

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082318\
 Data File : BF108422.D
 Acq On : 23 Aug 2018 12:47
 Operator : JU/SJ
 Sample : PB112242BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB112242BL

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.449	11	19	25	rVB	91534	142189	2.74%	0.364%
2	5.272	492	499	509	rBV	121580	178413	3.44%	0.457%
3	5.637	556	561	564	rBV	3580200	3825461	73.76%	9.790%
4	6.625	722	729	732	rBV	3693446	4029971	77.70%	10.313%
5	6.772	749	754	757	rBV	4018450	4038887	77.87%	10.336%
6	7.001	789	793	796	rBV	1007162	949397	18.30%	2.430%
7	7.154	814	819	823	rBV	3170320	3219674	62.08%	8.240%
8	7.560	882	888	892	rBV	2264516	2713999	52.33%	6.946%
9	8.283	1005	1011	1014	rBV	1258444	1196263	23.06%	3.061%
10	9.360	1187	1194	1197	rBV	4907347	5153961	99.37%	13.190%
11	10.042	1304	1310	1320	rBV2	1481527	1452420	28.00%	3.717%
12	10.842	1439	1446	1463	rBV	3019794	3238900	62.45%	8.289%
13	11.542	1559	1565	1568	rBV	1663491	1549423	29.87%	3.965%
14	13.130	1830	1835	1838	rBV	5417235	5186586	100.00%	13.273%
15	14.195	2012	2016	2024	rBV	1130275	1093679	21.09%	2.799%
16	15.748	2272	2280	2290	rVB	573026	848926	16.37%	2.173%
17	16.201	2352	2357	2362	rBV2	59050	91588	1.77%	0.234%
18	16.753	2445	2451	2458	rVB2	49352	97284	1.88%	0.249%
19	17.401	2555	2561	2569	rVB2	30766	68005	1.31%	0.174%

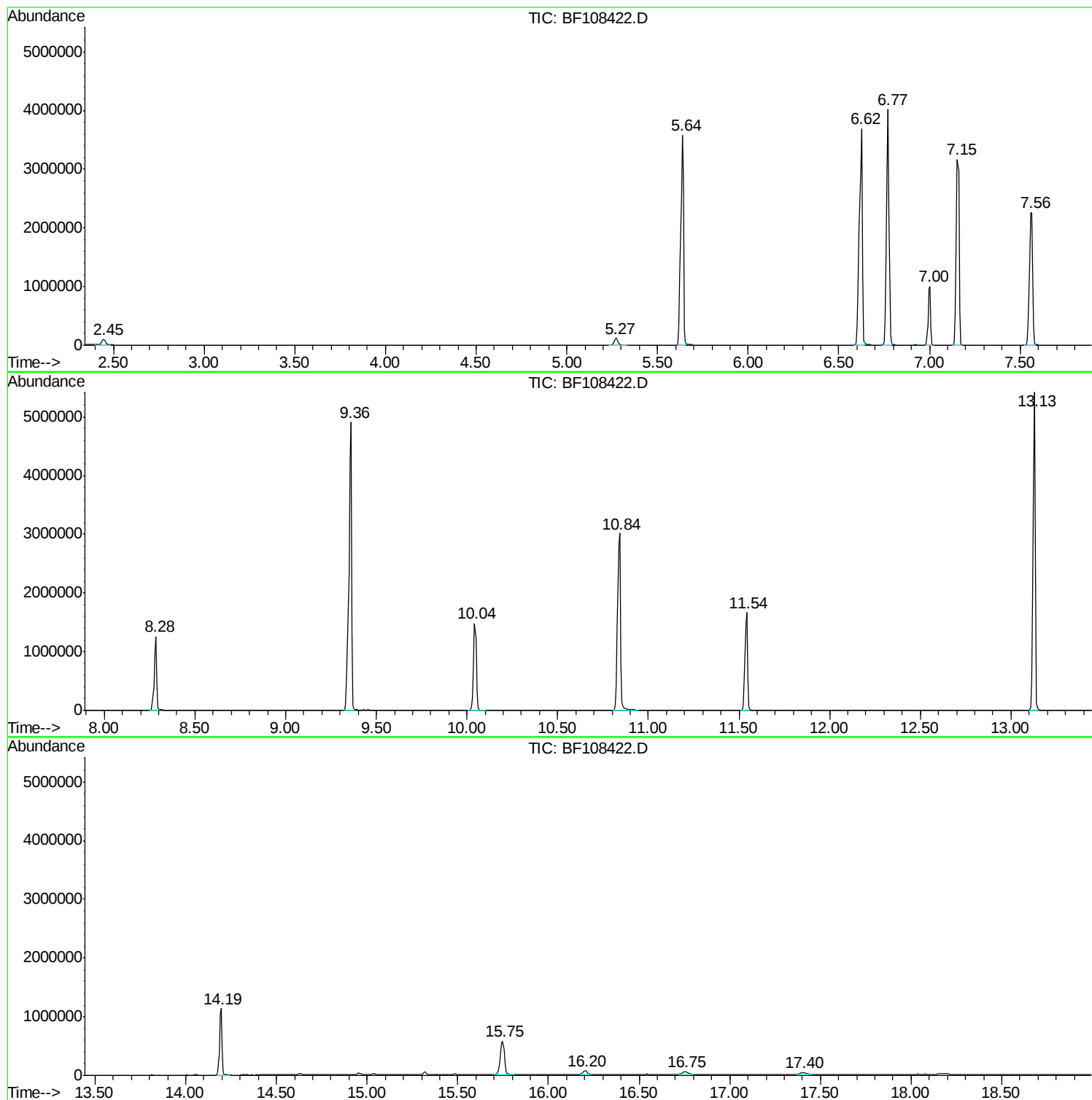
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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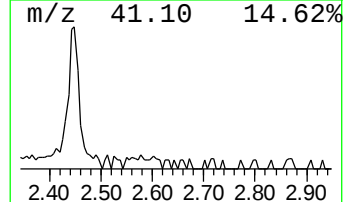
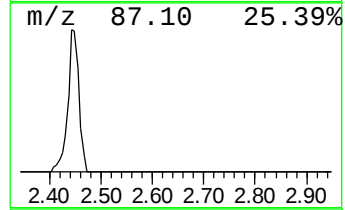
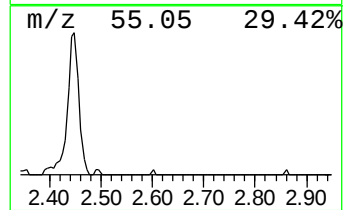
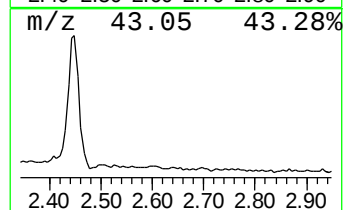
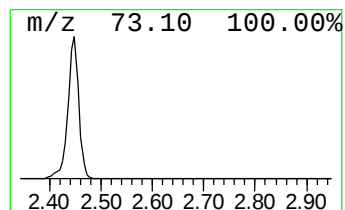
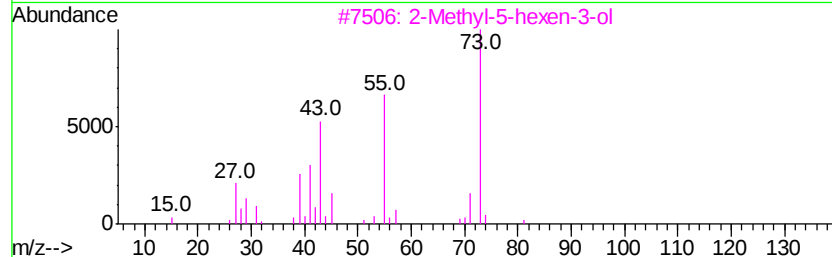
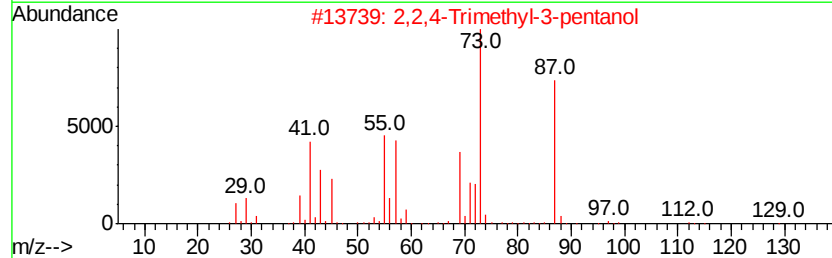
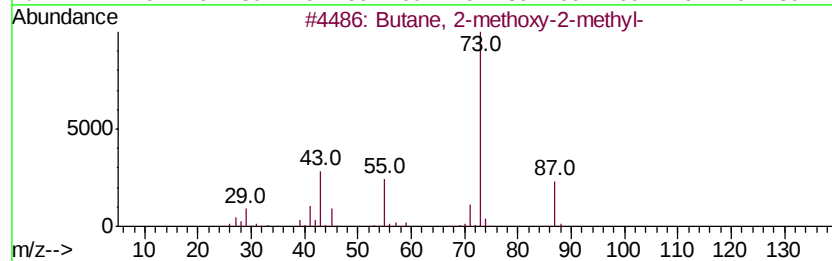
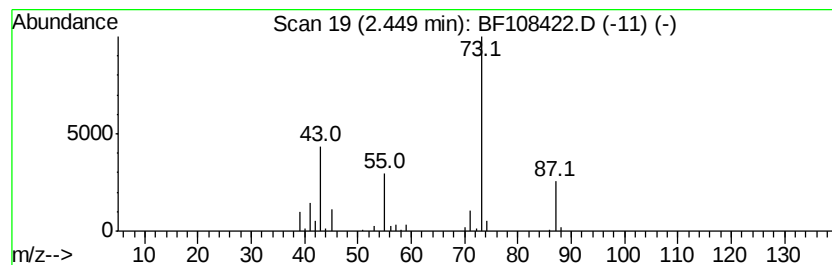
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.45	3.00 ng	142189	1,4-Dichlorobenzene-d4	7.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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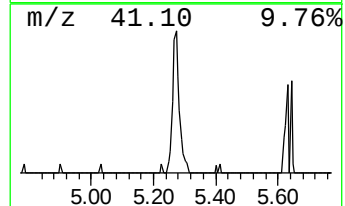
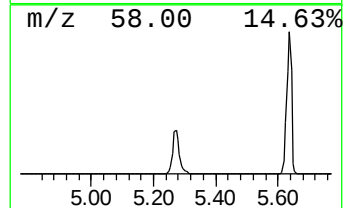
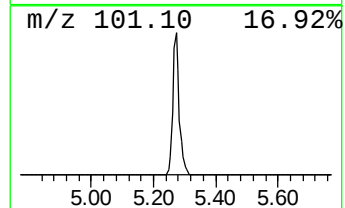
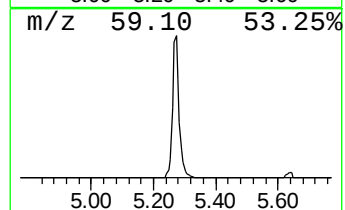
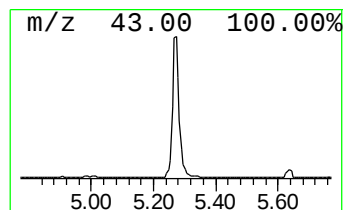
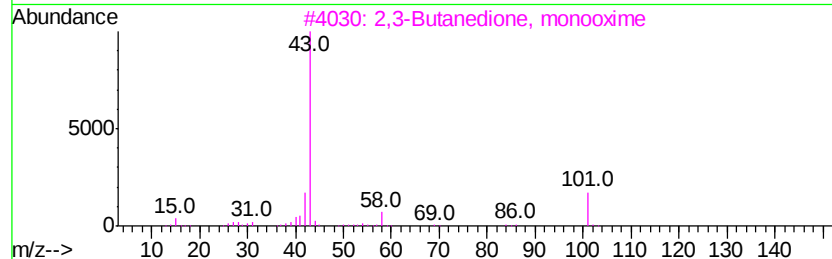
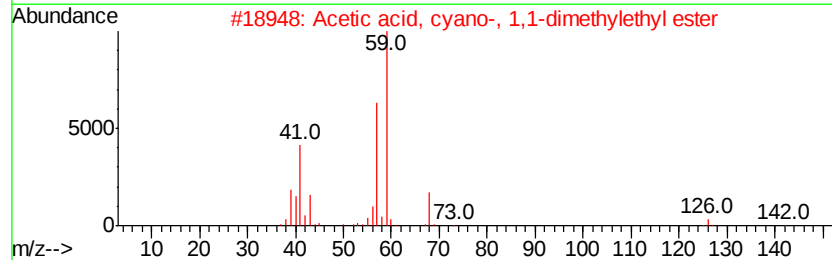
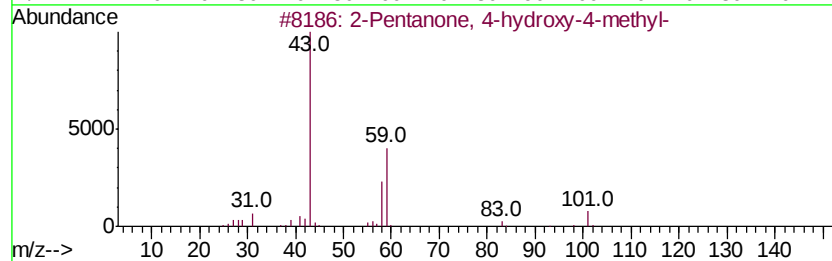
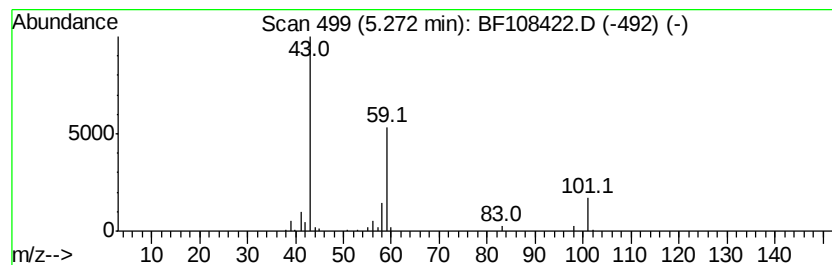
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TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.27	3.76 ng	178413	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5		Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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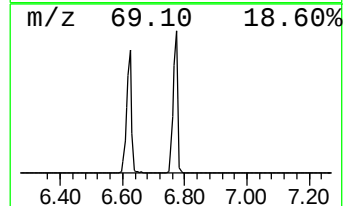
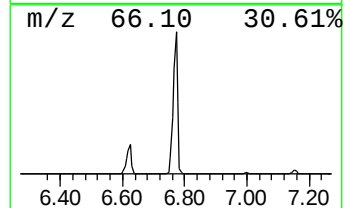
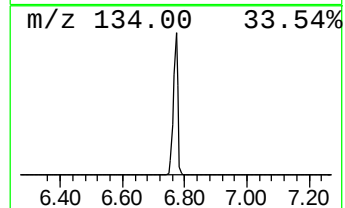
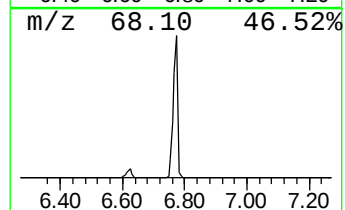
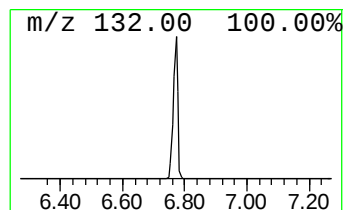
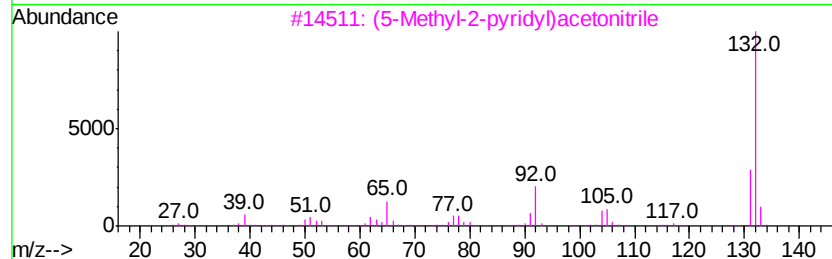
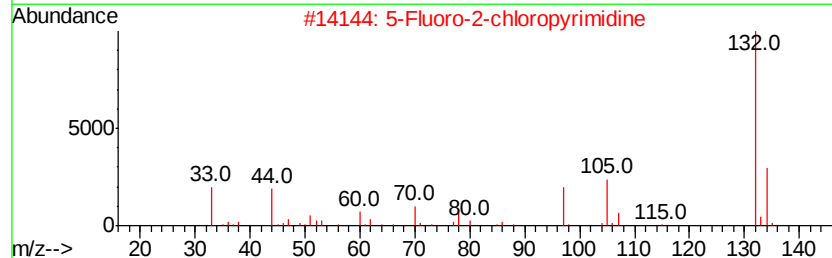
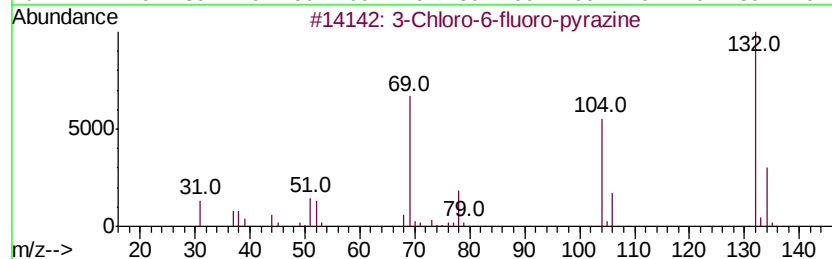
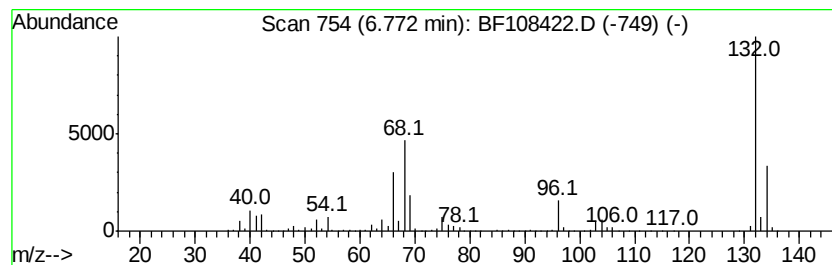
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Peak Number 3 unknown6.77 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.77	85.08 ng	4038890	1,4-Dichlorobenzene-d4	7.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Fluoro-2-chloropyrimidine			132	C4H2ClFN2	062802-42-0	12
3	(5-Methyl-2-pyridyl)acetonitrile			132	C8H8N2	1000241-93-9	10
4	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
5	1,3,5-Trifluorobenzene			132	C6H3F3	000372-38-3	9



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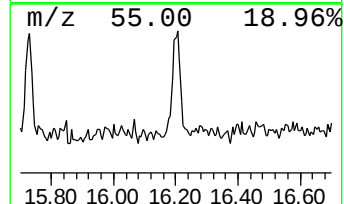
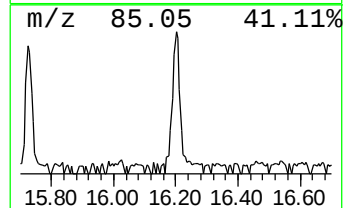
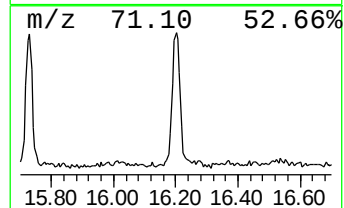
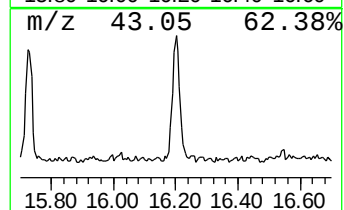
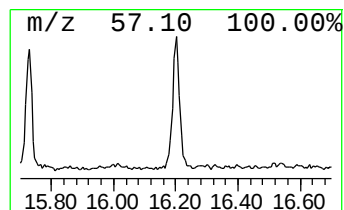
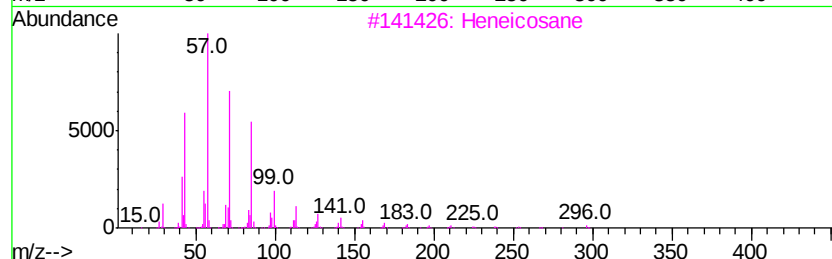
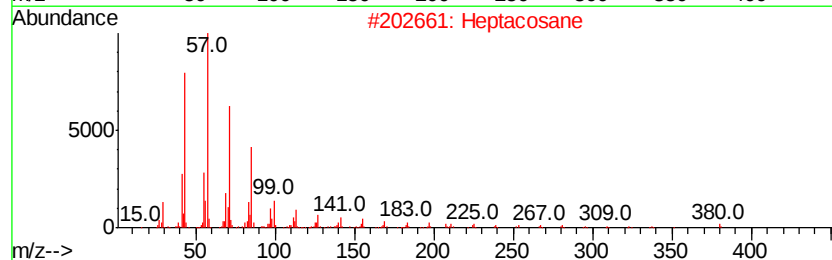
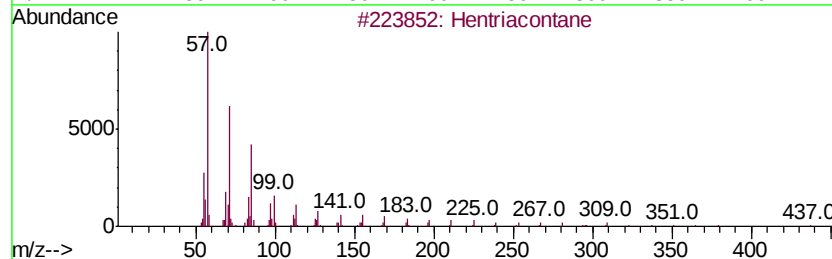
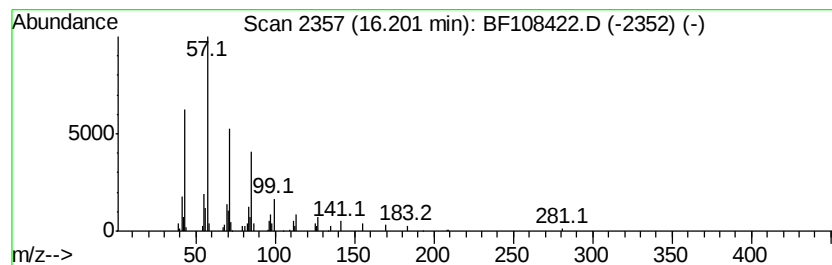
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Peak Number 5 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	2.16 ng	91588	Perylene-d12	15.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hentriacontane	437	C31H64	000630-04-6	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



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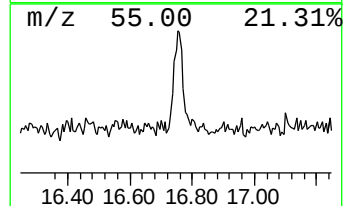
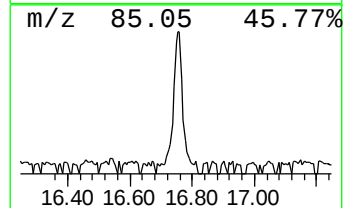
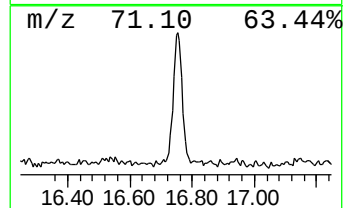
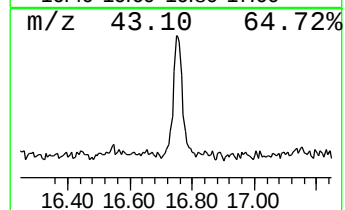
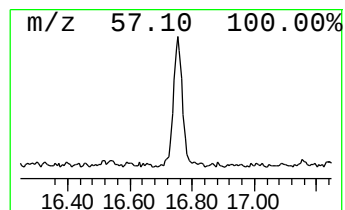
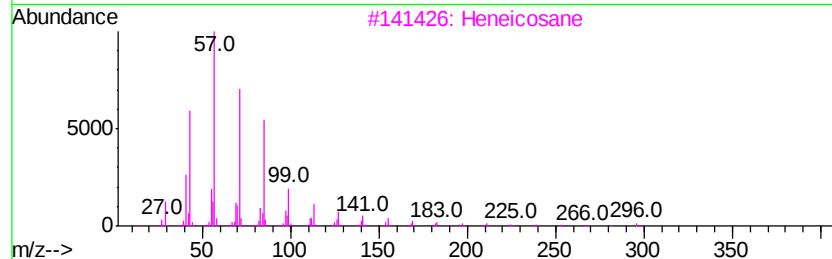
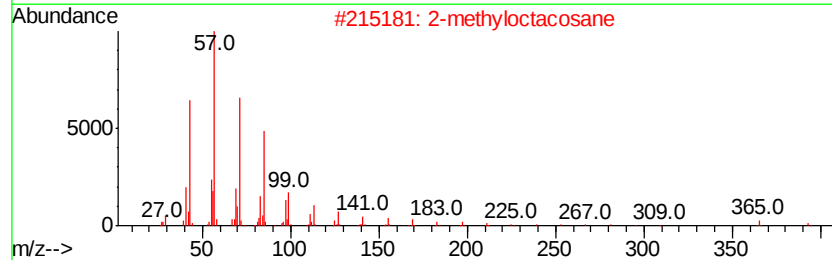
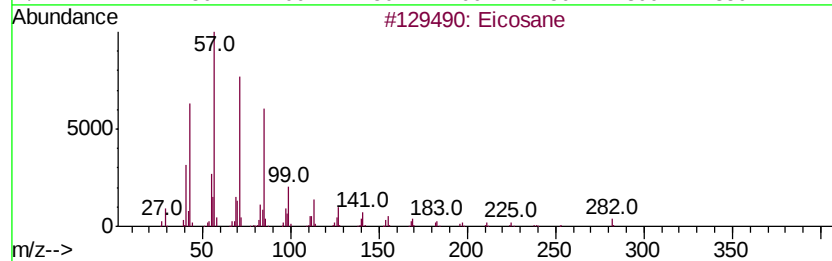
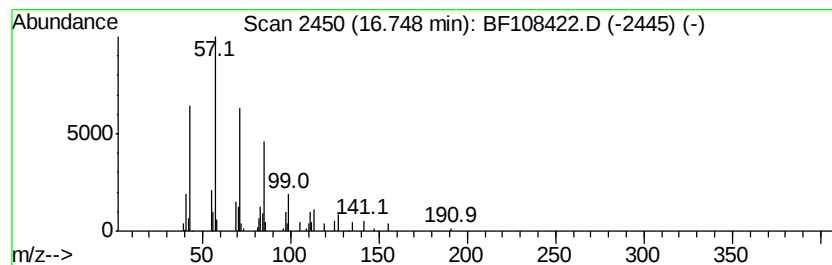
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.75	2.29 ng	97284	Perylene-d12	15.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	95
2	2-methyloctacosane		408	C29H60	1000376-72-8	90
3	Heneicosane		296	C21H44	000629-94-7	87
4	Heptadecane		240	C17H36	000629-78-7	86
5	Hexacosane		366	C26H54	000630-01-3	86



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.45	3.0	ng	142189	1	7.00	949397	20.0
2-Pentanone, 4-hy...	5.27	3.8	ng	178413	1	7.00	949397	20.0
unknown6.77	6.77	85.1	ng	4038890	1	7.00	949397	20.0
Hentriacontane	16.20	2.2	ng	91588	6	15.75	848926	20.0
Eicosane	16.75	2.3	ng	97284	6	15.75	848926	20.0