

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW090819\
 Data File : VW012769.D
 Acq On : 08 Sep 2019 15:10
 Operator : SY/VA
 Sample : K4603-06
 Misc : 3.28G/5ML/MSVOA W/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 PAR-E4-(0)

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_W\METHOD\82W090719S.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	2.148	35	40	50	rBV2	7768	20003	1.85%	0.309%
2	2.349	64	73	82	rBV	37028	68498	6.32%	1.059%
3	4.915	480	494	510	rBV3	17097	65685	6.06%	1.015%
4	5.434	566	579	590	rBV4	8857	31831	2.94%	0.492%
5	5.927	652	660	672	rVB5	10653	33197	3.06%	0.513%
6	7.543	915	925	937	rBV3	6416	18920	1.75%	0.293%
7	7.884	971	981	985	rBV	143640	337663	31.17%	5.220%
8	7.945	985	991	1001	rVB	306816	668821	61.73%	10.340%
9	8.305	1039	1050	1061	rBV	142424	325608	30.05%	5.034%
10	8.842	1129	1138	1149	rBV	419211	843015	77.81%	13.033%
11	10.323	1373	1381	1388	rBV	609429	1083452	100.00%	16.750%
12	10.384	1388	1391	1400	rVB	14992	28013	2.59%	0.433%
13	10.707	1437	1444	1448	rVB2	5543	11356	1.05%	0.176%
14	11.628	1587	1595	1607	rBV	541161	920008	84.91%	14.223%
15	12.615	1749	1757	1764	rBV	422729	740879	68.38%	11.454%
16	13.554	1904	1911	1920	rBV	459953	790580	72.97%	12.222%
17	13.749	1934	1943	1958	rVB6	43584	178013	16.43%	2.752%
18	13.871	1960	1963	1967	rBV4	8932	15118	1.40%	0.234%
19	13.926	1967	1972	1973	rBV4	11295	16753	1.55%	0.259%
20	14.066	1985	1995	2004	rBV4	28393	82103	7.58%	1.269%
21	14.523	2068	2070	2076	rVB6	9834	10978	1.01%	0.170%
22	14.725	2099	2103	2106	rVB	20154	26304	2.43%	0.407%
23	15.340	2200	2204	2207	rBV6	11453	21284	1.96%	0.329%
24	15.724	2259	2267	2275	rVB6	37850	130215	12.02%	2.013%

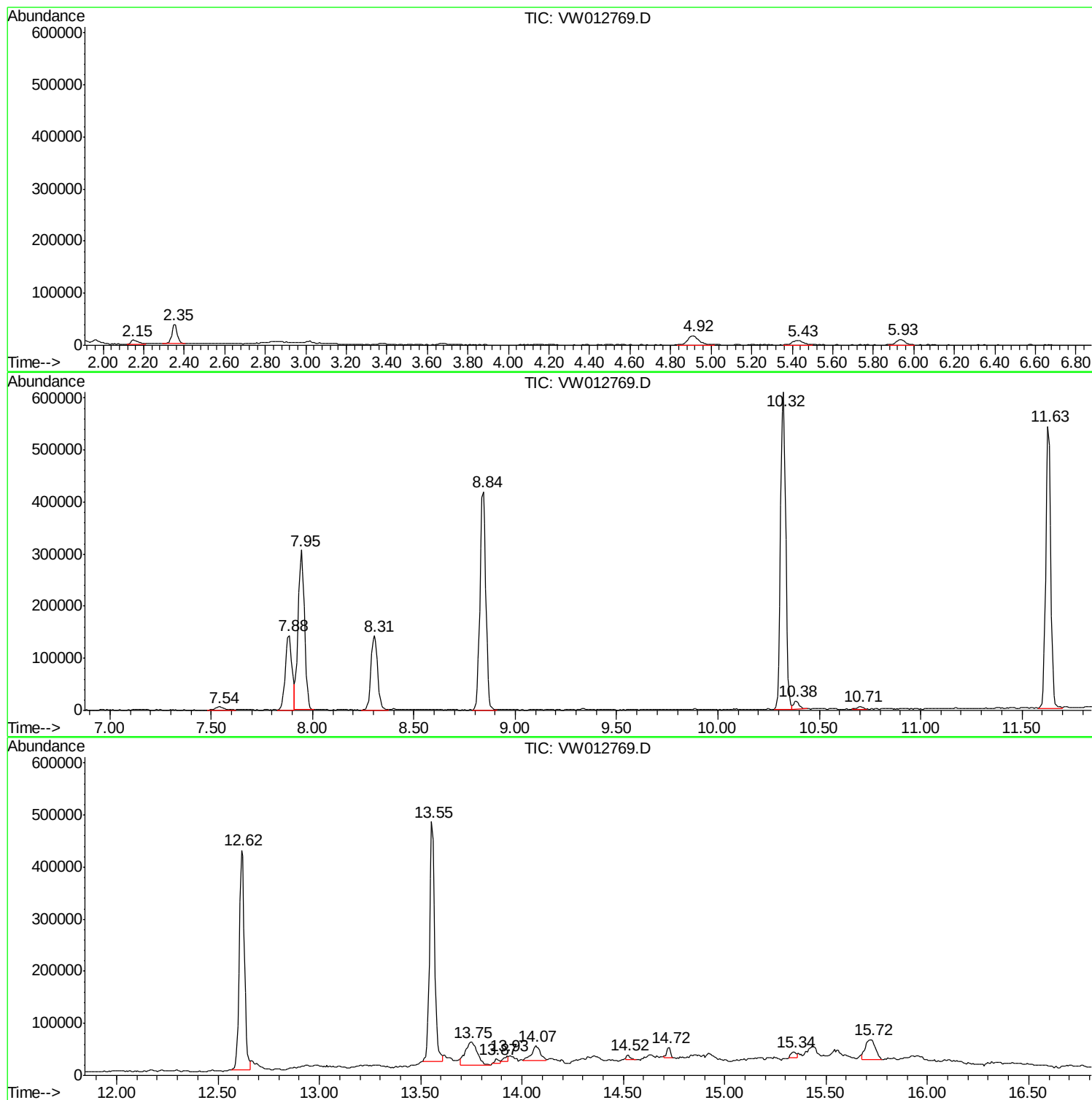
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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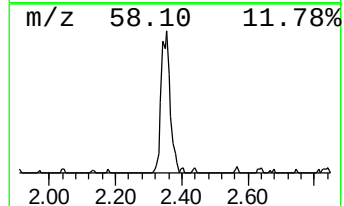
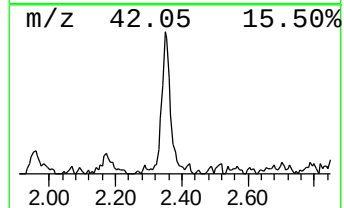
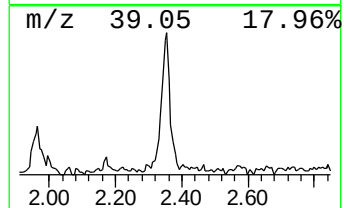
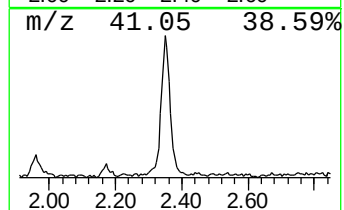
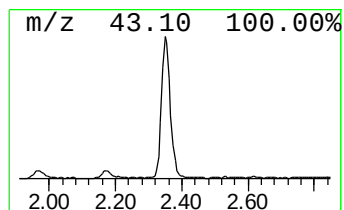
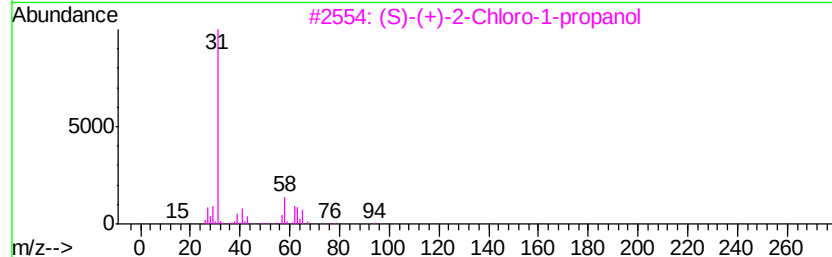
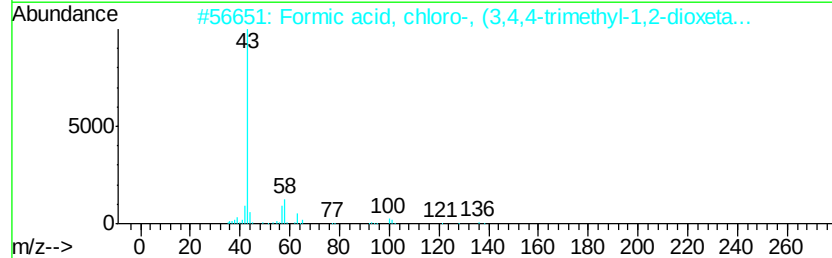
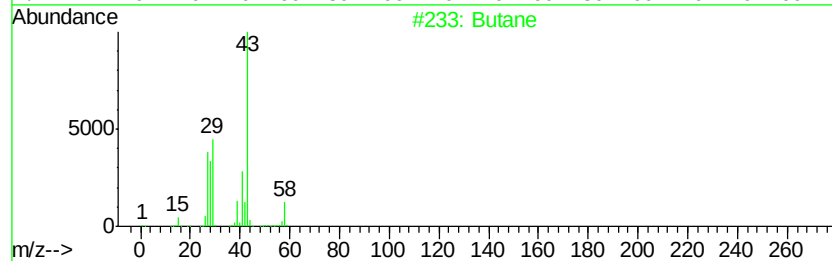
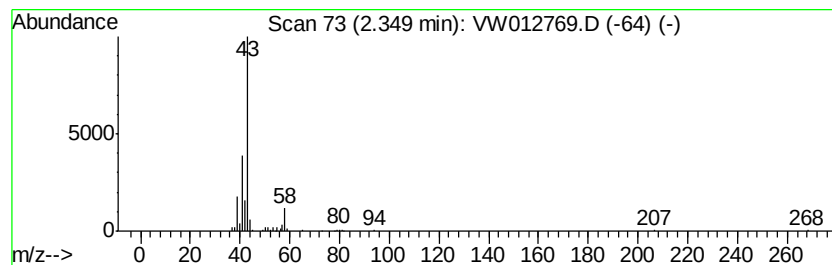
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W090719S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.35	5.12 ug/l	68498	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane	58	C4H10	000106-97-8	64
2		Formic acid, chloro-, (3,4,4-tri...	194	C7H11ClO4	107323-92-2	38
3		(S)-(+)-2-Chloro-1-propanol	94	C3H7ClO	019210-21-0	9
4		Isobutane	58	C4H10	000075-28-5	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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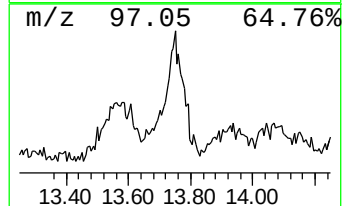
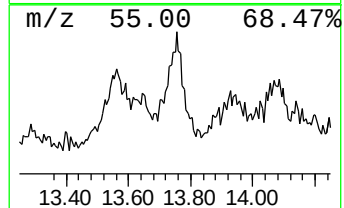
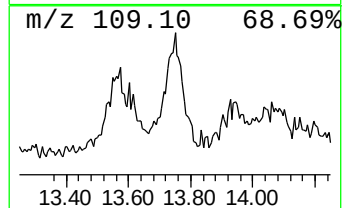
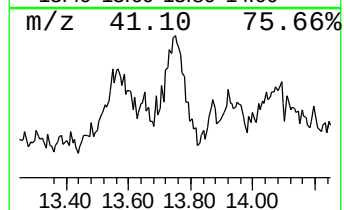
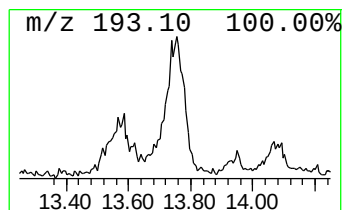
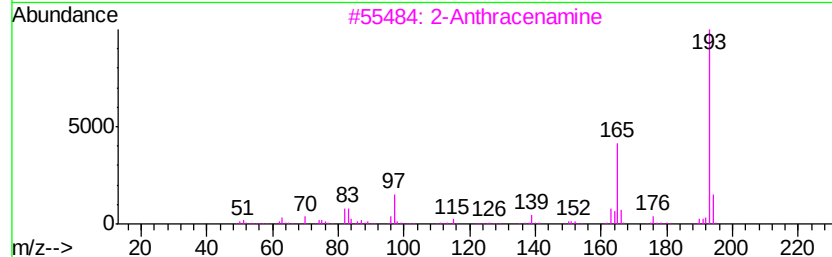
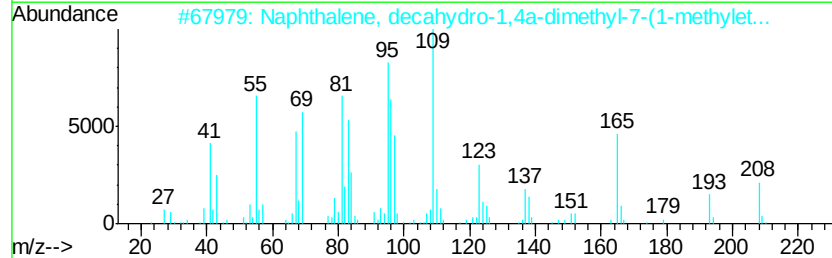
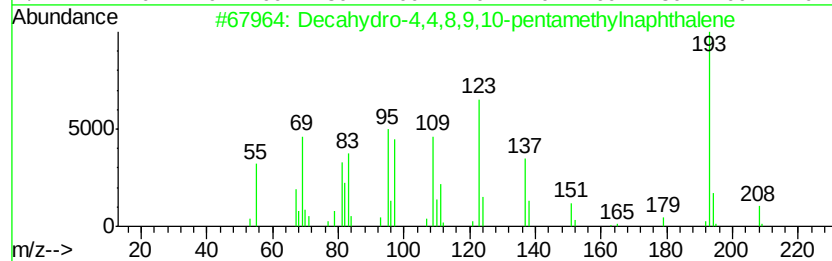
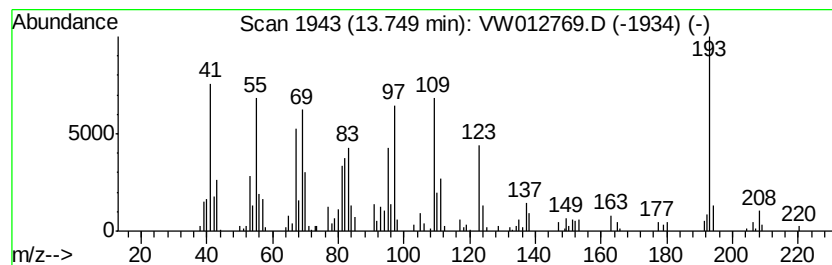
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 TIC Integration Parameters: LSCINT.P

 Peak Number 2 Decahydro-4,4,8,9,10-pentam... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	11.26 ug/l	178013	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	70
2		Naphthalene, decahydro-1,4a-dime...	208	C15H28	030824-81-8	38
3		2-Anthracenamine	193	C14H11N	000613-13-8	25
4		n-Octylidencyclohexane	194	C14H26	137595-12-1	25
5		Cyclohexene, 3,3,5-trimethyl-	124	C9H16	000503-45-7	25



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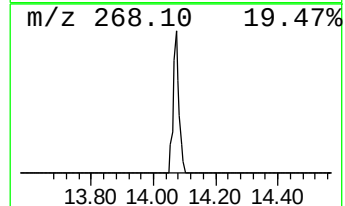
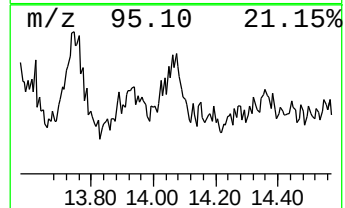
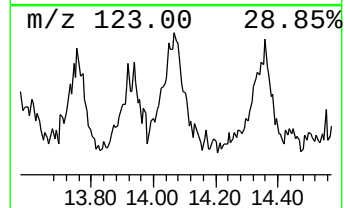
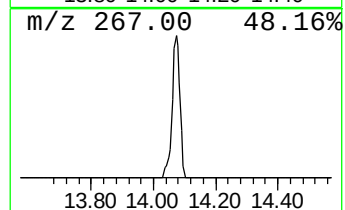
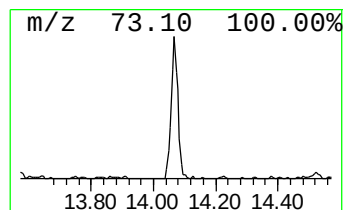
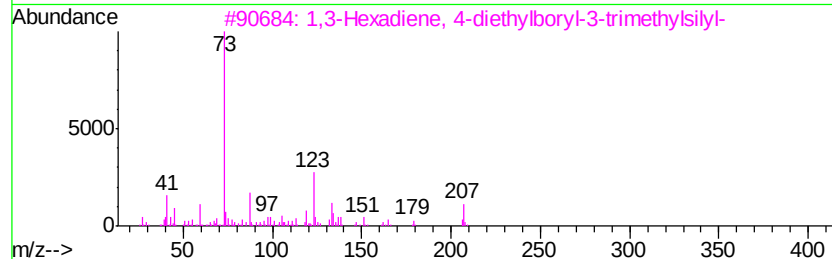
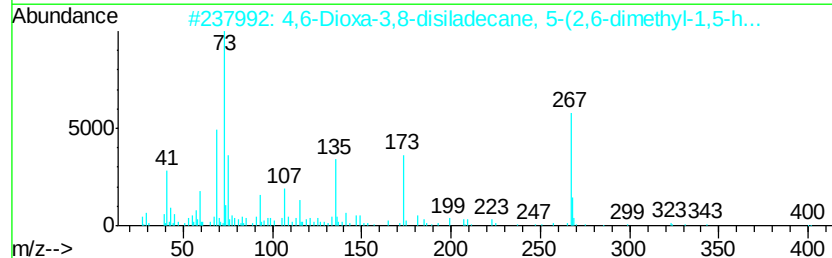
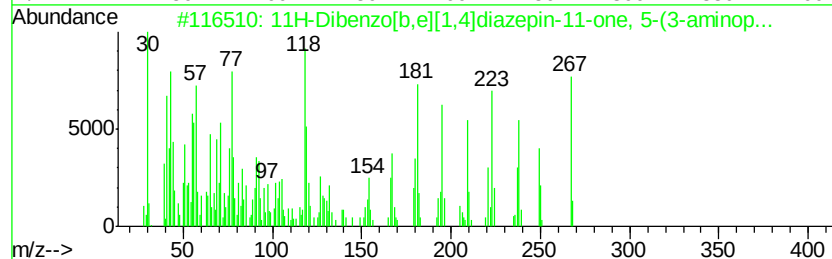
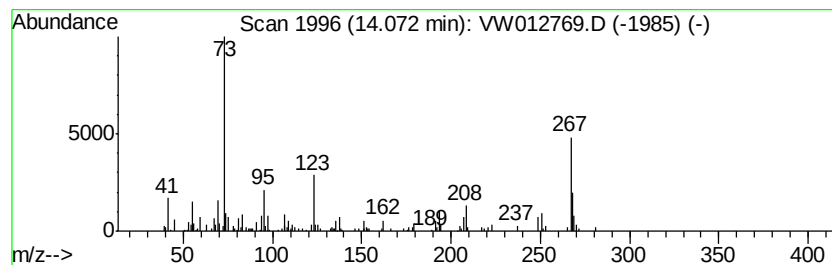
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Peak Number 3 unknown14.07 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.07	5.19 ug/l	82103	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	25
2		4,6-Dioxa-3,8-disiladecane, 5-(2...	532	C32H60O2Si2	109629-49-4	17
3		1,3-Hexadiene, 4-diethylborvl-3-...	236	C14H29BSi	1000155-03-7	12
4		Silane, (2-ethyl-5,5-dimethyl-4-...	208	C13H24Si	095798-05-3	9
5		Silane, trimethyl[[p-(trimethyls...	268	C13H24O2Si2	018401-58-6	9



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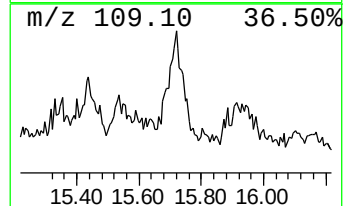
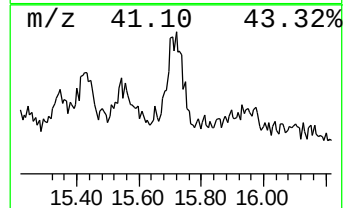
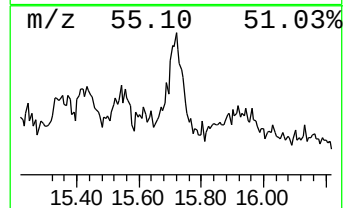
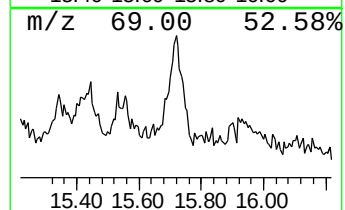
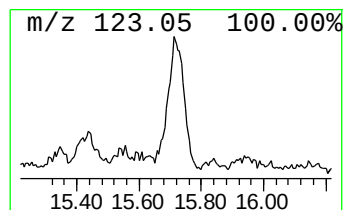
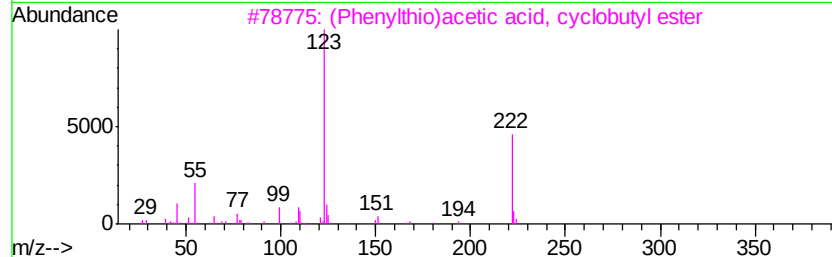
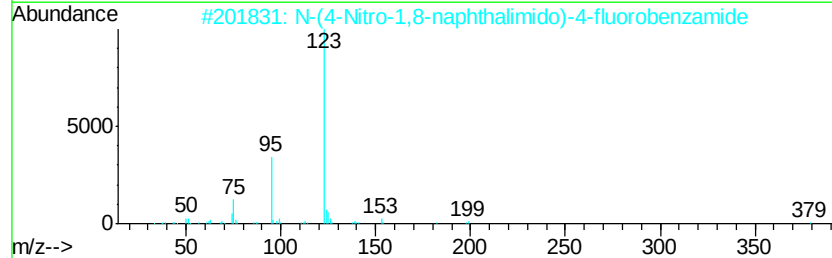
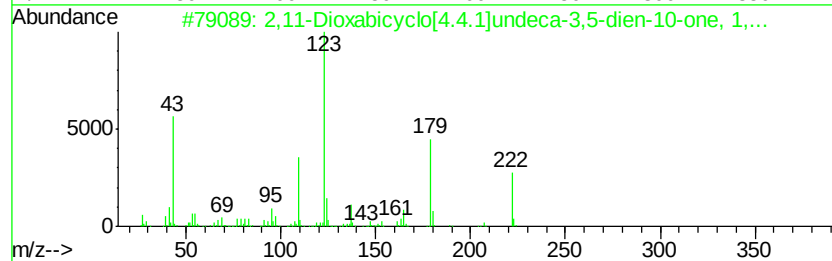
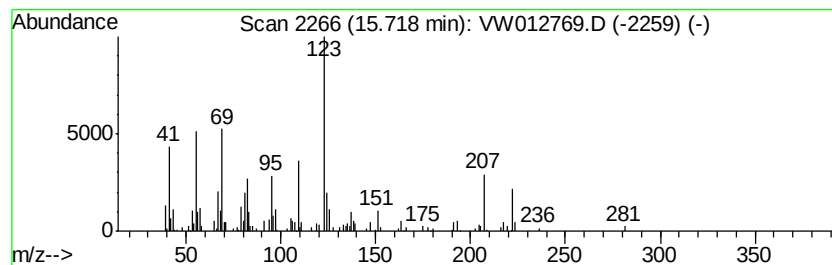
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Peak Number 4 unknown15.72 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	8.24 ug/l	130215	1,4-Dichlorobenzene-d4	13.55

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2	11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3	070412-52-1	43	
2	N-(4-Nitro-1,8-naphthalimido)-4-	379	C19H10FN3O5	300690-90-8	38		
3	(Phenylthio)acetic acid, cyclobu...	222	C12H14O2S	1000299-95-5	38		
4	3-Fluorobenzoic acid, propargyl ...	178	C10H7FO2	1000325-53-9	38		
5	trans-Chrysanthemal	152	C10H16O	020104-05-6	35		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane	2.35	5.1	ug/l	68498	1	7.95	668821	50.0
Decahydro-4,4,8,9...	13.75	11.3	ug/l	178013	4	13.55	790580	50.0
unknown14.07	14.07	5.2	ug/l	82103	4	13.55	790580	50.0
unknown15.72	15.72	8.2	ug/l	130215	4	13.55	790580	50.0