

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121218\
 Data File : BN003865.D
 Acq On : 12 Dec 2018 23:29
 Operator : JU/SJ
 Sample : J6299-10
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9F37

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	25	rVB	142279	218452	18.67%	3.082%
2	6.987	675	680	688	rBB2	14811	25148	2.15%	0.355%
3	7.164	705	710	720	rBB	63902	98861	8.45%	1.395%
4	7.358	738	743	750	rBB	63241	96457	8.24%	1.361%
5	7.834	818	824	828	rBV	50591	80131	6.85%	1.131%
6	8.522	936	941	950	rBB	50540	75644	6.46%	1.067%
7	8.922	1004	1009	1015	rBB	30287	45217	3.86%	0.638%
8	8.999	1015	1022	1027	rBV	81954	135882	11.61%	1.917%
9	9.711	1138	1143	1150	rBB	70524	106250	9.08%	1.499%
10	10.240	1227	1233	1242	rBV	97903	158864	13.58%	2.242%
11	10.628	1292	1299	1305	rBV	88979	145873	12.47%	2.058%
12	11.905	1511	1516	1521	rBB	8886	13374	1.14%	0.189%
13	12.804	1665	1669	1679	rBV	11434	20279	1.73%	0.286%
14	13.057	1707	1712	1717	rBB	27237	37657	3.22%	0.531%
15	13.869	1845	1850	1858	rBV	243953	327435	27.98%	4.620%
16	14.163	1893	1900	1910	rBV	271171	397720	33.99%	5.612%
17	14.469	1945	1952	1959	rVB2	150964	228265	19.51%	3.221%
18	14.663	1979	1985	1994	rBV	13958	21671	1.85%	0.306%
19	15.457	2114	2120	2129	rBV	390351	536327	45.84%	7.568%
20	15.581	2136	2141	2146	rBV	116414	151342	12.93%	2.135%
21	17.210	2412	2418	2426	rBV	228151	307932	26.32%	4.345%
22	17.310	2429	2435	2447	rVB	436136	613908	52.47%	8.662%
23	17.857	2525	2528	2532	rVB	33227	33333	2.85%	0.470%
24	19.604	2819	2825	2833	rBV	599022	798697	68.26%	11.270%
25	21.392	3123	3129	3135	rBV	375397	476975	40.76%	6.730%
26	22.427	3301	3305	3309	rBV2	14435	17721	1.51%	0.250%
27	22.945	3389	3393	3398	rVB3	19445	27804	2.38%	0.392%
28	23.557	3488	3497	3507	rVB	622464	1170120	100.00%	16.511%
29	23.704	3514	3522	3530	rBV	295181	564379	48.23%	7.964%
30	24.192	3600	3605	3613	rBV	21931	43562	3.72%	0.615%
31	24.963	3730	3736	3745	rVB	21919	48000	4.10%	0.677%
32	25.863	3884	3889	3896	rVB2	14472	31893	2.73%	0.450%
33	26.921	4063	4069	4078	rVB2	11325	31897	2.73%	0.450%

LSC Area Percent Report

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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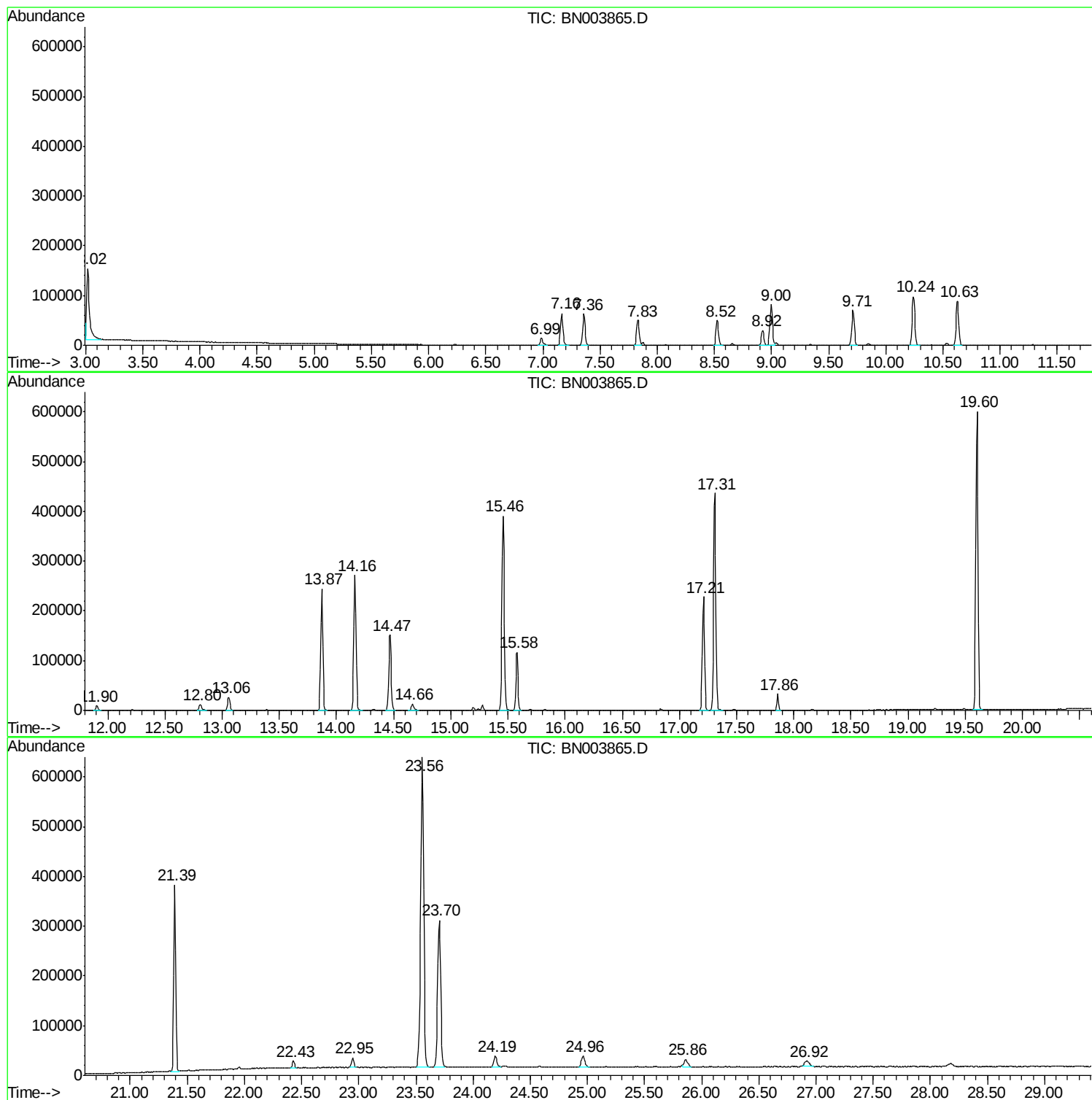
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Quant Title : SVOA CALIBRATION

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



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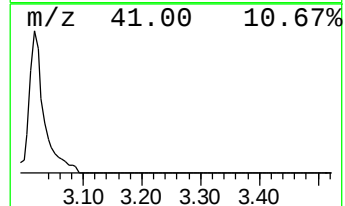
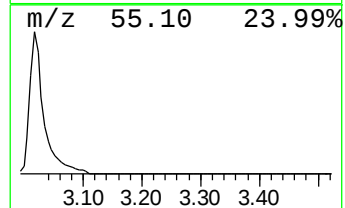
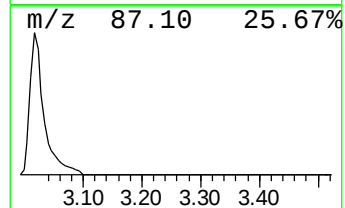
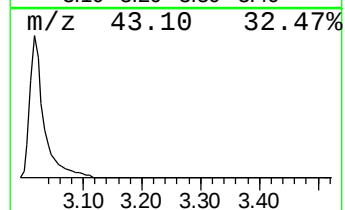
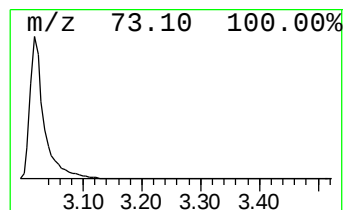
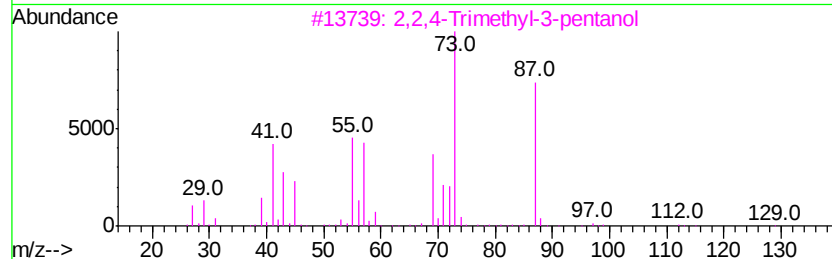
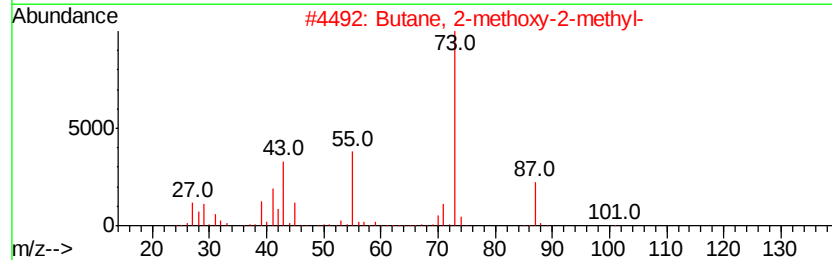
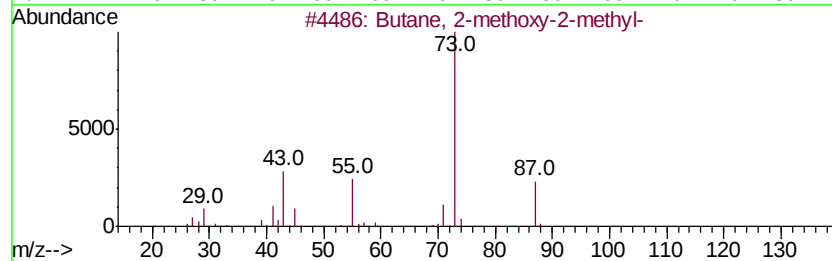
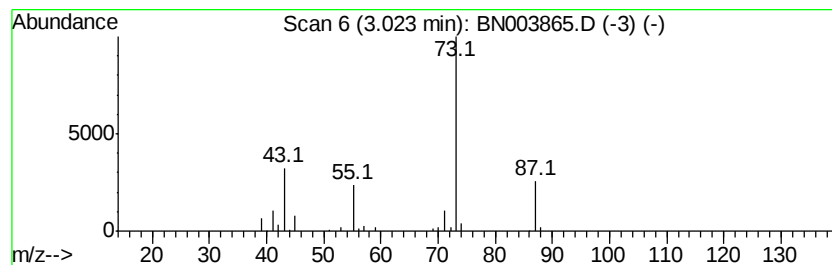
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	54.52 ng/ul	218452	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23



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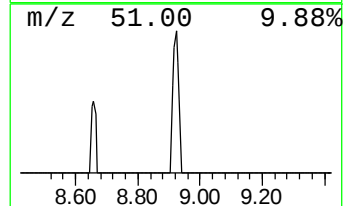
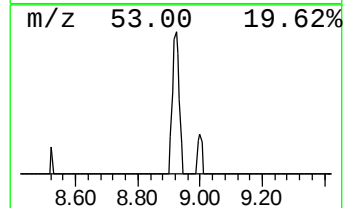
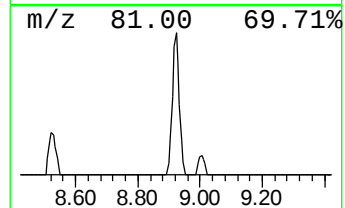
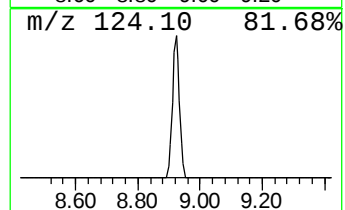
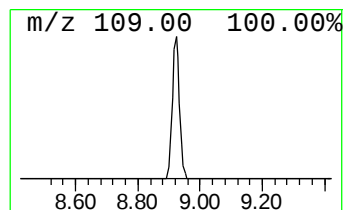
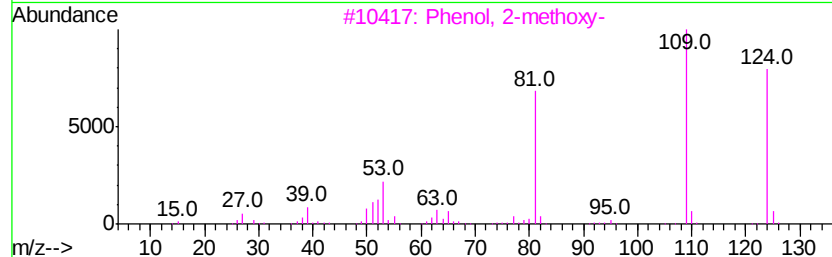
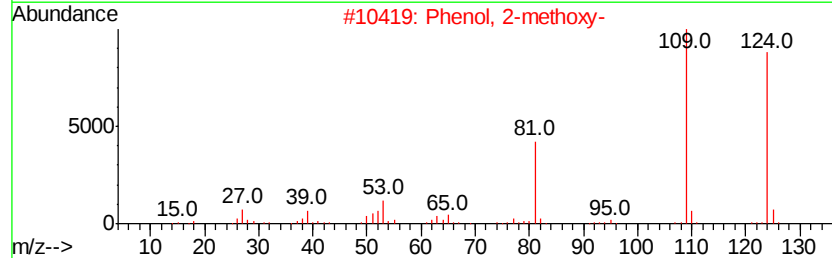
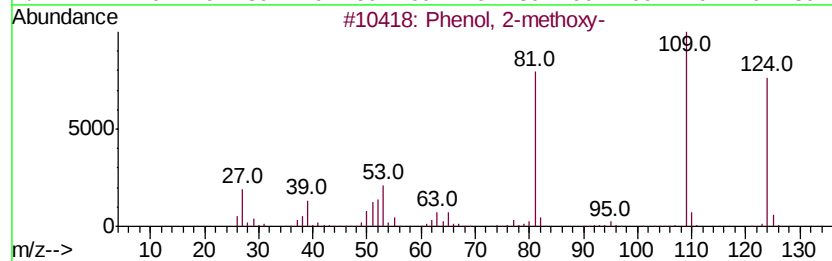
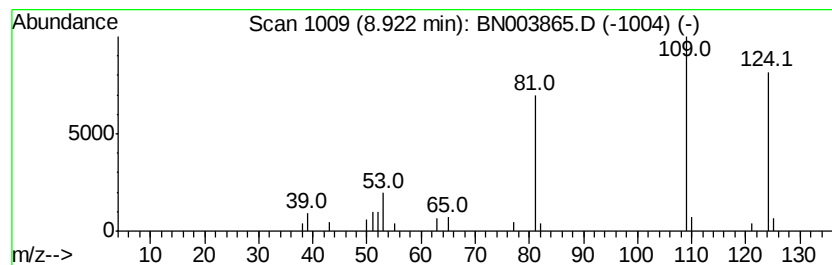
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 2-methoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.92	11.29 ng/ul	45217	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	91
2		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	91
3		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	91
4		Ethanone, 1-(2-methyl-1-cyclopropyl)-	124	C8H12O	003168-90-9	86
5		Mequinol	124	C7H8O2	000150-76-5	86



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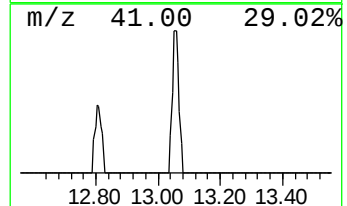
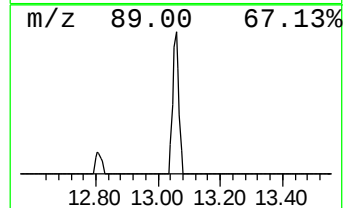
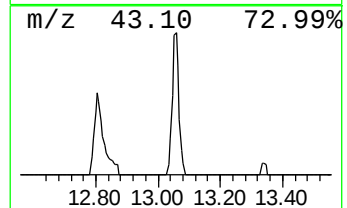
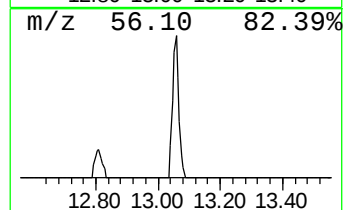
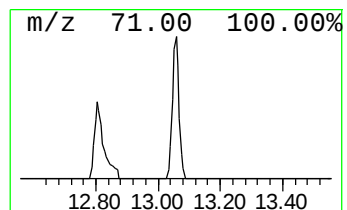
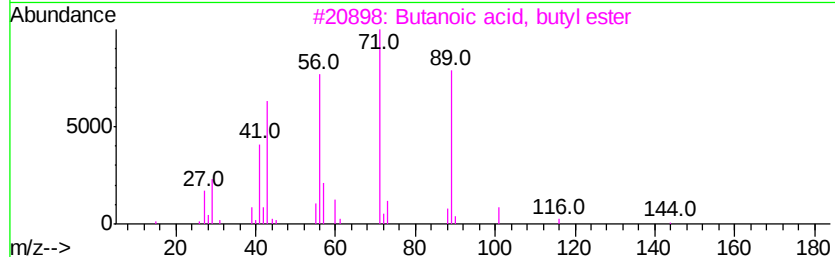
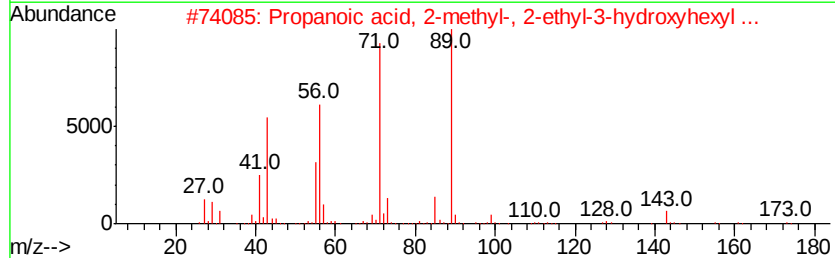
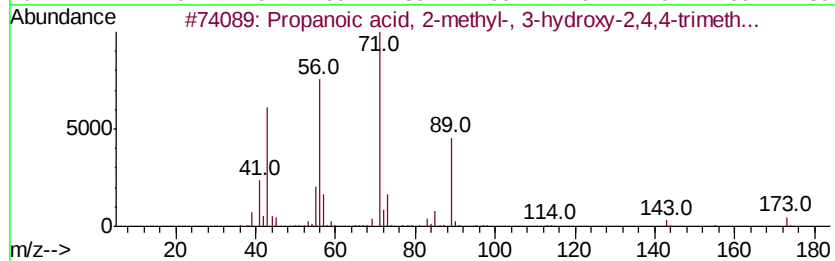
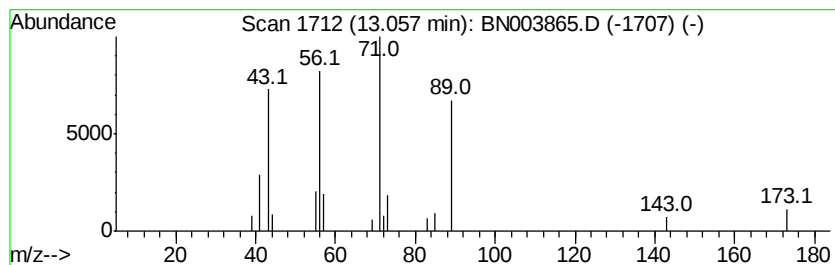
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Propanoic acid, 2-methyl-, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.06	3.30 ng/ul	37657	Acenaphthene-d10	14.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	074367-34-3	83
2		Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	78
3		Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	72
4		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	42
5		Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	40



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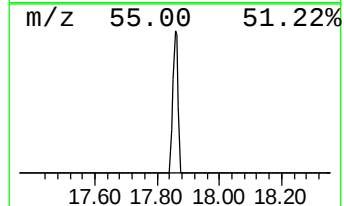
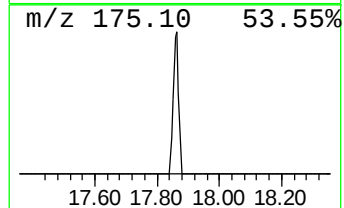
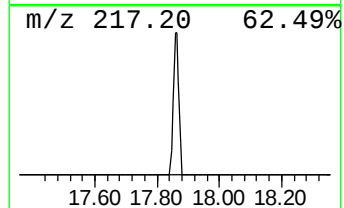
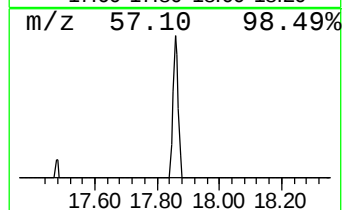
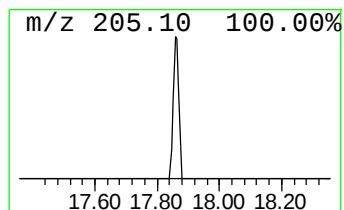
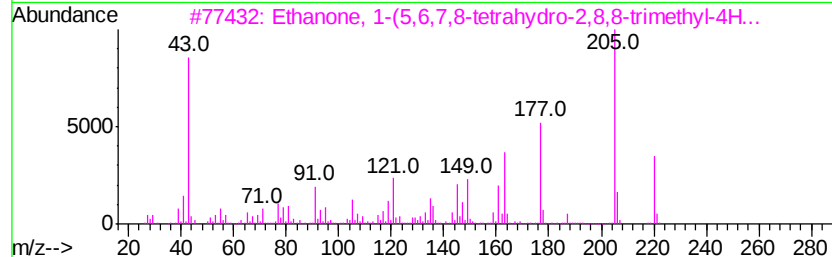
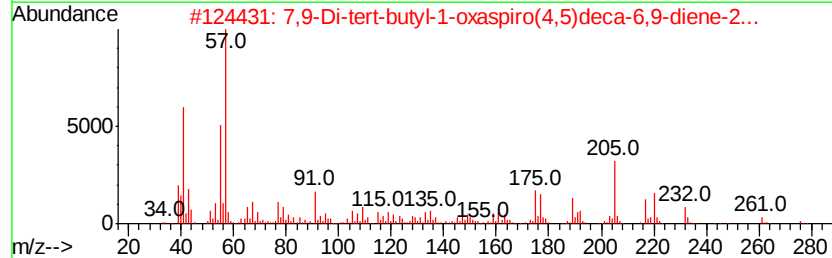
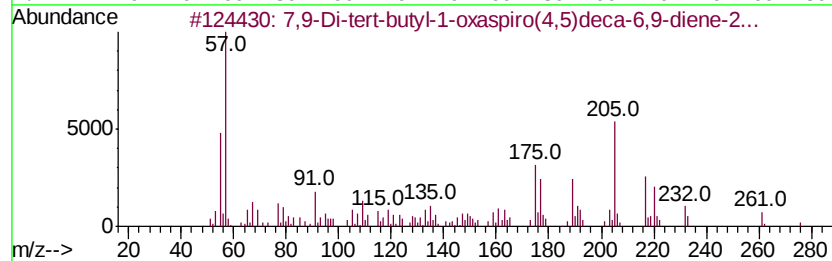
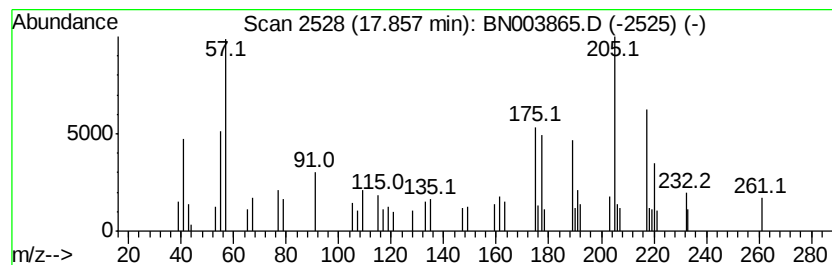
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.86	2.16 ng/ul	33333	Phenanthrene-d10	17.21		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	91	
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	72	
3	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	41	
4	1,3-Pentadiene, 1,1-diphenyl-, (Z)-	220	C17H16	015295-31-5	30	
5	2,4-Dimethyl-5,6-dimethoxy-8-ami...	232	C13H16N2O2	064993-02-8	22	



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.02	54.5	ng/ul	218452	1	7.83	80131	20.0
Phenol, 2-methoxy-	8.92	11.3	ng/ul	45217	1	7.83	80131	20.0
Propanoic acid, 2...	13.06	3.3	ng/ul	37657	3	14.46	228265	20.0
7,9-Di-tert-butyl...	17.86	2.2	ng/ul	33333	4	17.21	307932	20.0