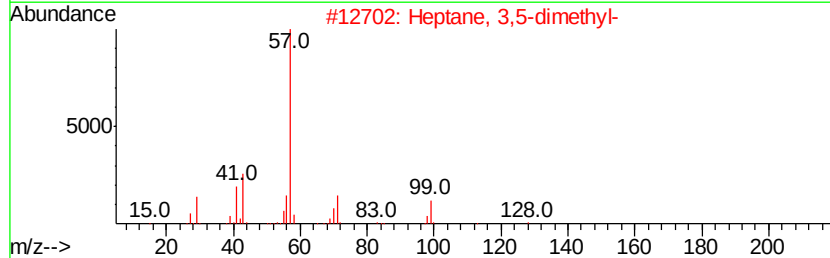
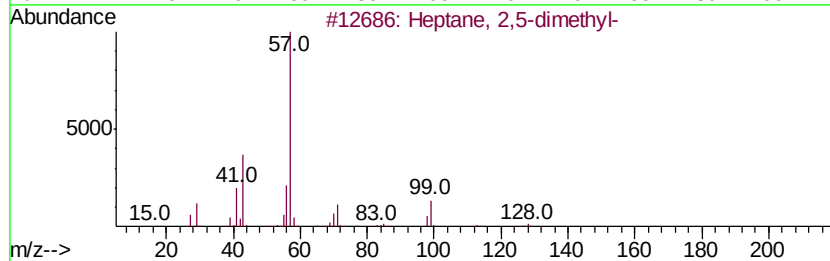
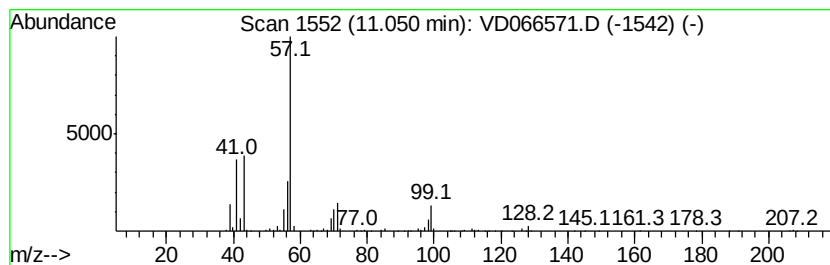
Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.05	20.22 ug/l	736137	Chlorobenzene-d5	11.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	90
2	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	72
3	Octane, 3-methyl-	128	C9H20	002216-33-3	64
4	Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	53
5	Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	53



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Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

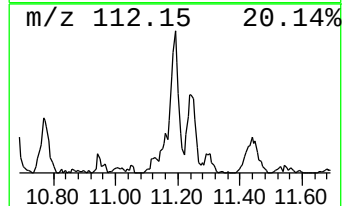
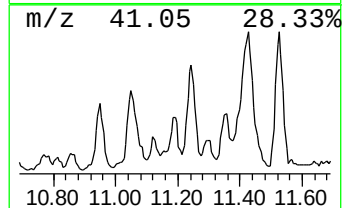
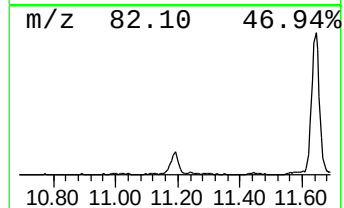
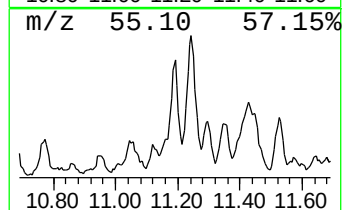
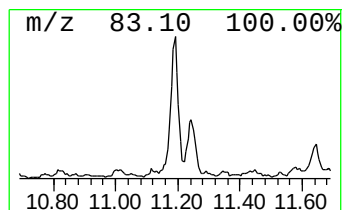
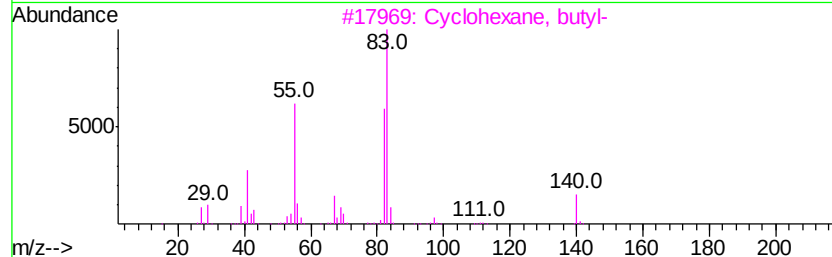
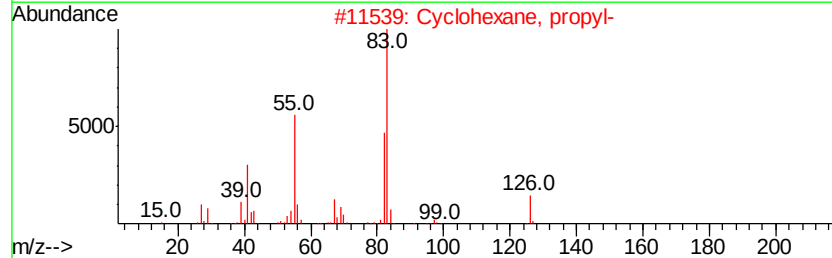
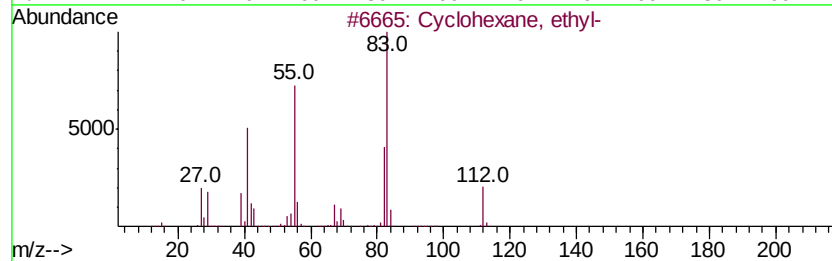
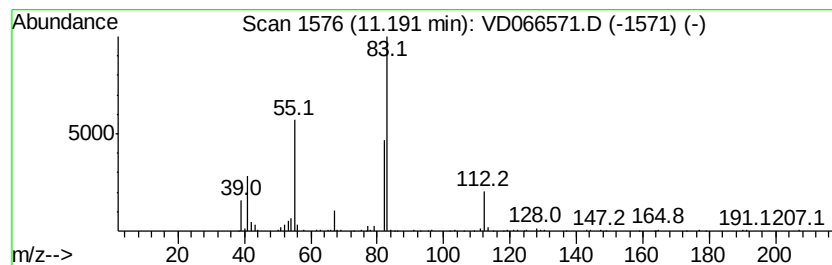
Instrument :
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ClientSampled :
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.19	9.90 ug/l	360288	Chlorobenzene-d5	11.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	90
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	72
3	Cyclohexane, butyl-	140	C10H20	001678-93-9	72
4	3-Cyclohexyl-1-chloropropane	160	C9H17Cl	001124-62-5	72
5	(E)-4-Oxohex-2-enal	112	C6H8O2	1000374-04-2	53



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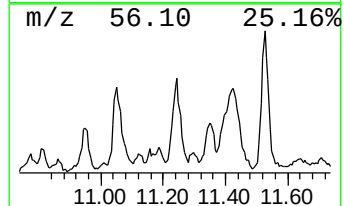
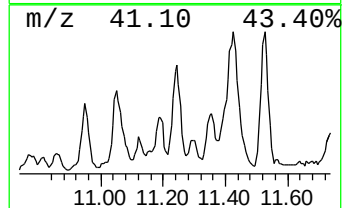
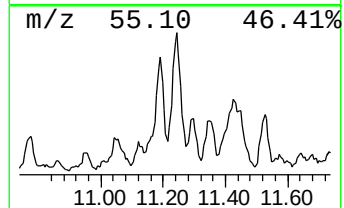
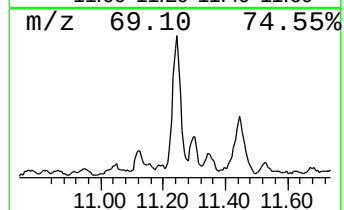
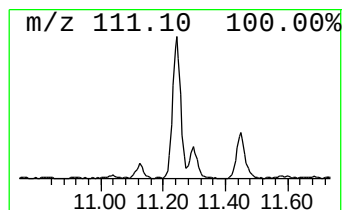
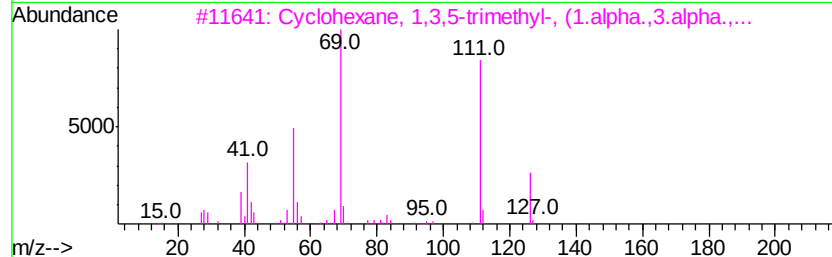
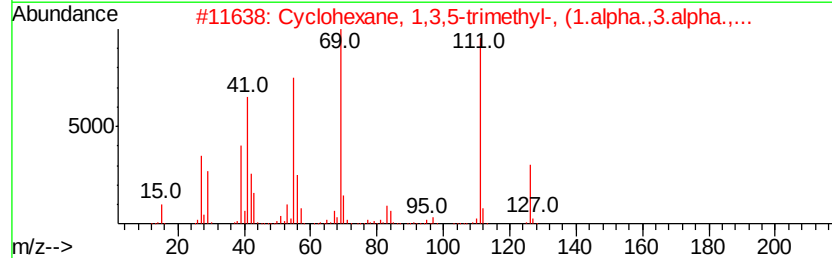
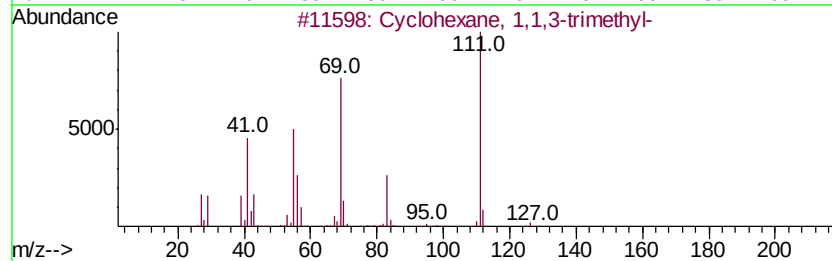
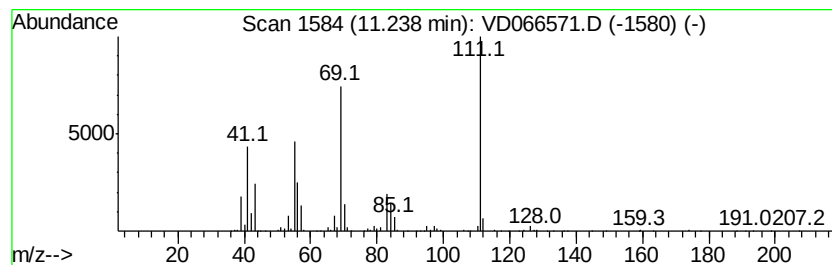
Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.24	23.89 ug/l	869605	Chlorobenzene-d5	11.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	91
2		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-26-2	72
3		Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-27-3	64
4		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	59
5		Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	59



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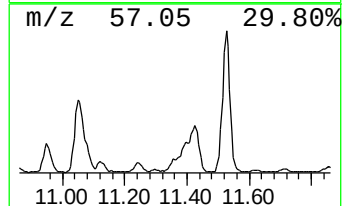
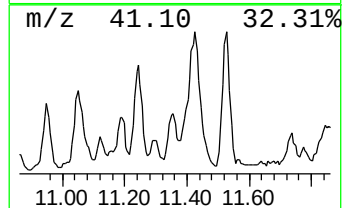
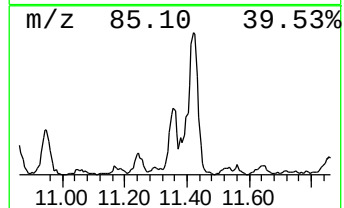
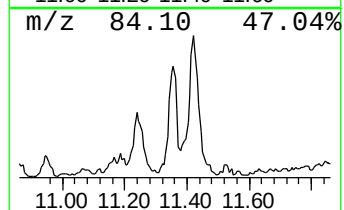
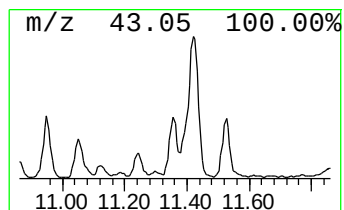
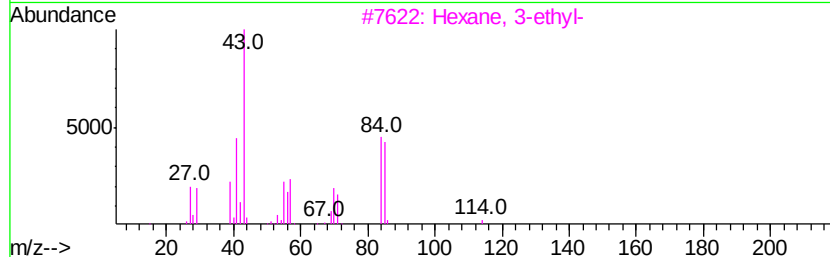
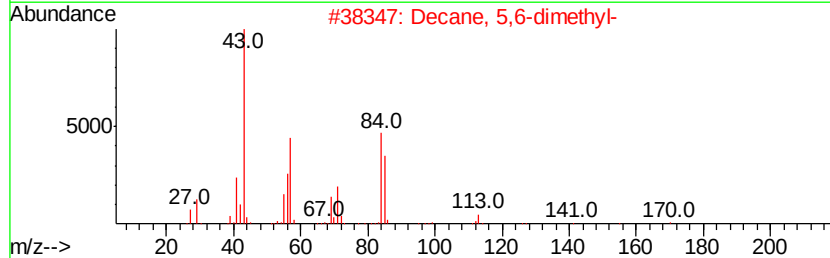
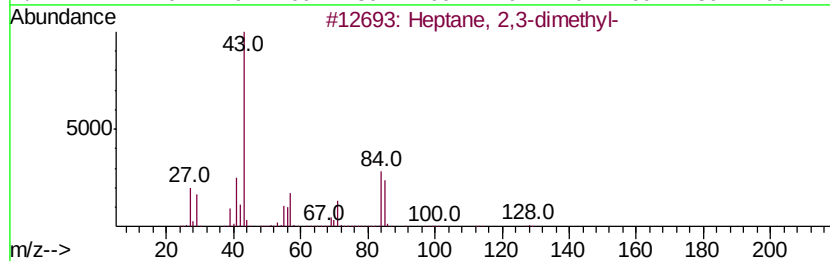
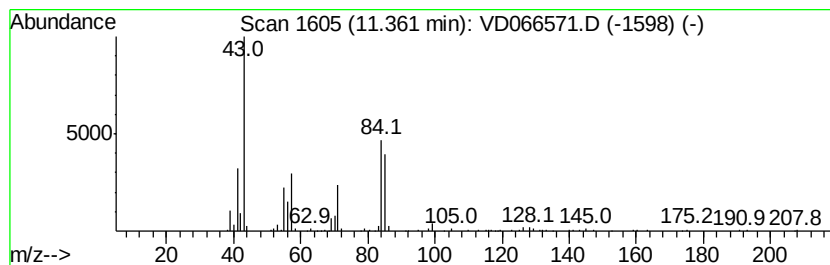
Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 5 Heptane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.36	10.96 ug/l	399020	Chlorobenzene-d5	11.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	81	
2	Decane, 5,6-dimethyl-	170	C12H26	001636-43-7	72	
3	Hexane, 3-ethyl-	114	C8H18	000619-99-8	64	
4	Octane, 4,5-diethyl-	170	C12H26	001636-41-5	64	
5	Hexane, 3-ethyl-2-methyl-	128	C9H20	016789-46-1	53	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066571.D
Acq On : 04 Sep 2020 18:57
Operator : VA/SY
Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

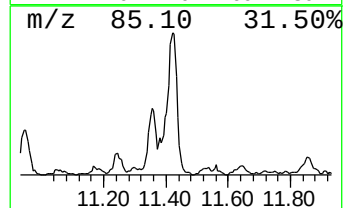
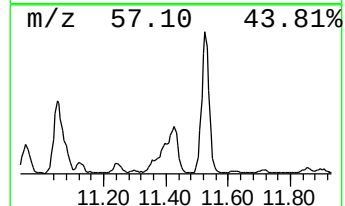
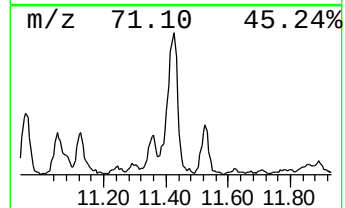
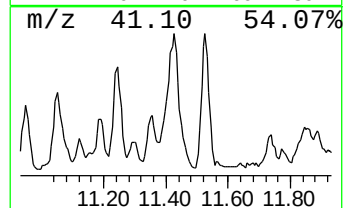
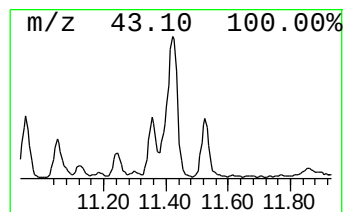
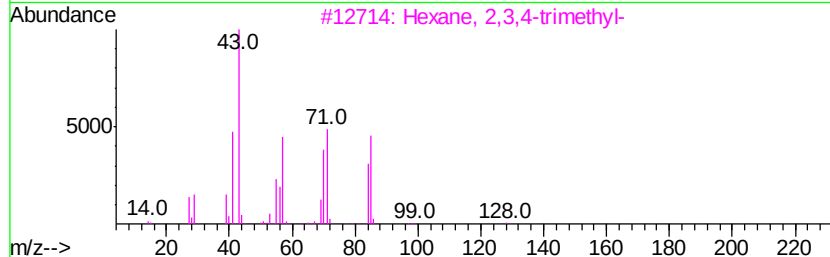
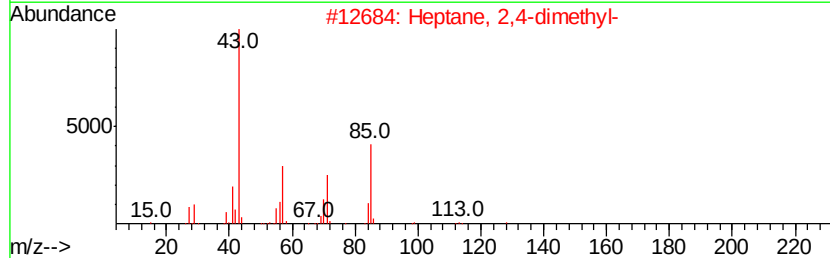
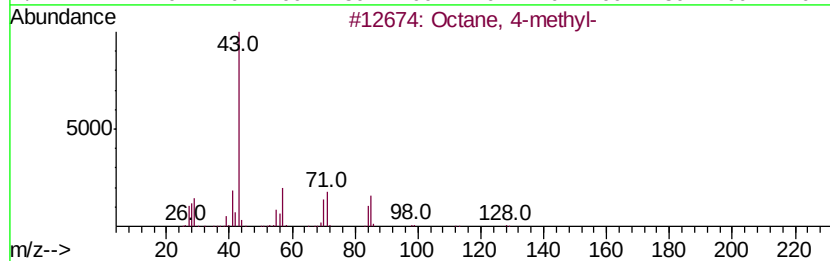
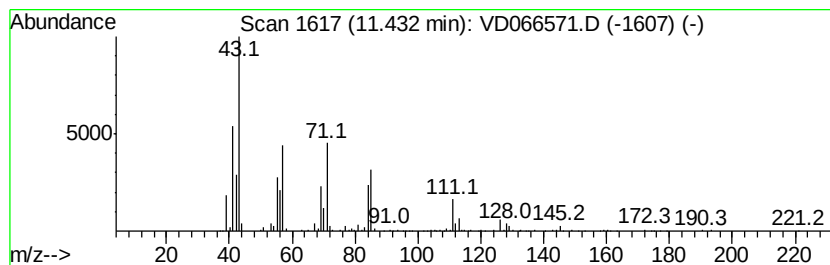
Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 6 Octane, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.43	47.95 ug/l	1745380	Chlorobenzene-d5	11.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 4-methyl-	128	C9H20	002216-34-4	59
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	53
3		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	50
4		Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	50
5		Ether, hexyl pentyl	172	C11H24O	032357-83-8	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066571.D
Acq On : 04 Sep 2020 18:57
Operator : VA/SY
Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-1

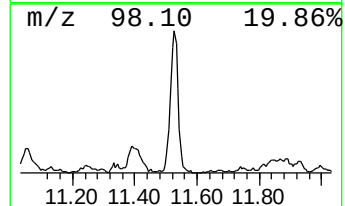
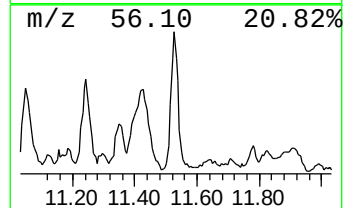
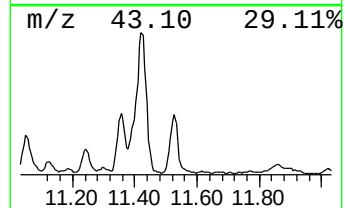
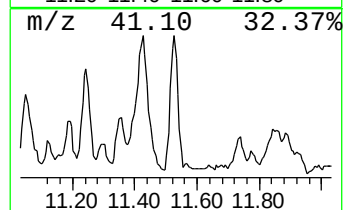
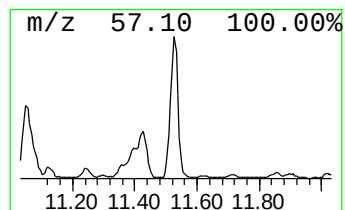
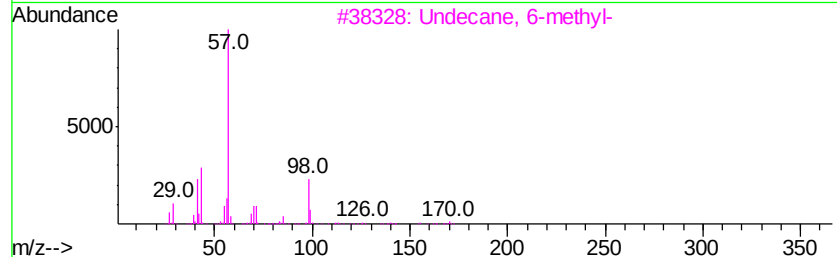
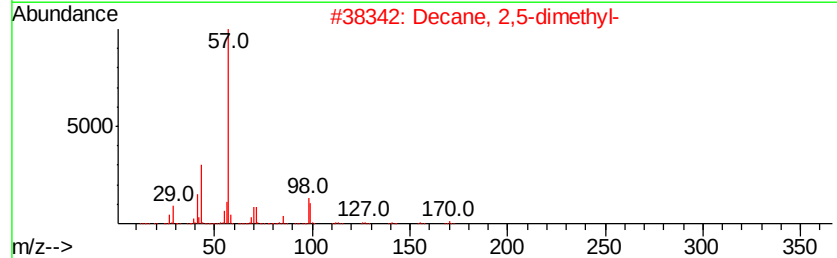
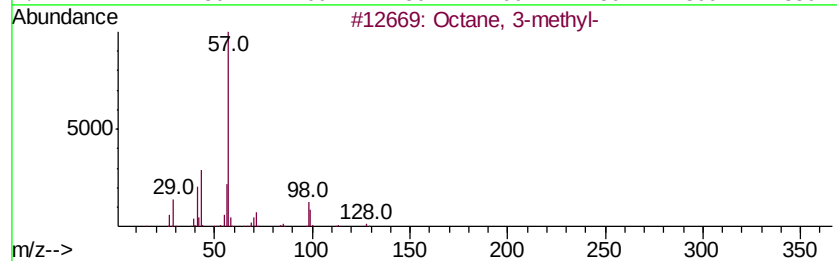
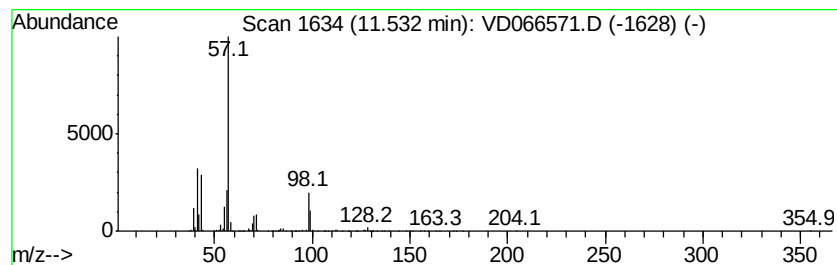
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Quant Title : SW846 8260

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

Peak Number 7 Octane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	25.93 ug/l	943799	Chlorobenzene-d5	11.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 3-methyl-	128	C9H20	002216-33-3	72
2		Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	59
3		Undecane, 6-methyl-	170	C12H26	017302-33-9	59
4		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	59
5		Heptane, 3-ethyl-	128	C9H20	015869-80-4	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066571.D
Acq On : 04 Sep 2020 18:57
Operator : VA/SY
Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

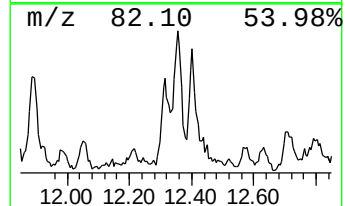
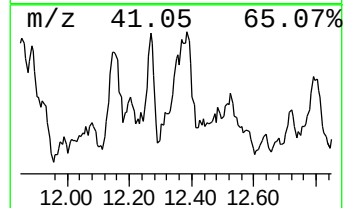
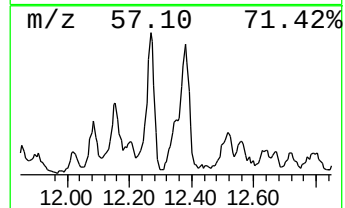
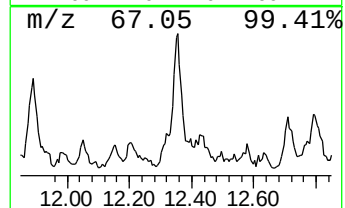
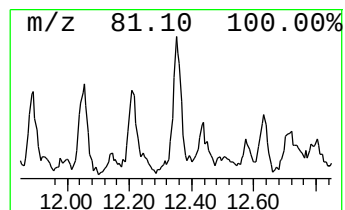
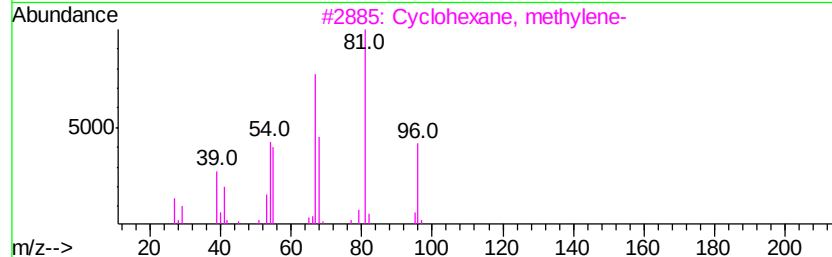
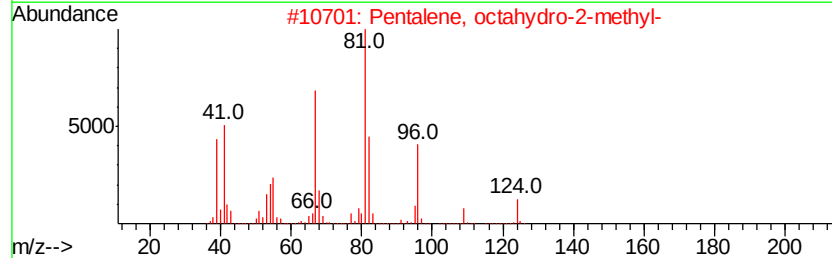
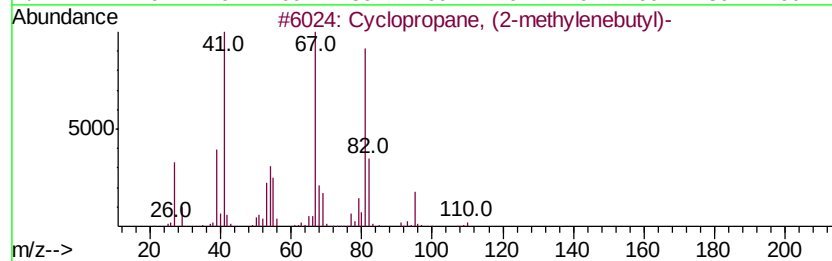
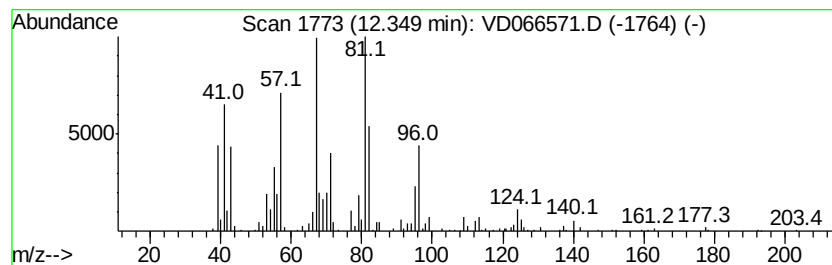
Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 8 Cyclopropane, (2-methyleneb... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.35	14.36 ug/l	522787	Chlorobenzene-d5	11.64		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, (2-methylenebutyl)-	110	C8H14	074685-56-6	64
2		Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	64
3		Cyclohexane, methylene-	96	C7H12	001192-37-6	47
4		Cyclohexene, 1-propyl-	124	C9H16	002539-75-5	43
5		Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066571.D
Acq On : 04 Sep 2020 18:57
Operator : VA/SY
Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-1

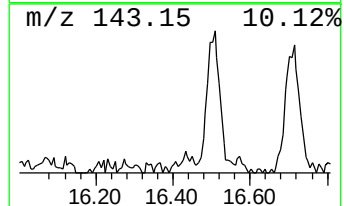
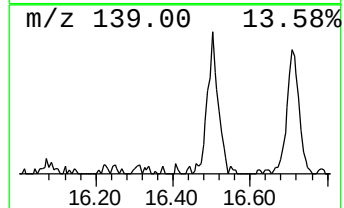
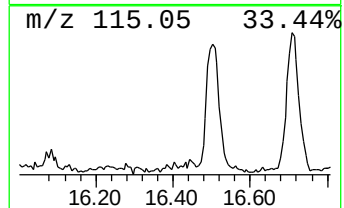
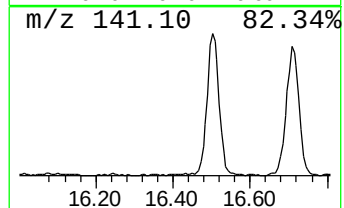
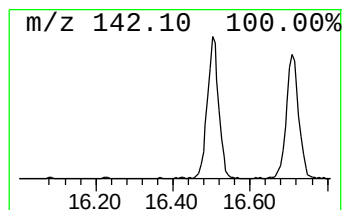
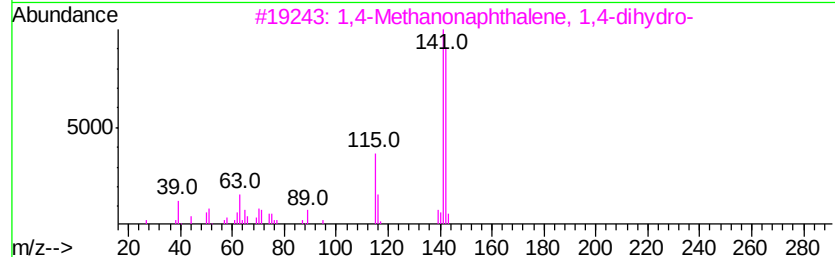
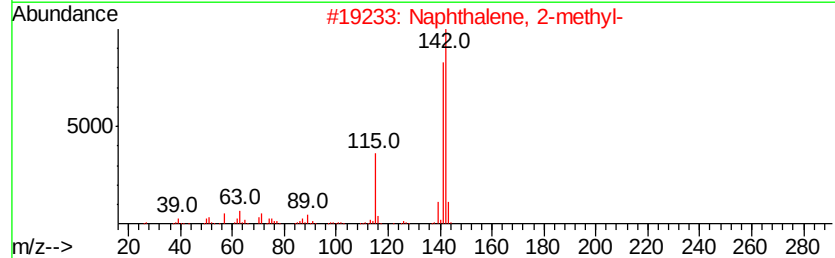
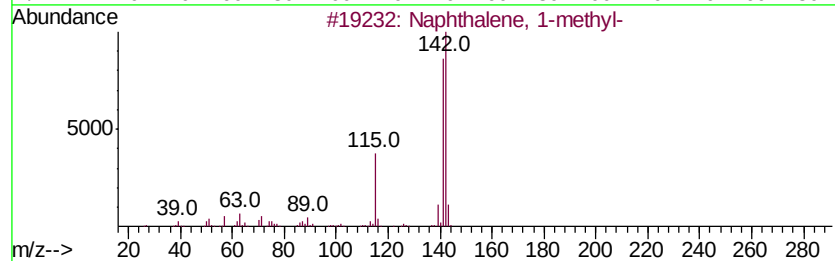
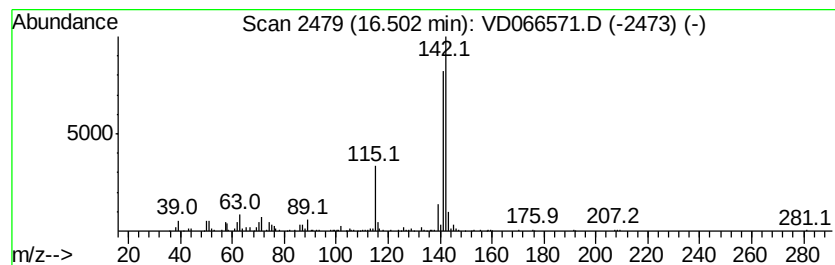
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	18.76 ug/l	438269	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
3		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59
5		Benzocycloheptatriene	142	C11H10	000264-09-5	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066571.D
Acq On : 04 Sep 2020 18:57
Operator : VA/SY
Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-1

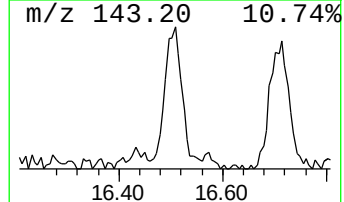
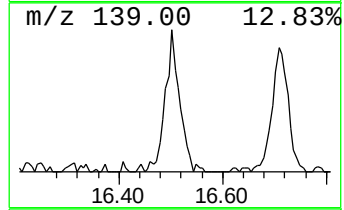
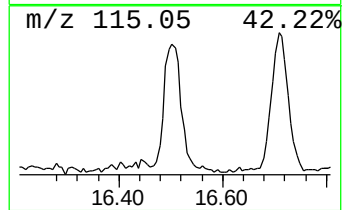
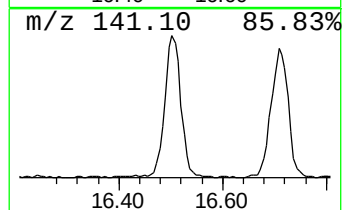
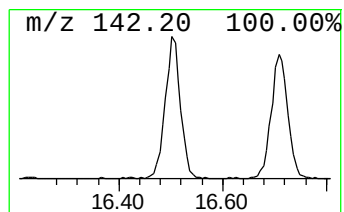
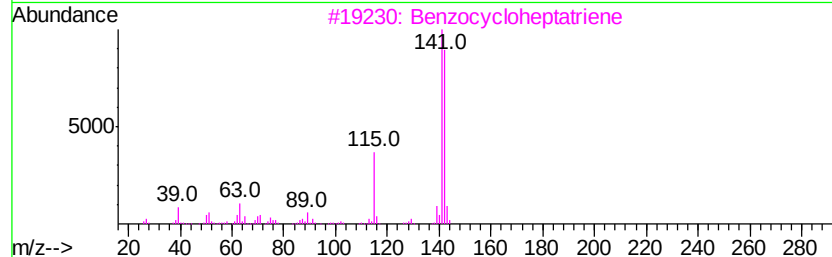
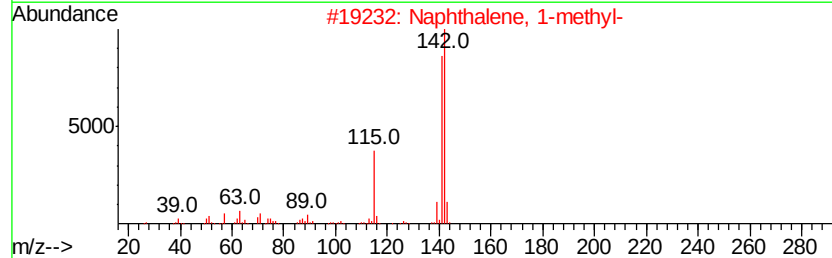
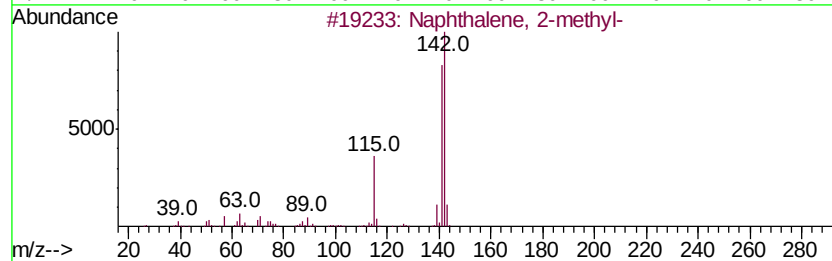
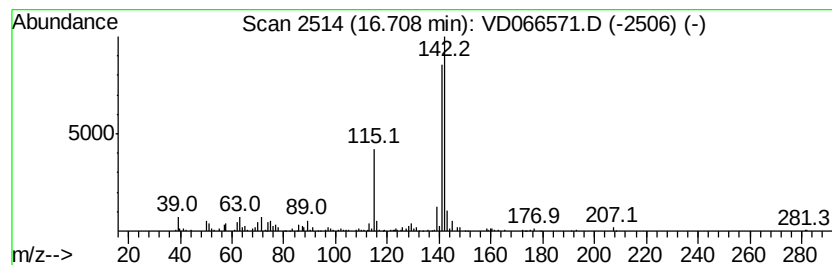
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.71	22.39 ug/l	523163	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5		1-Naphthaleneacetamide	185	C12H11NO	000086-86-2	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD090420\
Data File : VD066571.D
Acq On : 04 Sep 2020 18:57
Operator : VA/SY
Sample : L3896-01
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
WC-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.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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Heptane, 2,5-dime...	11.05	20.2	ug/l	736137	3	11.64	1820170	50.0
Cyclohexane, ethyl-	11.19	9.9	ug/l	360288	3	11.64	1820170	50.0
Cyclohexane, 1,1,...	11.24	23.9	ug/l	869605	3	11.64	1820170	50.0
Heptane, 2,3-dime...	11.36	11.0	ug/l	399020	3	11.64	1820170	50.0
Octane, 4-methyl-	11.43	48.0	ug/l	1745380	3	11.64	1820170	50.0
Octane, 3-methyl-	11.53	25.9	ug/l	943799	3	11.64	1820170	50.0
Cyclopropane, (2-...	12.35	14.4	ug/l	522787	3	11.64	1820170	50.0
Naphthalene, 1-me...	16.50	18.8	ug/l	438269	4	13.58	1168090	50.0
Benzocycloheptatr...	16.71	22.4	ug/l	523163	4	13.58	1168090	50.0