

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052319\
 Data File : VX009863.D
 Acq On : 23 May 2019 17:04
 Operator : JC/SP
 Sample : K2976-02 10X
 Misc : 5.76g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SB-E2-3R(17-20)

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_X\METHOD\82X042919W.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	5.501	701	713	725	rBV	95727	279032	2.78%	0.219%
2	5.659	727	739	757	rVB2	177381	507622	5.07%	0.398%
3	6.062	794	805	819	rBV2	120715	317801	3.17%	0.249%
4	6.860	924	936	951	rBV	273488	664068	6.63%	0.521%
5	8.665	1224	1232	1235	rBV	92570	167196	1.67%	0.131%
6	8.714	1235	1240	1251	rVB	620344	1002639	10.01%	0.786%
7	9.043	1288	1294	1302	rVB	104746	178365	1.78%	0.140%
8	9.348	1338	1344	1350	rVB	60363	100965	1.01%	0.079%
9	9.445	1350	1360	1369	rBV	207530	339454	3.39%	0.266%
10	9.561	1373	1379	1385	rBV2	216632	464200	4.63%	0.364%
11	9.622	1385	1389	1393	rVB2	227083	318581	3.18%	0.250%
12	9.677	1393	1398	1402	rBV	464186	696401	6.95%	0.546%
13	9.817	1414	1421	1425	rBV2	87399	133559	1.33%	0.105%
14	9.890	1425	1433	1437	rVV3	592312	1413453	14.11%	1.108%
15	9.957	1440	1444	1449	rVB	552168	738427	7.37%	0.579%
16	10.061	1455	1461	1465	rBV	345506	531665	5.31%	0.417%
17	10.110	1465	1469	1474	rBV	604576	806197	8.05%	0.632%
18	10.177	1474	1480	1485	rVB4	141541	336878	3.36%	0.264%
19	10.238	1485	1490	1496	rBV2	372801	680495	6.79%	0.533%
20	10.335	1496	1506	1509	rBV2	823479	1721771	17.18%	1.350%
21	10.378	1509	1513	1516	rVV	1585888	2064310	20.60%	1.618%
22	10.408	1516	1518	1530	rVB	1142358	1932914	19.29%	1.515%
23	10.512	1530	1535	1540	rBV	645590	957907	9.56%	0.751%
24	10.585	1540	1547	1550	rBV2	306812	507606	5.07%	0.398%
25	10.640	1550	1556	1564	rBV3	2900387	5128974	51.19%	4.021%
26	10.725	1564	1570	1574	rBV2	1613505	2579726	25.74%	2.022%
27	10.768	1574	1577	1580	rVB2	848374	1033892	10.32%	0.811%
28	10.835	1580	1588	1592	rBV2	6428967	10020343	100.00%	7.855%
29	10.914	1592	1601	1604	rBV3	5866703	9816377	97.96%	7.696%
30	10.951	1604	1607	1612	rVB	3086025	4094501	40.86%	3.210%
31	11.006	1612	1616	1621	rVB3	959441	1680102	16.77%	1.317%
32	11.061	1621	1625	1626	rBV2	1234786	1638937	16.36%	1.285%
33	11.085	1626	1629	1633	rVB2	1887356	2314687	23.10%	1.815%
34	11.140	1633	1638	1642	rVB3	2371515	3882628	38.75%	3.044%

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35	11.183	1642	1645	1648	rBV	888171	1100631	10.98%	0.863%
36	11.219	1648	1651	1653	rBV	639985	710492	7.09%	0.557%
37	11.274	1653	1660	1669	rVB3	3569094	6560214	65.47%	5.143%
38	11.359	1670	1674	1676	rBV	885361	1044995	10.43%	0.819%
39	11.396	1676	1680	1684	rBV3	1197264	1898695	18.95%	1.488%
40	11.445	1684	1688	1691	rBV2	1556843	2063575	20.59%	1.618%
41	11.487	1691	1695	1699	rVB	3479592	4126295	41.18%	3.235%
42	11.536	1699	1703	1707	rVB3	2087764	3356272	33.49%	2.631%
43	11.579	1707	1710	1715	rVB4	552687	756868	7.55%	0.593%
44	11.628	1715	1718	1720	rBV	538677	580332	5.79%	0.455%
45	11.683	1720	1727	1732	rVB5	1717425	3431102	34.24%	2.690%
46	11.737	1732	1736	1738	rBV2	1280793	1532661	15.30%	1.202%
47	11.768	1738	1741	1744	rBV	4348777	4704843	46.95%	3.688%
48	11.804	1744	1747	1749	rBV2	640237	748066	7.47%	0.586%
49	11.890	1757	1761	1763	rBV	1022565	1387232	13.84%	1.088%
50	11.920	1763	1766	1767	rVV	1267817	1626961	16.24%	1.275%
51	11.945	1767	1770	1775	rVV2	2417870	3986335	39.78%	3.125%
52	12.000	1775	1779	1782	rVV	3890636	5041013	50.31%	3.952%
53	12.030	1782	1784	1789	rVV3	825920	1252087	12.50%	0.982%
54	12.091	1789	1794	1797	rVV3	1065480	1909081	19.05%	1.497%
55	12.128	1797	1800	1803	rVB2	711453	796240	7.95%	0.624%
56	12.164	1803	1806	1808	rBV2	250288	303871	3.03%	0.238%
57	12.195	1808	1811	1812	rBV4	165135	157163	1.57%	0.123%
58	12.298	1824	1828	1833	rBV2	1211285	1759825	17.56%	1.380%
59	12.365	1835	1839	1846	rVV2	2294204	4172015	41.64%	3.271%
60	12.475	1851	1857	1862	rVV3	723685	1748551	17.45%	1.371%
61	12.524	1862	1865	1870	rVB4	761933	1282290	12.80%	1.005%
62	12.609	1873	1879	1881	rBV4	563241	996319	9.94%	0.781%
63	12.664	1885	1888	1892	rVB	930516	1052337	10.50%	0.825%
64	12.762	1897	1904	1910	rVV6	522039	1566228	15.63%	1.228%
65	12.816	1910	1913	1918	rVV4	225644	469250	4.68%	0.368%
66	12.871	1918	1922	1929	rVB4	775963	1323256	13.21%	1.037%
67	12.944	1929	1934	1937	rBV3	689618	1182454	11.80%	0.927%
68	12.975	1937	1939	1943	rVV2	592268	670773	6.69%	0.526%
69	13.042	1946	1950	1953	rVV2	451079	600055	5.99%	0.470%
70	13.079	1953	1956	1963	rVB2	219510	407725	4.07%	0.320%
71	13.146	1963	1967	1972	rBV2	169366	302844	3.02%	0.237%

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Title : SW846 8260

72	13.341	1995	1999	2006	rVB2	651923	973071	9.71%	0.763%
73	13.426	2010	2013	2016	rVV2	96879	149918	1.50%	0.118%
74	13.469	2017	2020	2031	rVB7	161451	411256	4.10%	0.322%
75	13.639	2044	2048	2051	rBV2	97631	142918	1.43%	0.112%
76	13.688	2051	2056	2059	rVV5	43702	103136	1.03%	0.081%
77	13.725	2059	2062	2069	rVB4	65220	117671	1.17%	0.092%

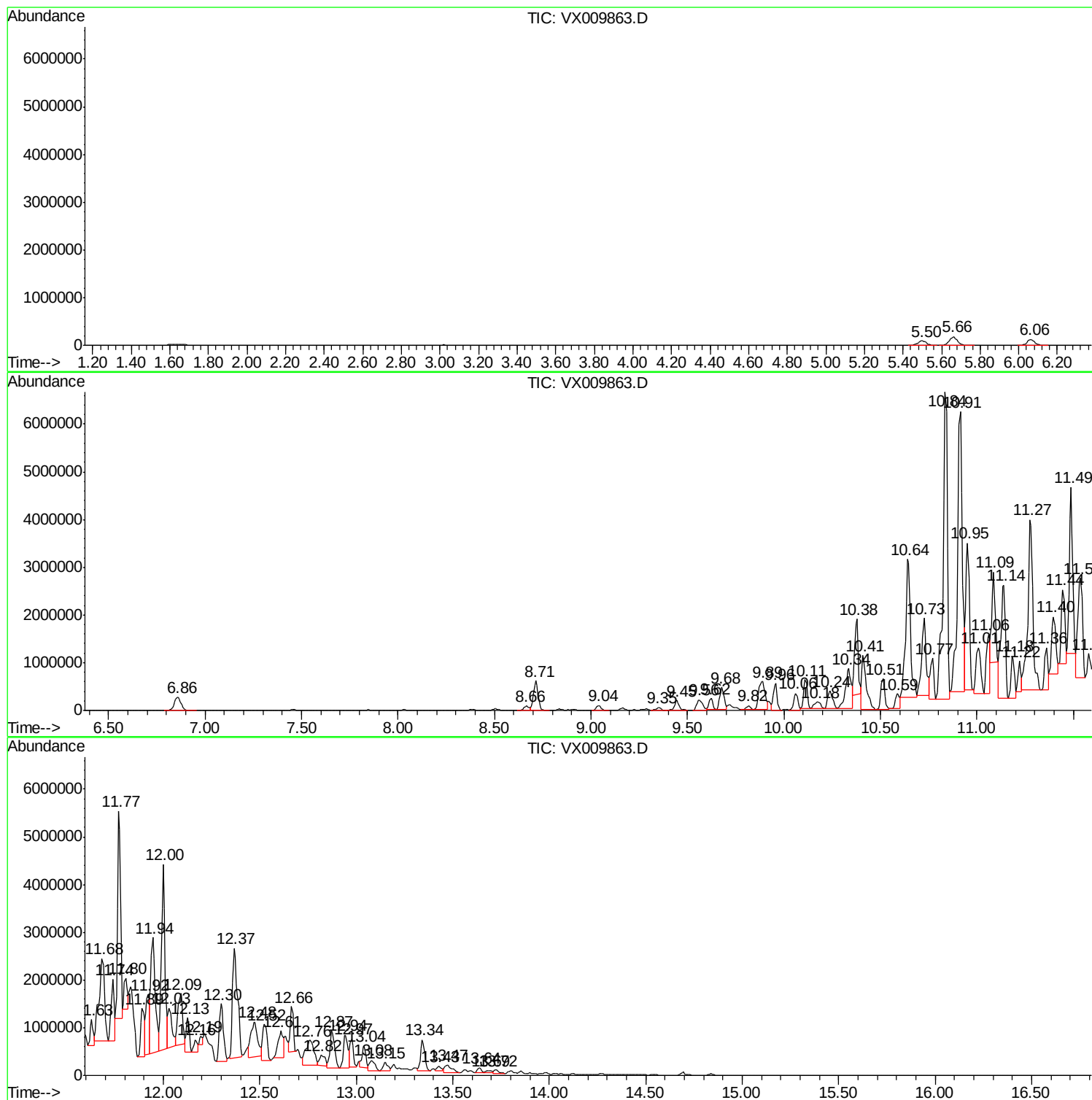
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



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Instrument :
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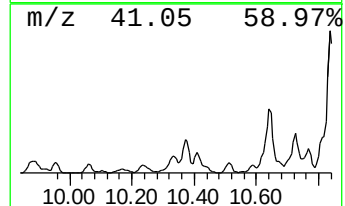
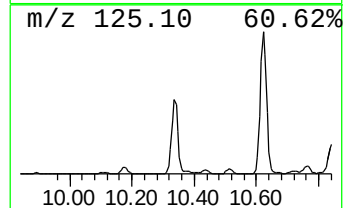
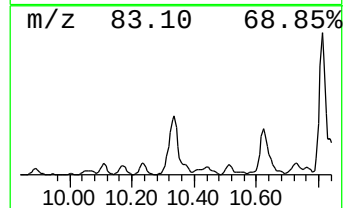
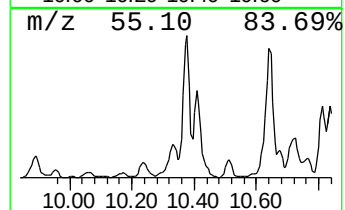
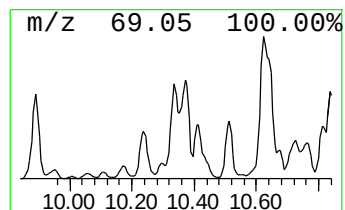
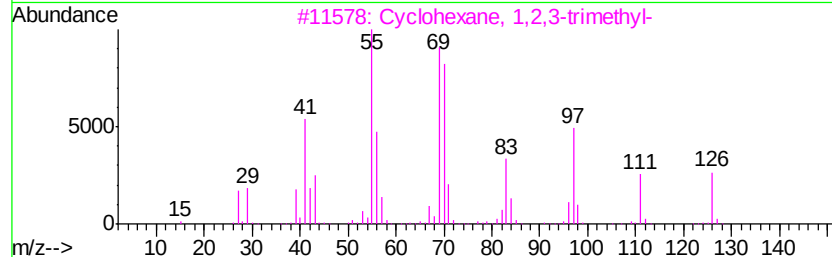
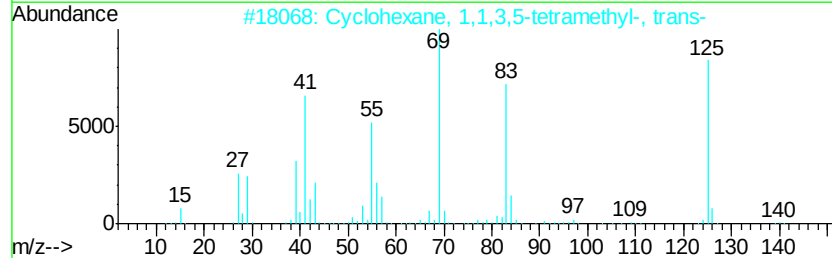
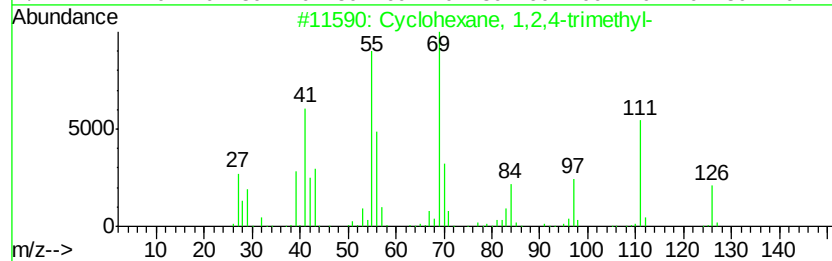
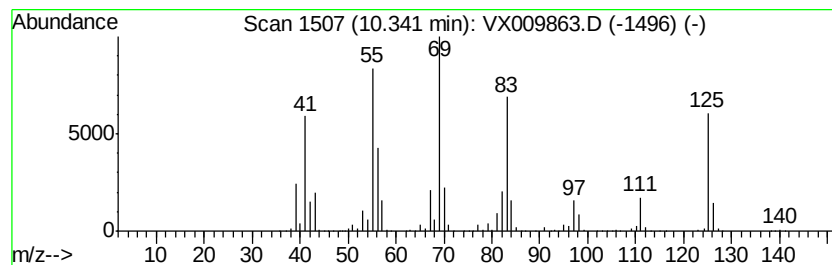
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1,2,4-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.34	106.78 ug/l	1721770	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	70
2		Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	49
3		Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	45
4		Cyclopentane, 1-hexyl-3-methyl-	168	C12H24	061142-68-5	43
5		2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	38



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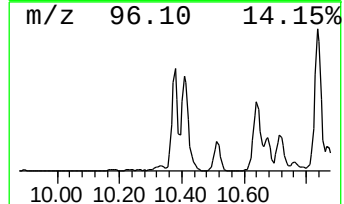
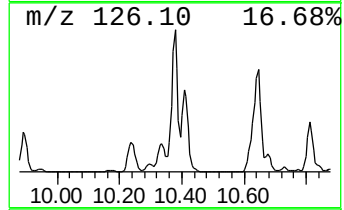
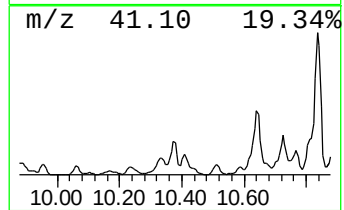
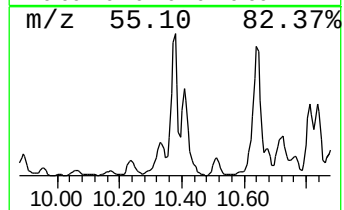
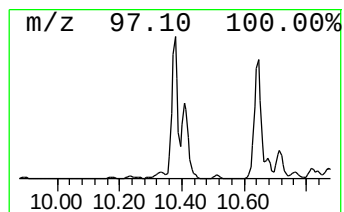
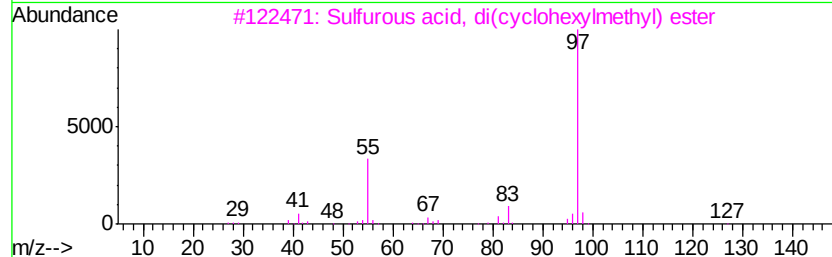
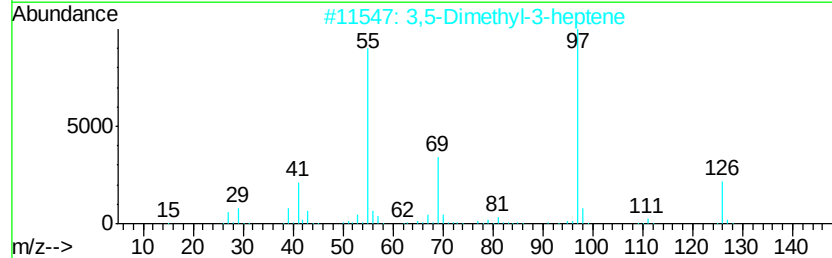
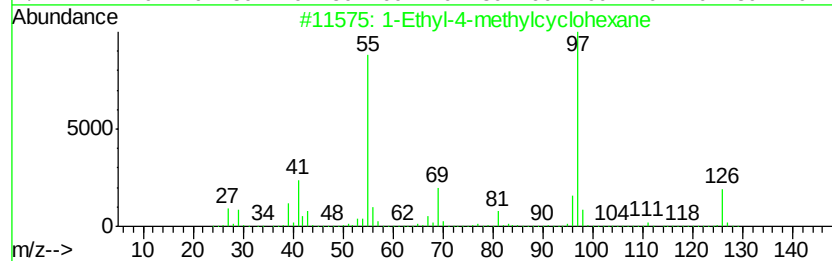
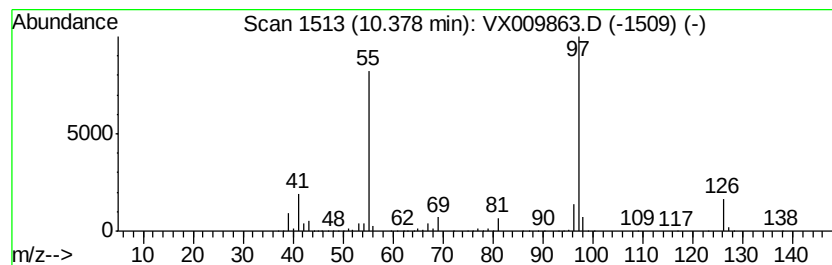
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Ethyl-4-methylcyclohexane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.38	128.03 ug/l	2064310	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	91
2		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	72
3		Sulfurous acid, di(cyclohexylmet...	274	C14H26O3S	1000309-22-7	72
4		Sulfurous acid, cyclohexylmethyl...	262	C13H26O3S	1000309-21-5	72
5		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	64



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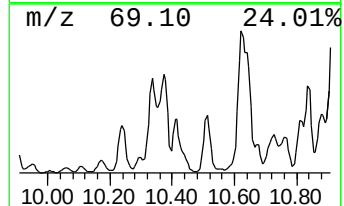
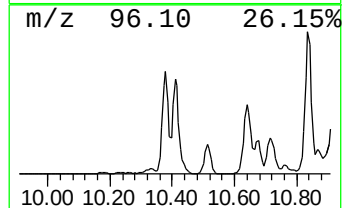
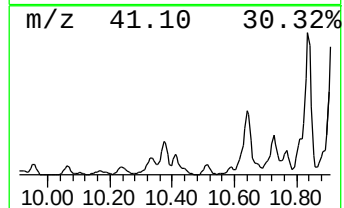
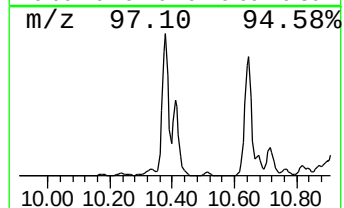
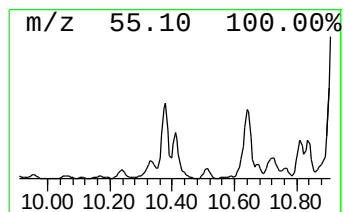
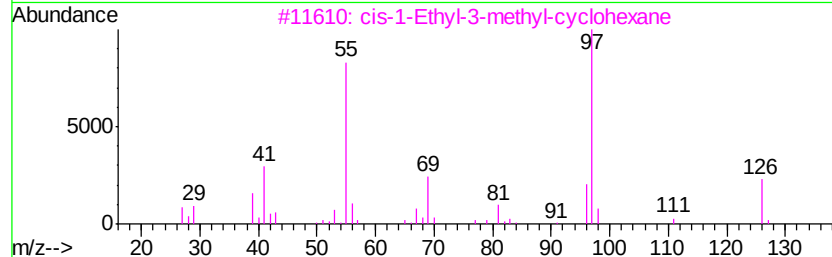
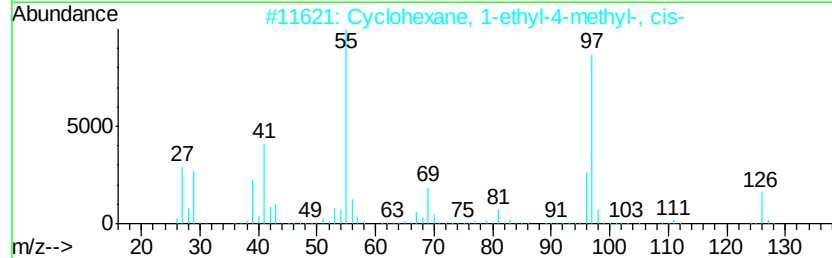
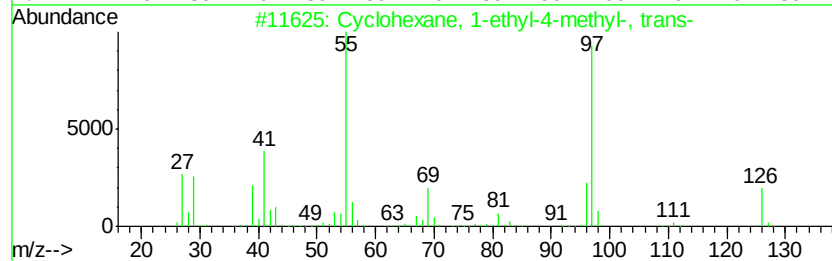
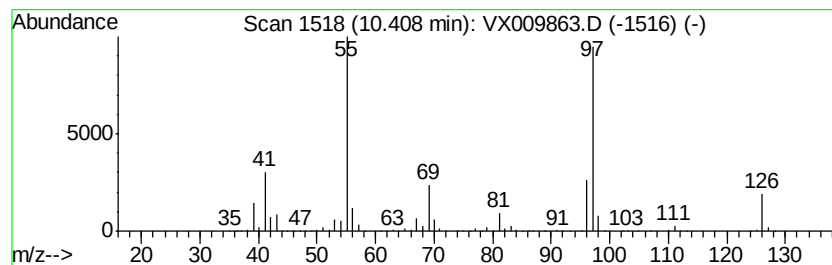
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1-ethyl-4-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.41	119.88 ug/l	1932910	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	94
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	94
3		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	93
4		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	90
5		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	90



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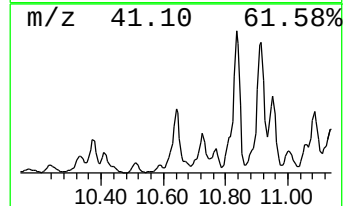
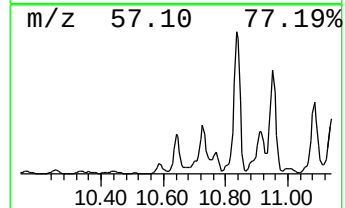
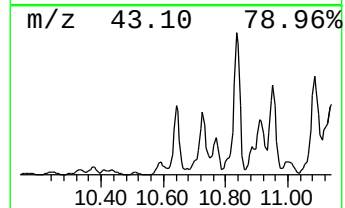
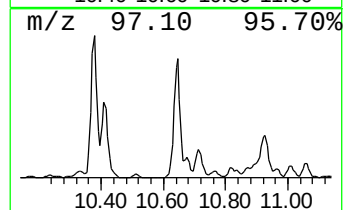
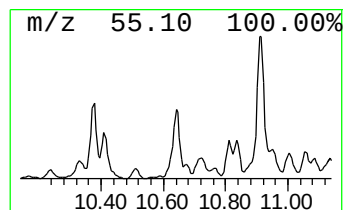
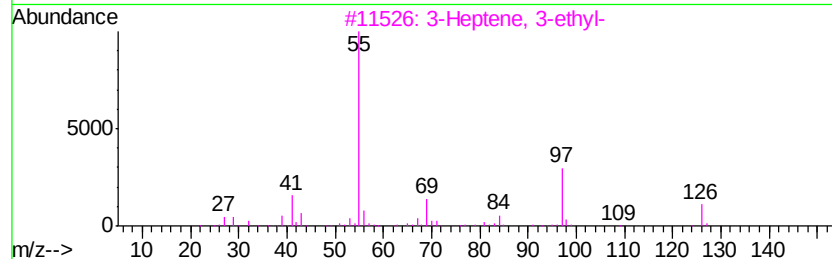
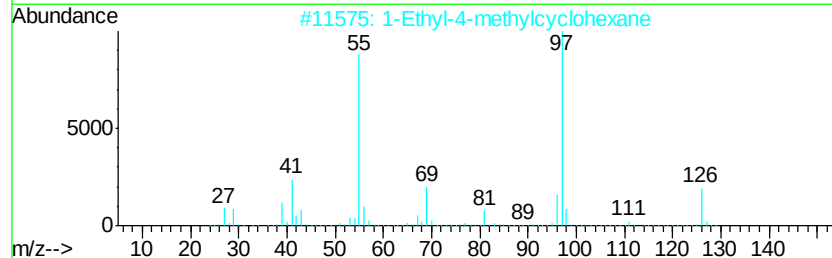
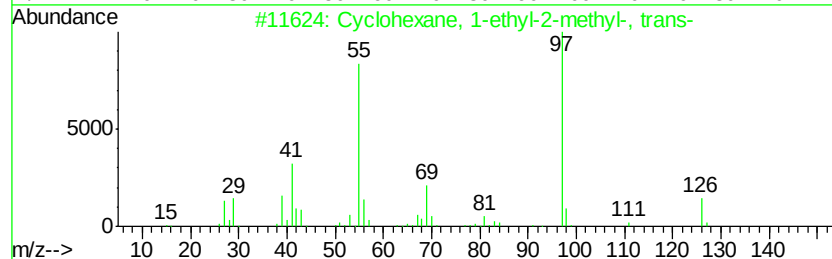
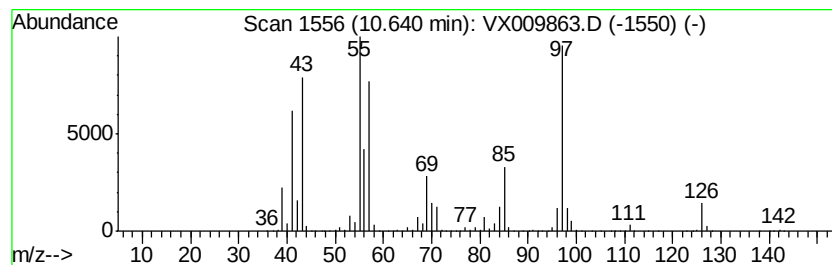
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.64	318.10 ug/l	5128970	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	55
2		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	49
3		3-Heptene, 3-ethyl-	126	C9H18	074764-46-8	49
4		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	49
5		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

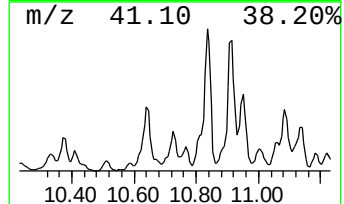
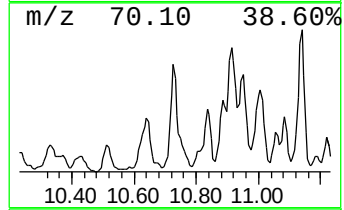
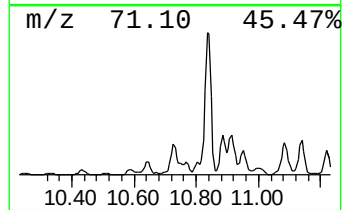
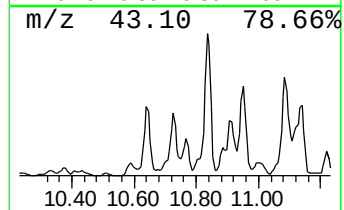
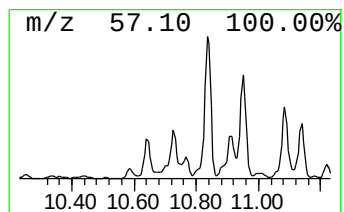
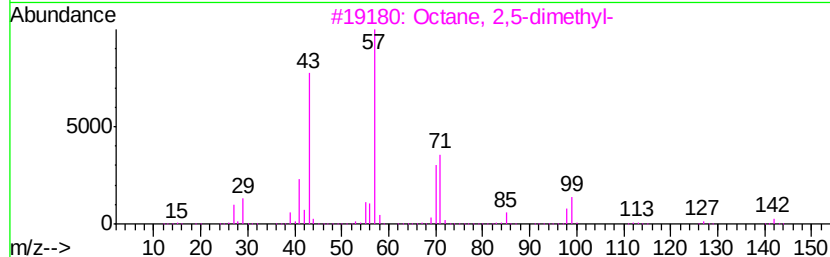
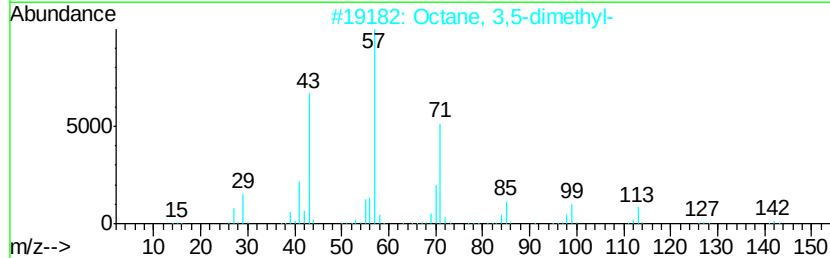
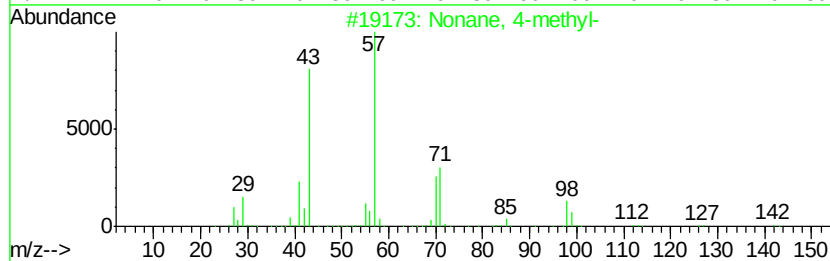
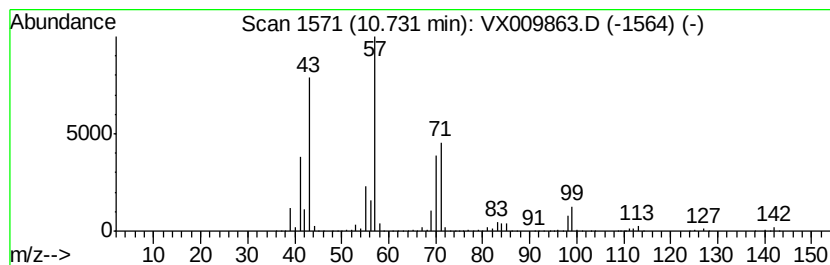
Instrument :
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ClientSampleId :
EP1-SB-E2-3R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 5 Nonane, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	159.99 ug/l	2579730	Chlorobenzene-d5	10.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Nonane, 4-methyl-	142 C10H22	017301-94-9 87
2		Octane, 3,5-dimethyl-	142 C10H22	015869-93-9 80
3		Octane, 2,5-dimethyl-	142 C10H22	015869-89-3 74
4		Heptane, 2,3,4-trimethyl-	142 C10H22	052896-95-4 72
5		Heptadecane, 2,6-dimethyl-	268 C19H40	054105-67-8 53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

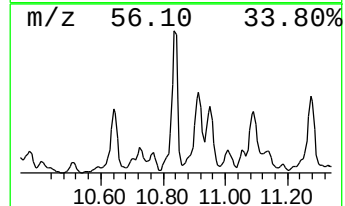
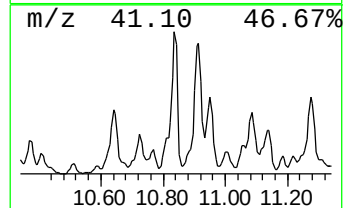
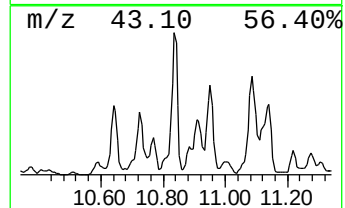
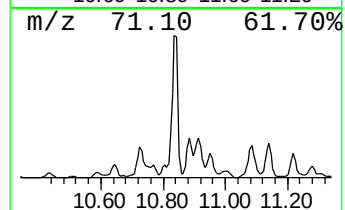
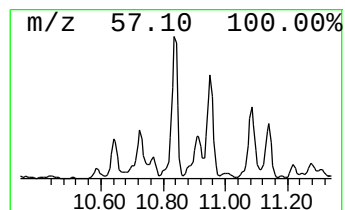
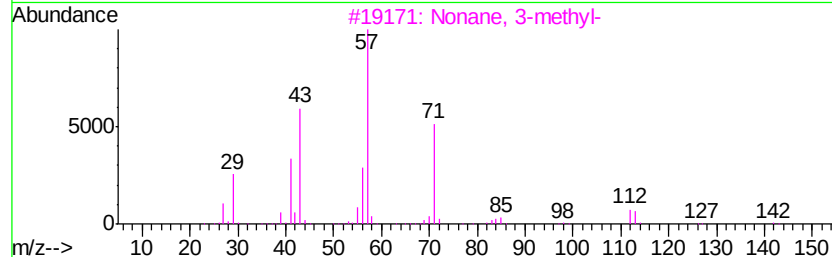
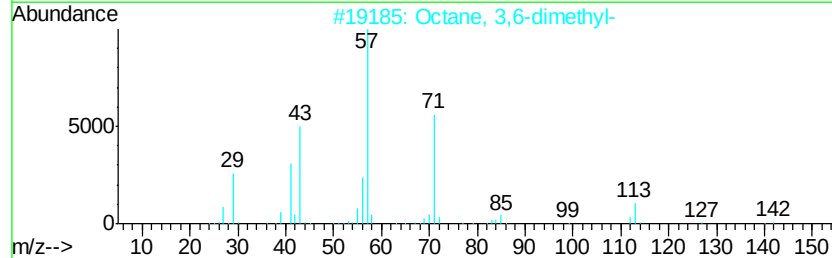
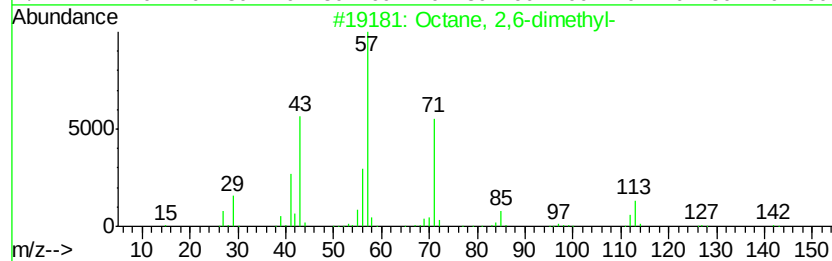
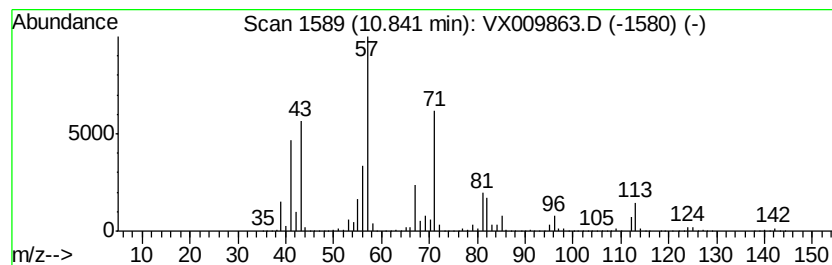
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.84	621.46 ug/l	10020300	Chlorobenzene-d5	10.11	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	93
2	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	76
3	Nonane, 3-methyl-	142	C10H22	005911-04-6	70
4	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	47
5	Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(17-20)

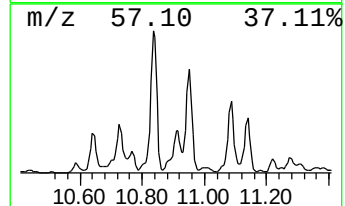
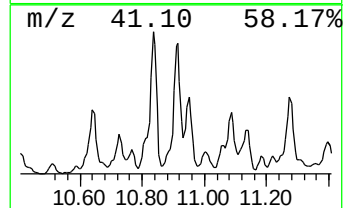
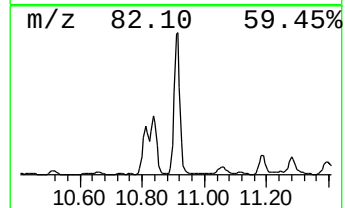
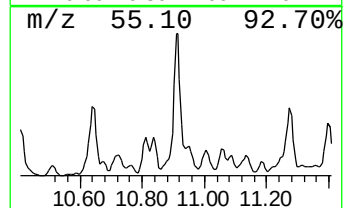
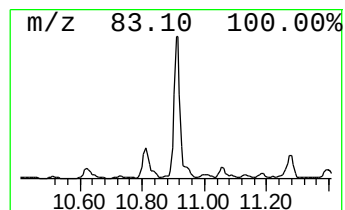
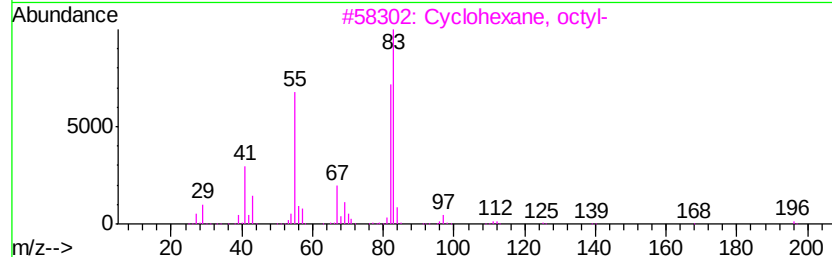
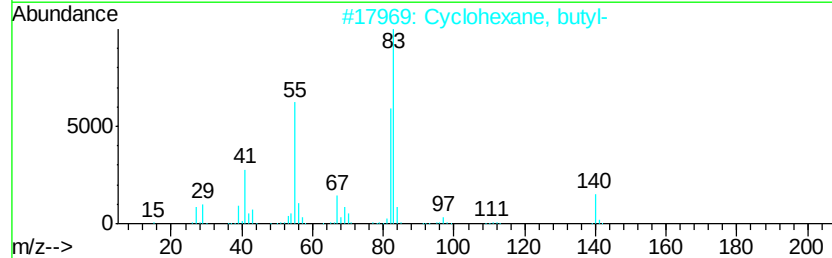
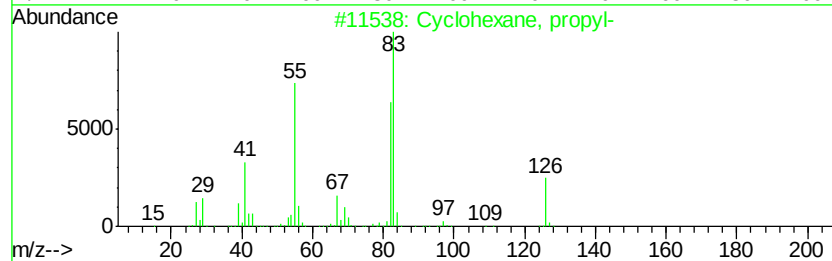
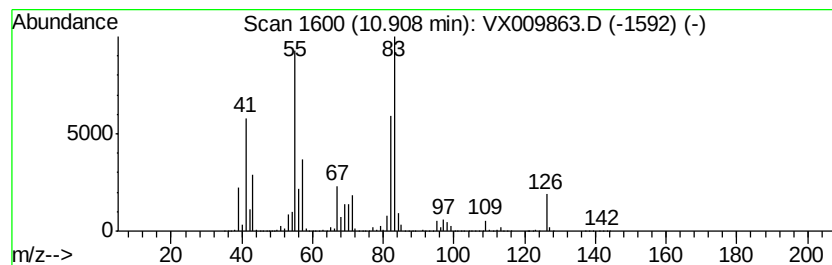
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	608.81 ug/l	9816380	Chlorobenzene-d5	10.11

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, propyl-	126	C9H18	001678-92-8	64
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	64
3		Cyclohexane, octyl-	196	C14H28	001795-15-9	59
4		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	58
5		Cyclopentane, 1-methyl-3-(1-meth...	126	C9H18	053771-88-3	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

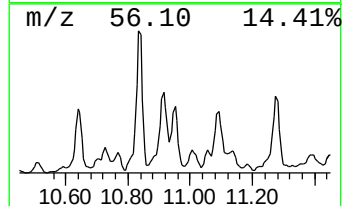
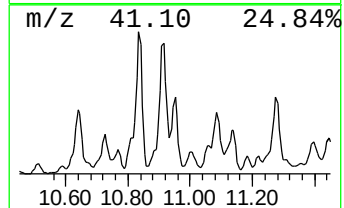
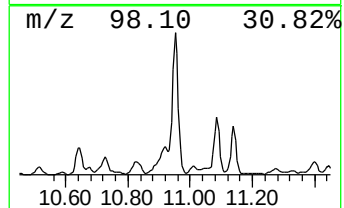
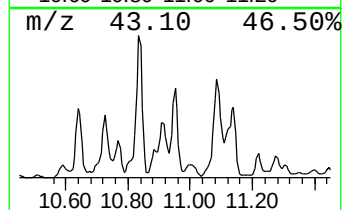
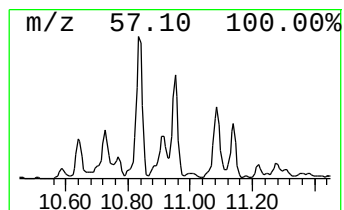
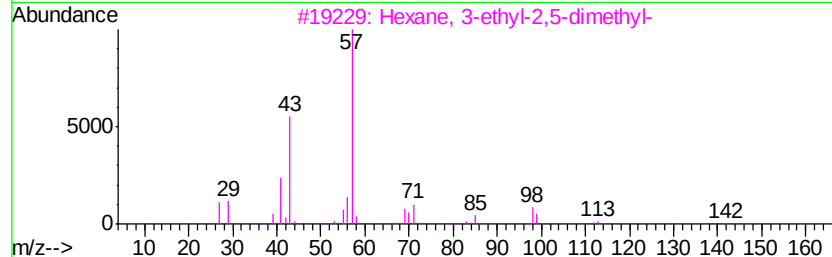
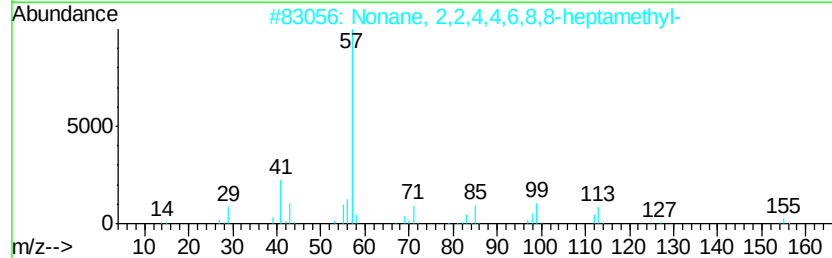
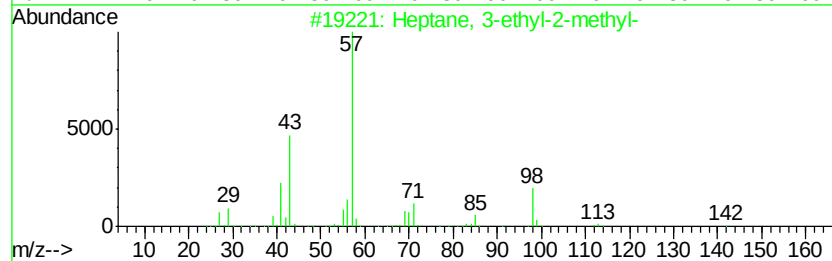
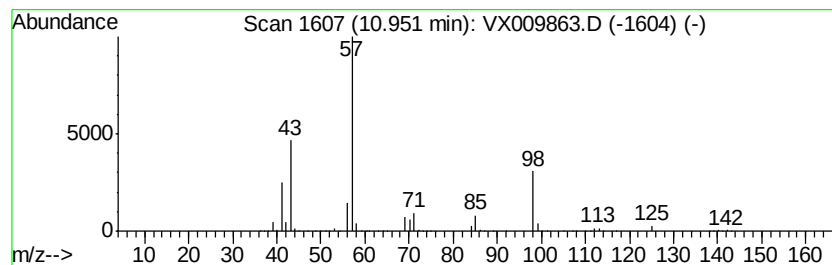
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 8 Heptane, 3-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.95	253.94 ug/l	4094500	Chlorobenzene-d5	10.11	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	64
2	Nonane, 2,2,4,4,6,8,8-heptamethyl-	226	C16H34	004390-04-9	53
3	Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	53
4	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	50
5	Undecane, 6-methyl-	170	C12H26	017302-33-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

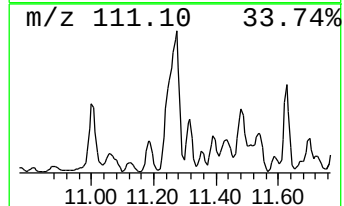
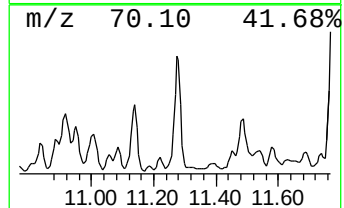
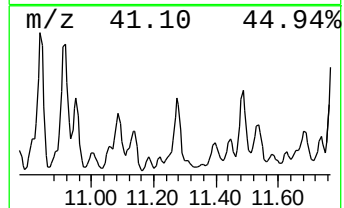
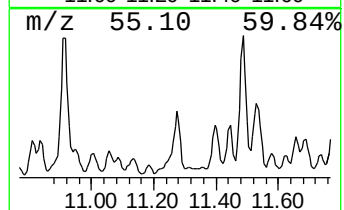
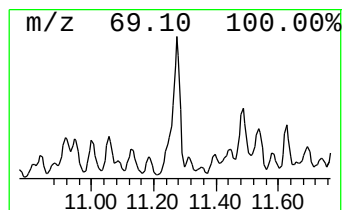
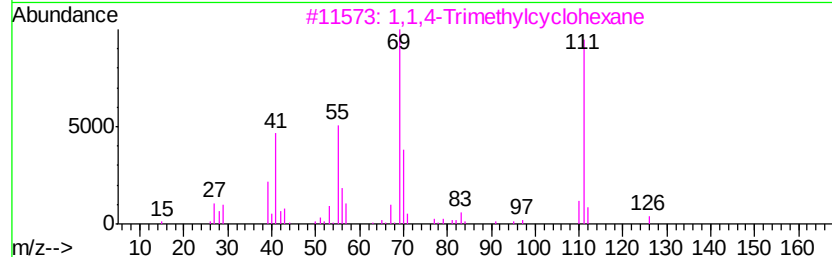
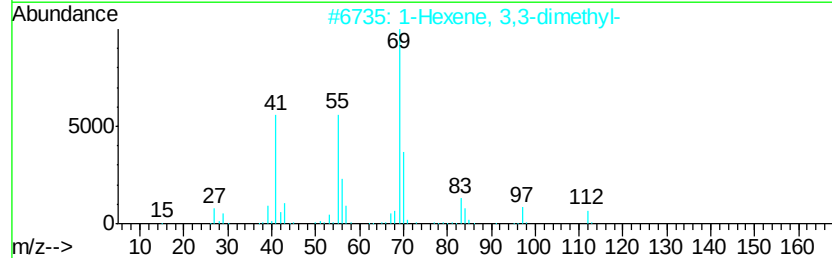
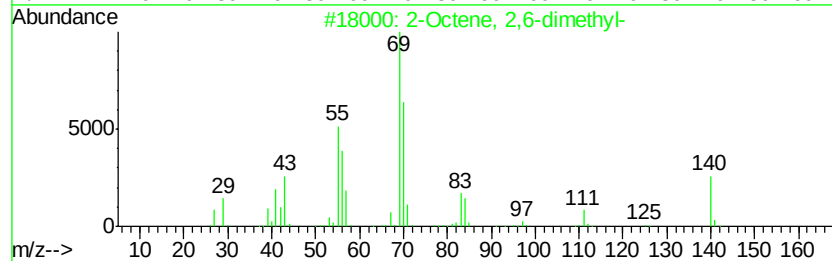
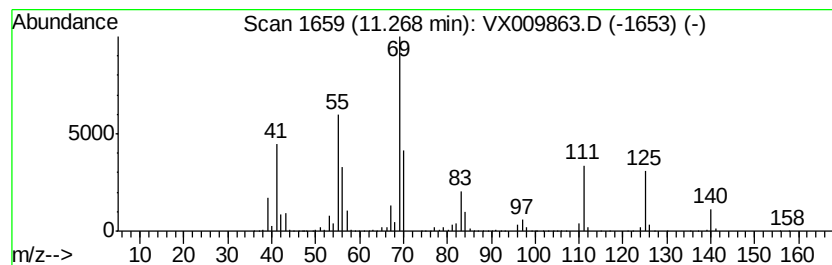
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ClientSampleId :
EP1-SB-E2-3R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 9 2-Octene, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.27	171.82 ug/l	6560210	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	64
2	1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	49
3	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	47
4	2-Octene, 2,7-dimethyl-	140	C10H20	002050-81-9	45
5	3-Ethyl-4-octene	140	C10H20	019781-32-9	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(17-20)

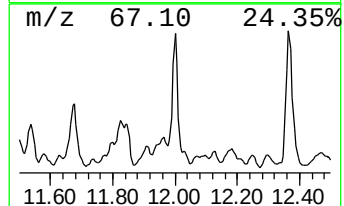
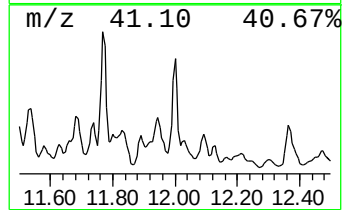
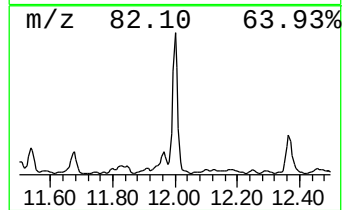
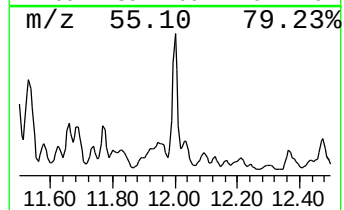
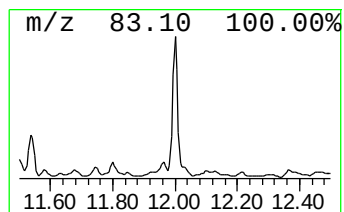
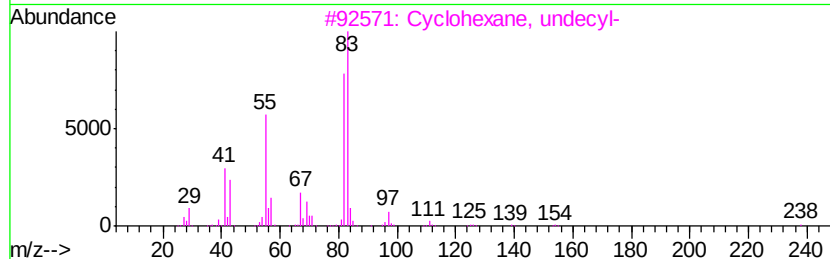
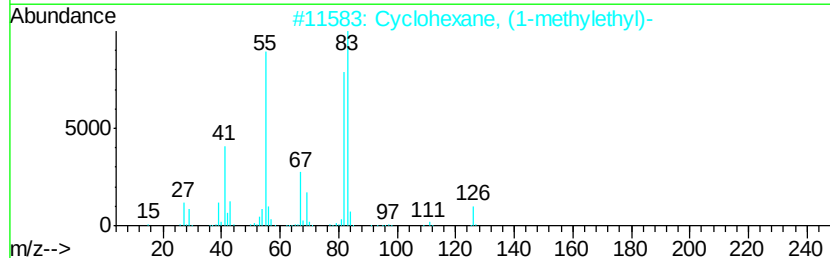
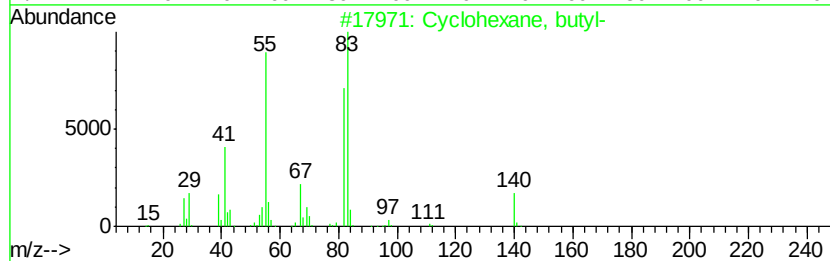
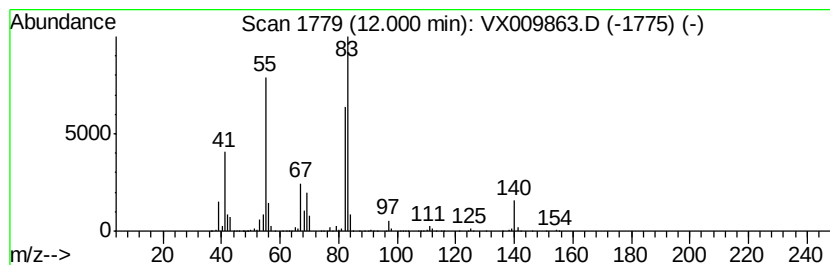
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 10 Cyclohexane, butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	132.03 ug/l	5041010	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, butyl-	140	C10H20	001678-93-9	91
2		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	80
3		Cyclohexane, undecyl-	238	C17H34	054105-66-7	78
4		Cyclohexane, pentyl-	154	C11H22	004292-92-6	64
5		Cyclohexane, ethyl-	112	C8H16	001678-91-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX052319\
Data File : VX009863.D
Acq On : 23 May 2019 17:04
Operator : JC/SP
Sample : K2976-02 10X
Misc : 5.76g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.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
Cyclohexane, 1,2,...	10.34	106.8	ug/l	1721770	3	10.11	806197	50.0
1-Ethyl-4-methylc...	10.38	128.0	ug/l	2064310	3	10.11	806197	50.0
Cyclohexane, 1-et...	10.41	119.9	ug/l	1932910	3	10.11	806197	50.0
Cyclohexane, 1-et...	10.64	318.1	ug/l	5128970	3	10.11	806197	50.0
Nonane, 4-methyl-	10.73	160.0	ug/l	2579730	3	10.11	806197	50.0
Octane, 2,6-dimet...	10.84	621.5	ug/l	10020300	3	10.11	806197	50.0
Cyclohexane, propyl-	10.91	608.8	ug/l	9816380	3	10.11	806197	50.0
Heptane, 3-ethyl-...	10.95	253.9	ug/l	4094500	3	10.11	806197	50.0
2-Octene, 2,6-dim...	11.27	171.8	ug/l	6560210	4	12.08	1909080	50.0
Cyclohexane, butyl-	12.00	132.0	ug/l	5041010	4	12.08	1909080	50.0