

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041816\
 Data File : BG021725.D
 Acq On : 18 Apr 2016 5:12
 Operator : UM/SJ
 Sample : H2513-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ACF-NW-2

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:\HPCHEM1\BNA_G\METHODS\8270-BG041316.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	3.098	39	44	66	rVB	1494300	2174923	100.00%	16.493%
2	5.290	411	417	424	rVB	21227	34394	1.58%	0.261%
3	5.766	490	498	507	rBV	433050	725522	33.36%	5.502%
4	7.364	761	770	780	rBV	491988	860909	39.58%	6.528%
5	7.746	826	835	845	rBV	462199	802332	36.89%	6.084%
6	8.210	908	914	921	rBV	102407	178577	8.21%	1.354%
7	8.539	962	970	978	rVB	334697	583224	26.82%	4.423%
8	9.379	1105	1113	1121	rBV	304468	543201	24.98%	4.119%
9	11.030	1384	1394	1405	rVB	167605	297167	13.66%	2.253%
10	13.451	1797	1806	1817	rBV	885873	1390616	63.94%	10.545%
11	14.262	1938	1944	1951	rBV	99010	142777	6.56%	1.083%
12	14.691	2011	2017	2023	rBV	26187	34486	1.59%	0.262%
13	14.820	2033	2039	2047	rVB2	287496	473140	21.75%	3.588%
14	15.637	2173	2178	2183	rVB	45683	59083	2.72%	0.448%
15	16.042	2238	2247	2251	rBV4	13050	30262	1.39%	0.229%
16	16.301	2285	2291	2302	rVB	669824	1006006	46.25%	7.629%
17	16.494	2319	2324	2332	rBV	44422	75723	3.48%	0.574%
18	17.282	2454	2458	2462	rBV	21184	25945	1.19%	0.197%
19	17.558	2496	2505	2512	rBV2	333538	534781	24.59%	4.055%
20	20.143	2939	2945	2951	rBV	1408405	1898630	87.30%	14.398%
21	21.436	3161	3165	3174	rBV5	15101	27526	1.27%	0.209%
22	21.829	3225	3232	3240	rVB	354576	616043	28.32%	4.672%
23	23.539	3518	3523	3531	rVB	30312	60788	2.79%	0.461%
24	25.155	3787	3798	3814	rVB3	201651	611007	28.09%	4.633%

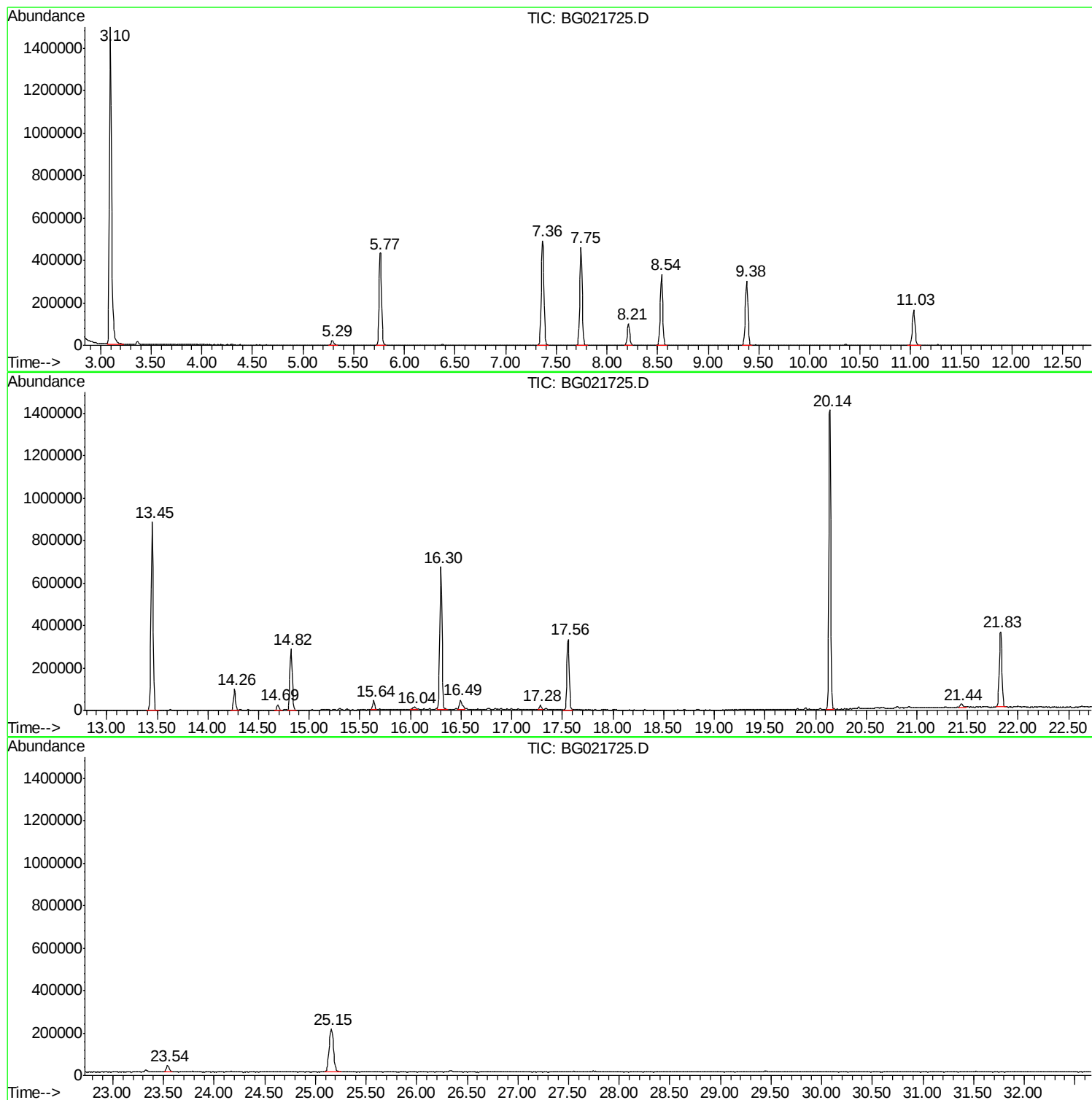
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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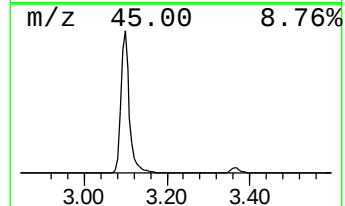
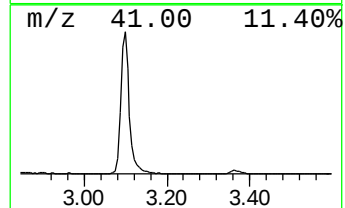
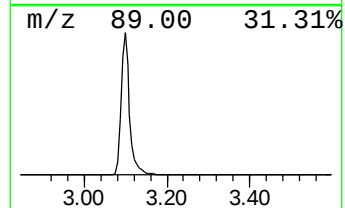
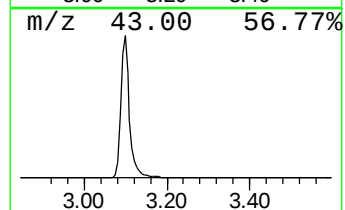
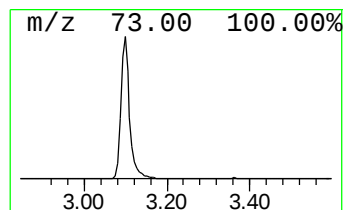
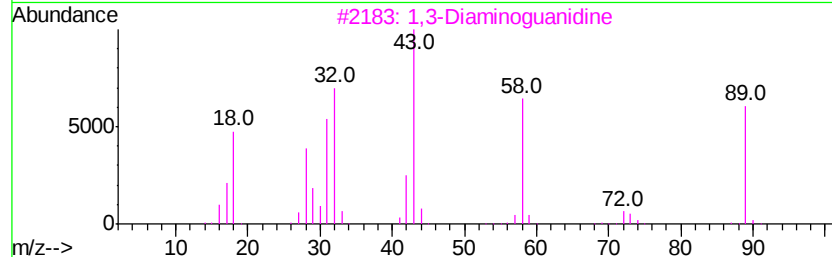
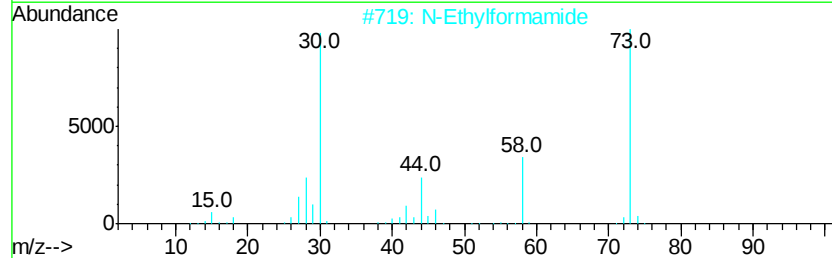
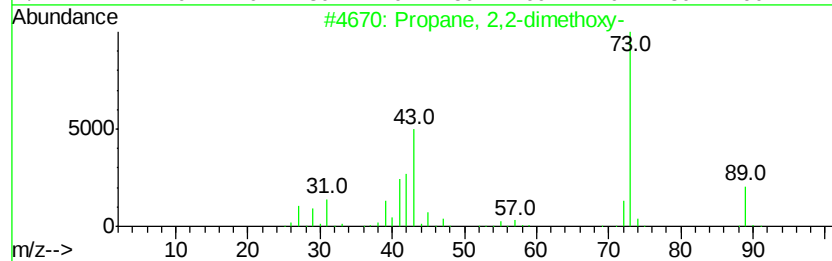
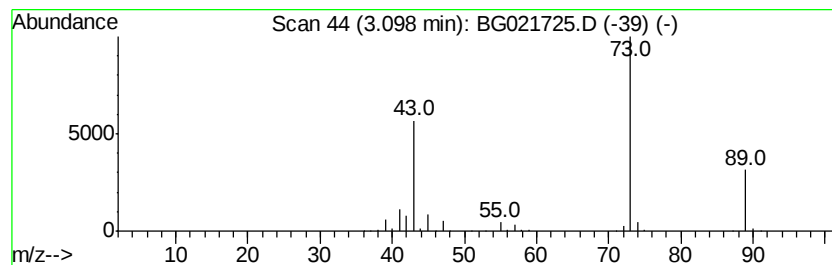
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	243.58 ng	2174920	1,4-Dichlorobenzene-d4	8.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2			N-Ethylformamide	73	C3H7NO	000627-45-2	9
3			1,3-Diaminoquanidine	89	CH7N5	004364-78-7	9
4			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5			1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9



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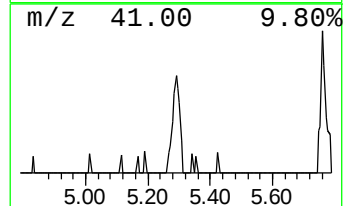
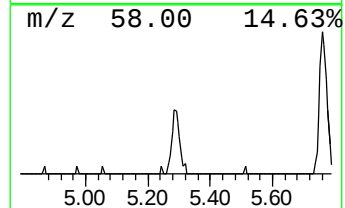
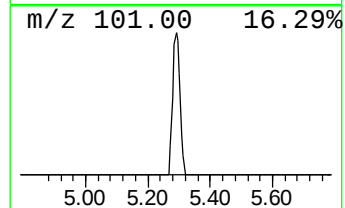
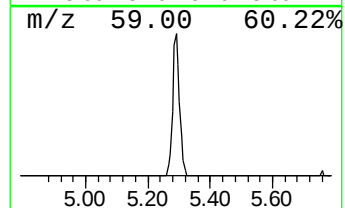
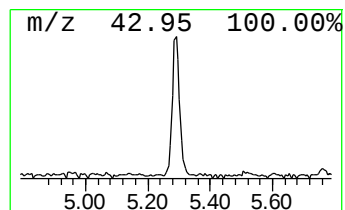
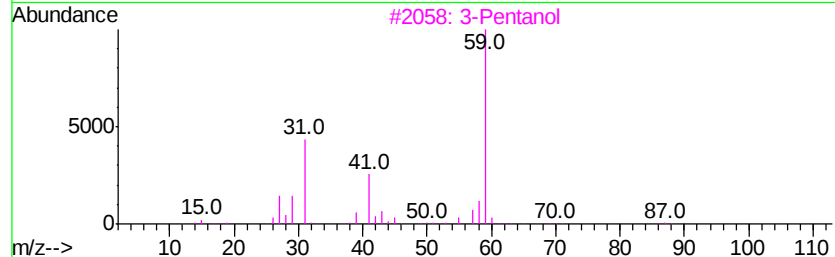
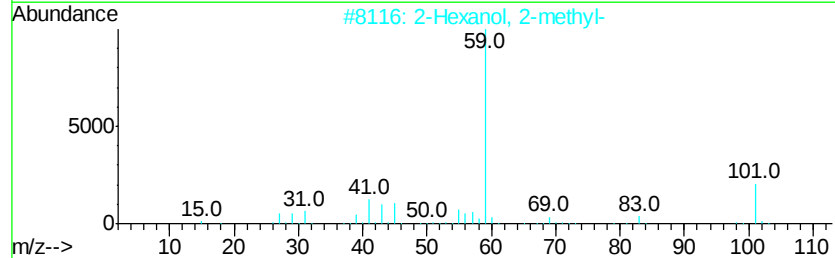
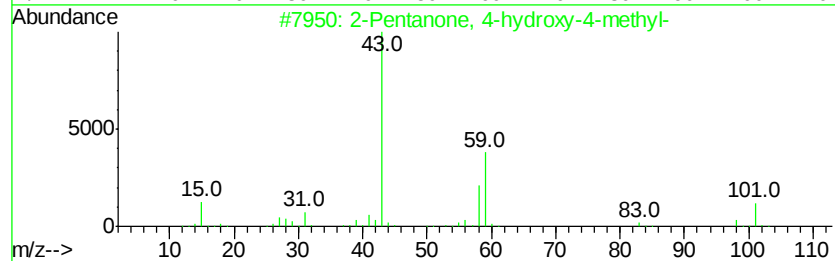
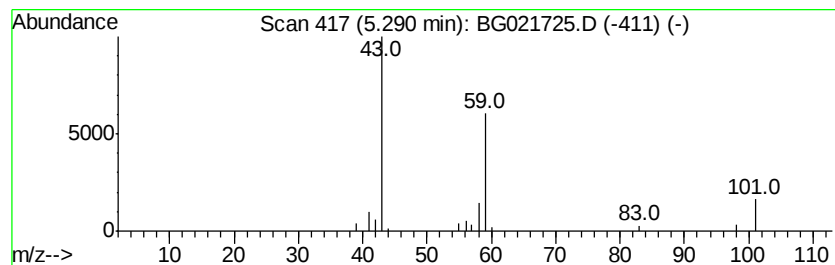
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.29	3.85 ng	34394	1,4-Dichlorobenzene-d4	8.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	47
3		3-Pentanol	88	C5H12O	000584-02-1	9
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9



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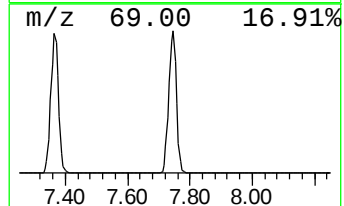
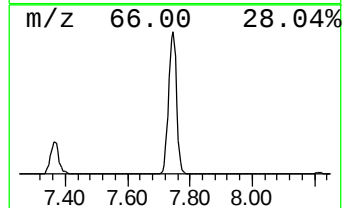
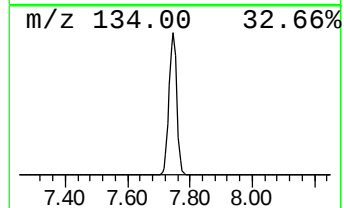
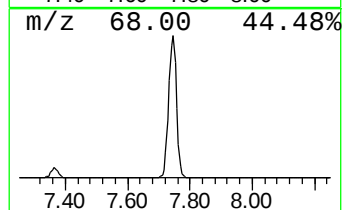
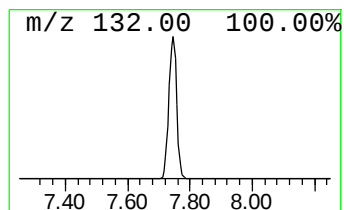
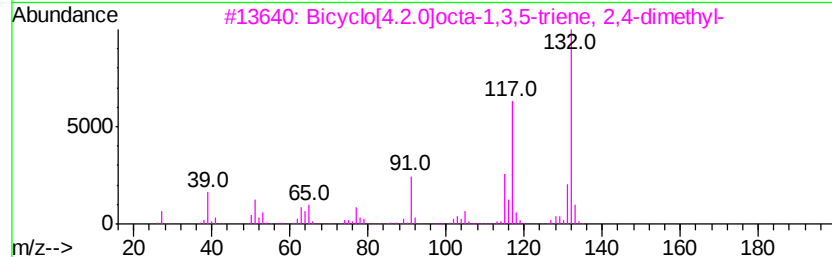
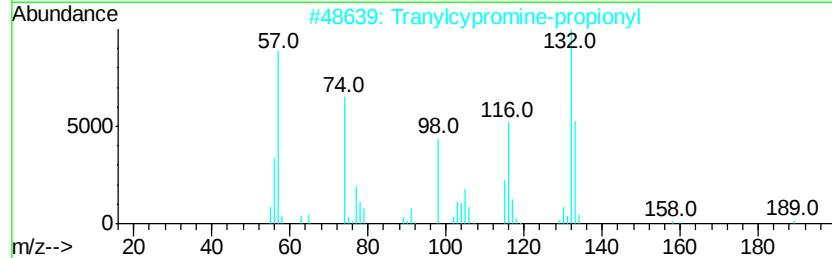
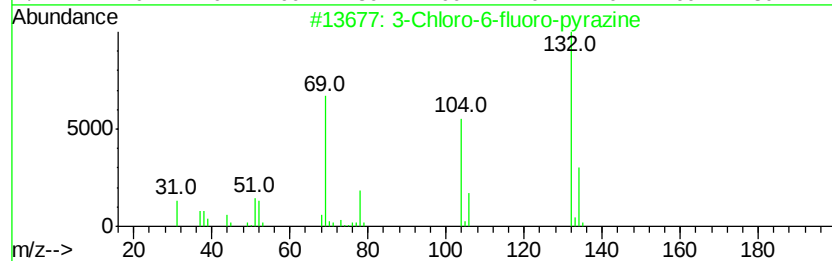
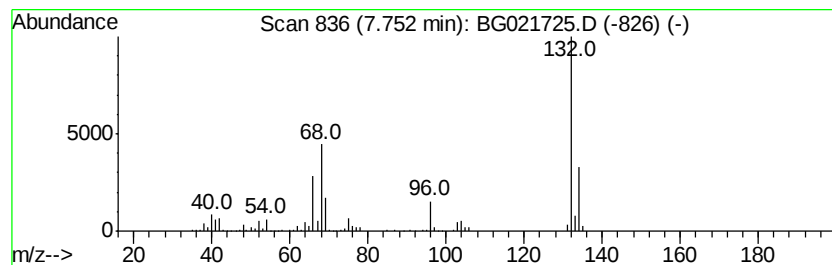
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Peak Number 3 unknown7.75 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	89.86 ng	802332	1,4-Dichlorobenzene-d4	8.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		1H-Tetrazaborole, 5-chloro-4,5-d...	132	C2H6BClN4	021960-49-6	9



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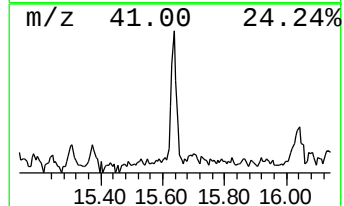
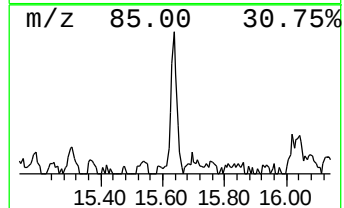
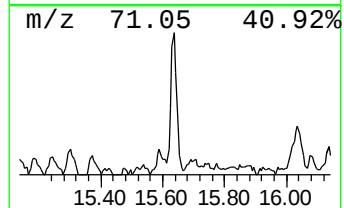
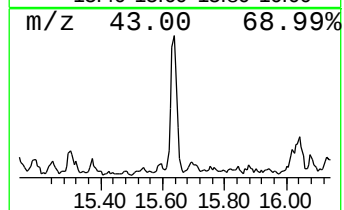
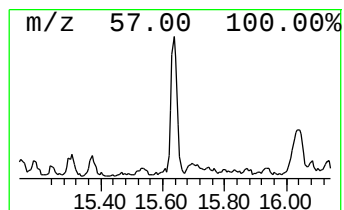
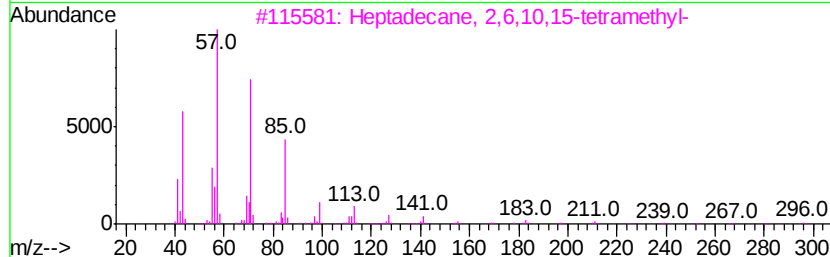
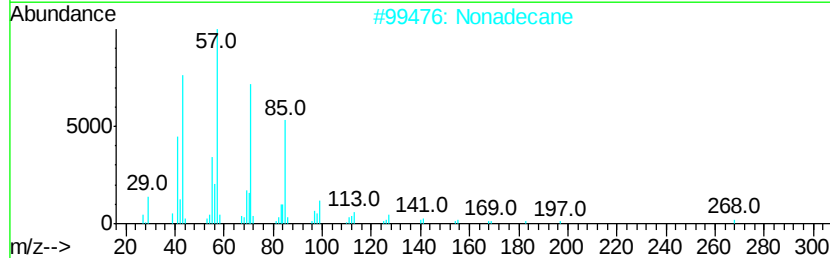
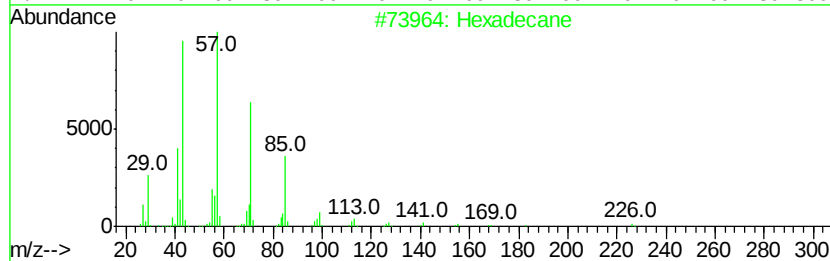
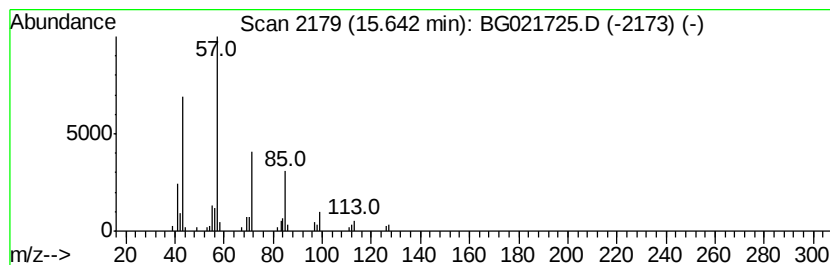
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Peak Number 5 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	2.50 ng	59083	Acenaphthene-d10	14.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	96
2	Nonadecane	268	C19H40	000629-92-5	86
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	83
4	Eicosane	282	C20H42	000112-95-8	80
5	Tridecane	184	C13H28	000629-50-5	80



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-dime...	3.10	243.6	ng	2174920	1	8.21	178577	20.0
2-Pentanone, 4-hy...	5.29	3.9	ng	34394	1	8.21	178577	20.0
unknown7.75	7.75	89.9	ng	802332	1	8.21	178577	20.0
Hexadecane	15.64	2.5	ng	59083	3	14.82	473140	20.0