

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102367.D
 Acq On : 24 Jan 2018 2:42
 Operator : SJ/JU
 Sample : J1249-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AB-1(0-2)

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_F\METHODS\8270-BF012218.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	4.916	478	481	490	rVB	100938	144052	4.82%	0.580%
2	5.328	546	551	555	rBV	2677141	2816047	94.27%	11.333%
3	6.363	721	727	739	rBV	2655470	2987126	100.00%	12.021%
4	6.487	743	748	751	rBV	3011051	2901218	97.12%	11.676%
5	6.716	783	787	796	rBV	931772	786072	26.32%	3.163%
6	6.869	809	813	817	rBV	2288411	2146250	71.85%	8.637%
7	7.281	878	883	886	rBV	1738463	1793205	60.03%	7.217%
8	7.998	1001	1005	1008	rBV	1231548	1030844	34.51%	4.148%
9	8.557	1096	1100	1104	rBV	59675	50651	1.70%	0.204%
10	9.075	1183	1188	1191	rBV	3180858	2835119	94.91%	11.410%
11	9.469	1251	1255	1263	rBV	362485	307658	10.30%	1.238%
12	9.751	1299	1303	1311	rBV	1262616	1084406	36.30%	4.364%
13	10.463	1419	1424	1431	rBV	147671	131539	4.40%	0.529%
14	10.539	1433	1437	1447	rVB	1533417	1415140	47.37%	5.695%
15	11.233	1551	1555	1563	rBV	1030060	942046	31.54%	3.791%
16	12.639	1789	1794	1798	rBV	60544	56040	1.88%	0.226%
17	12.822	1821	1825	1828	rBV	2012845	1768005	59.19%	7.115%
18	13.345	1911	1914	1918	rBV	38014	35801	1.20%	0.144%
19	13.716	1973	1977	1981	rBV	152937	163387	5.47%	0.658%
20	13.874	1996	2004	2011	rBV	780372	761739	25.50%	3.066%
21	15.298	2241	2246	2253	rBV	521384	650139	21.76%	2.616%
22	16.198	2395	2399	2405	rBV3	24801	42185	1.41%	0.170%

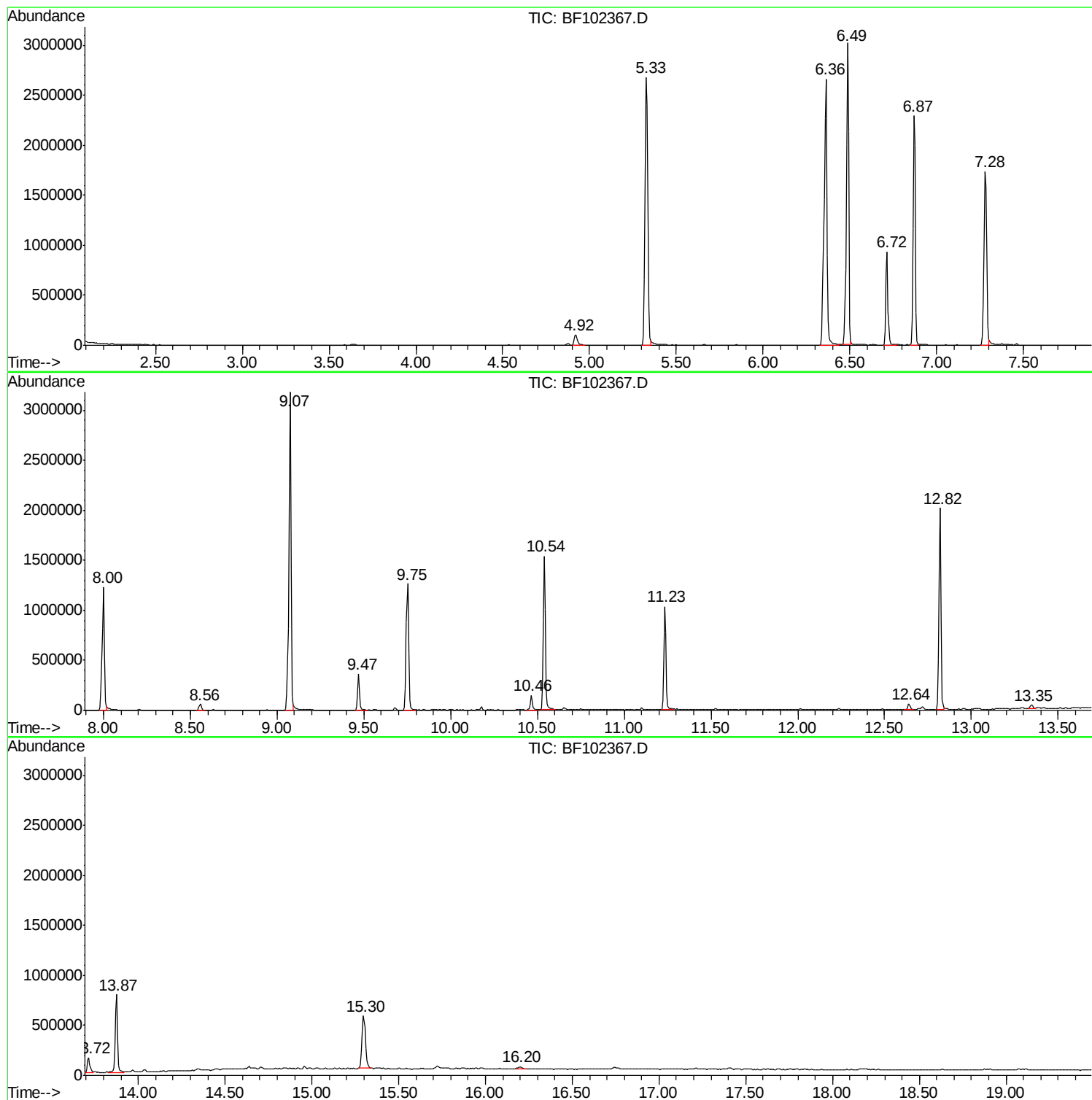
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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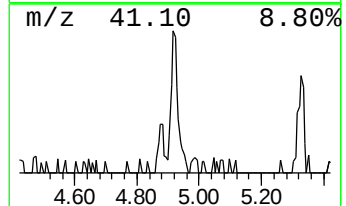
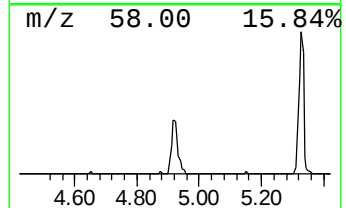
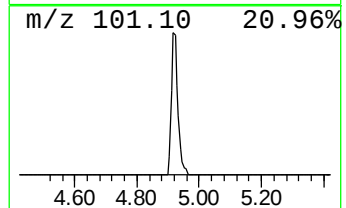
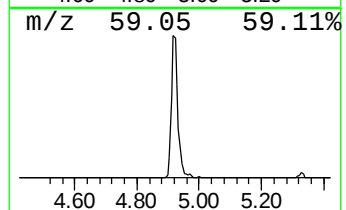
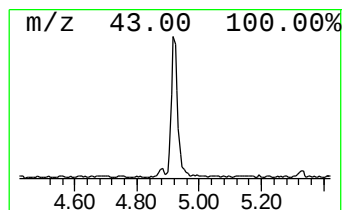
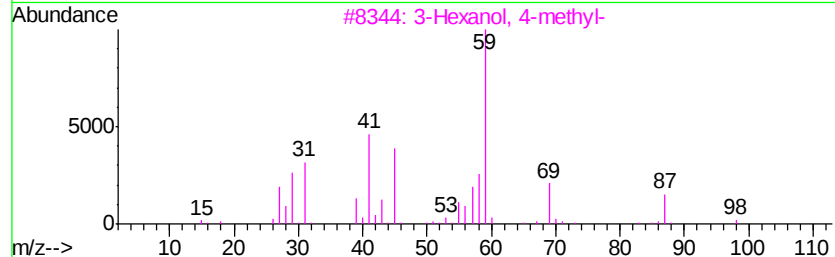
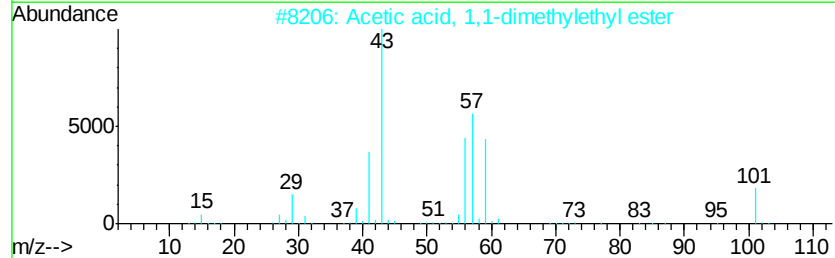
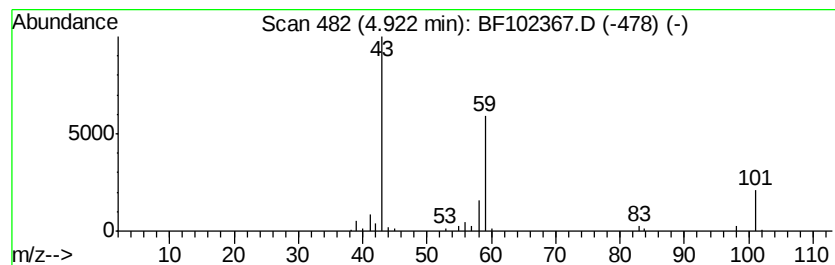
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	3.67 ng	144052	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23



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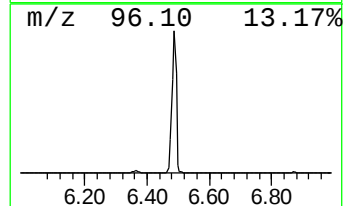
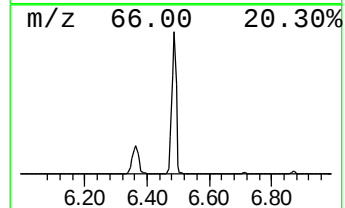
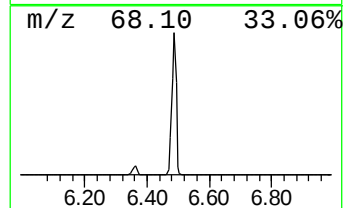
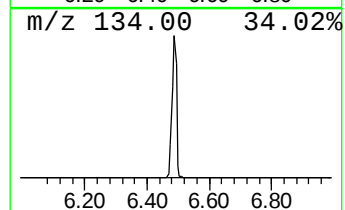
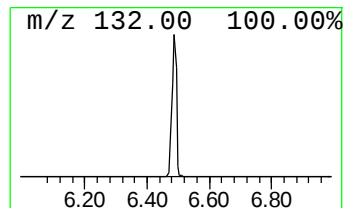
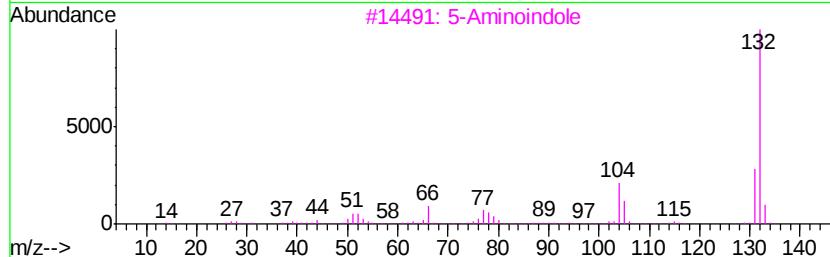
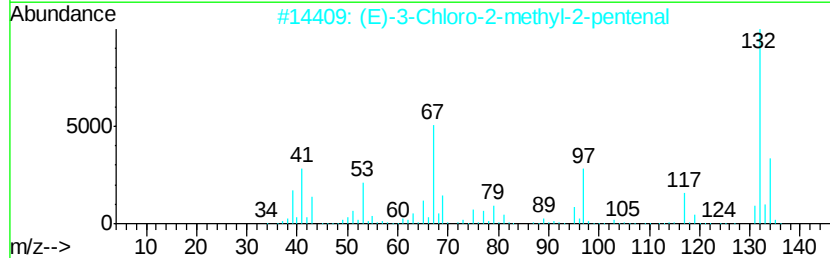
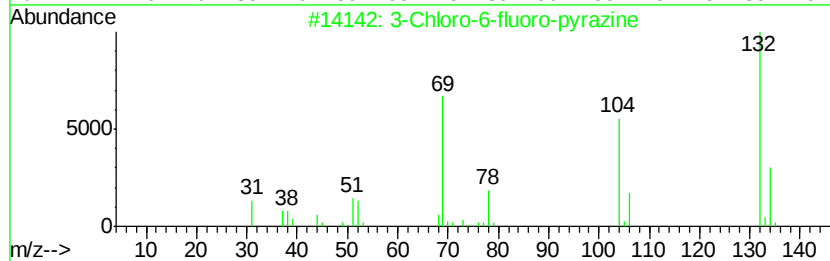
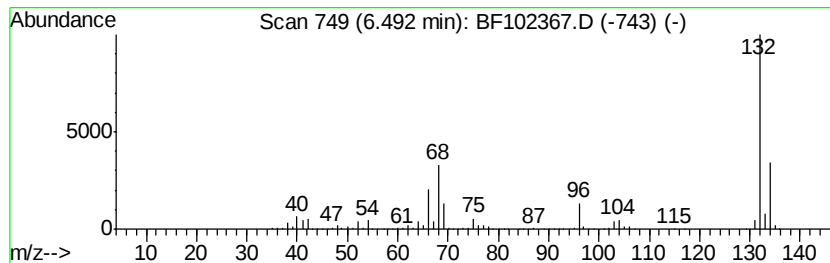
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Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	73.82 ng	2901220	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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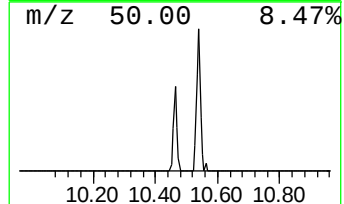
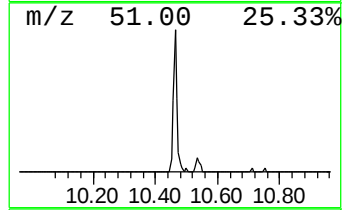
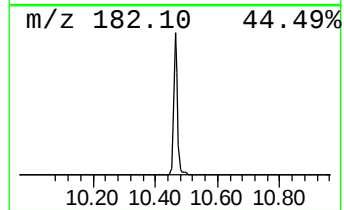
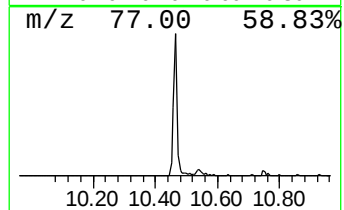
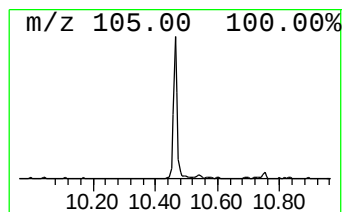
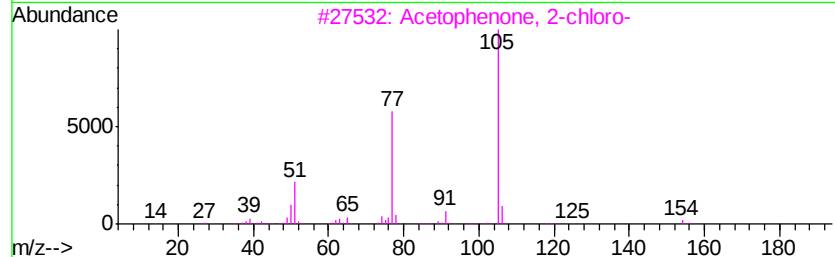
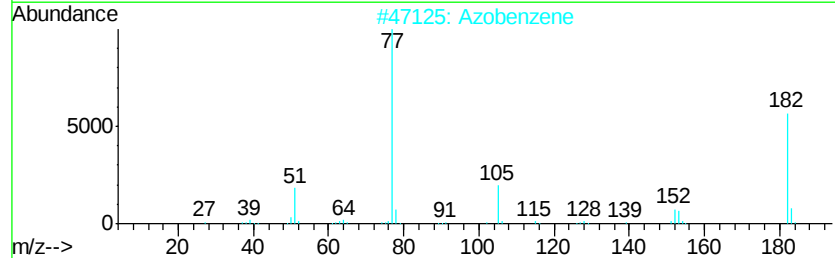
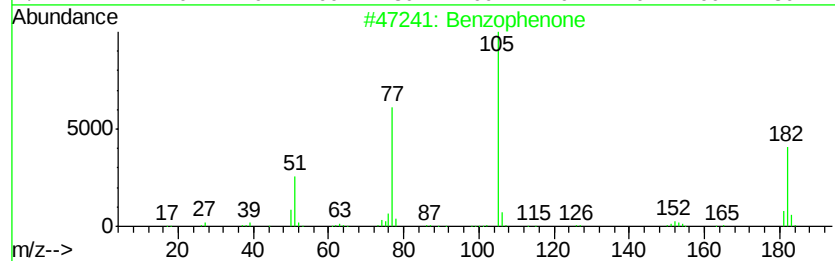
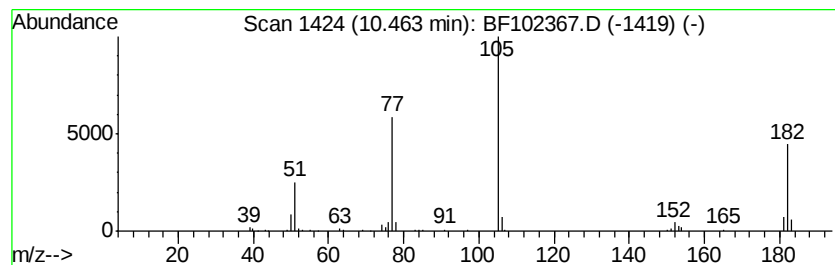
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Peak Number 4 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.46	2.43 ng	131539	Acenaphthene-d10	9.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		Azobenzene	182	C12H10N2	000103-33-3	64
3		Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
4		Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	47
5		N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	47



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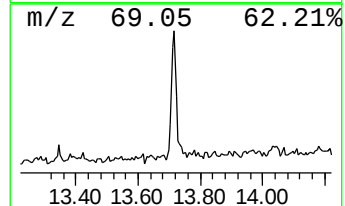
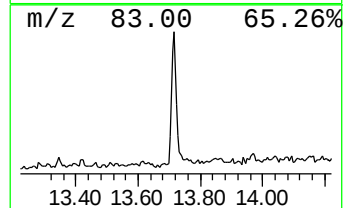
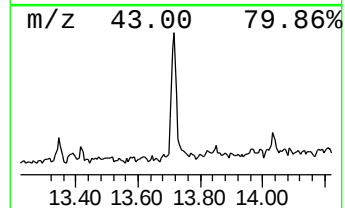
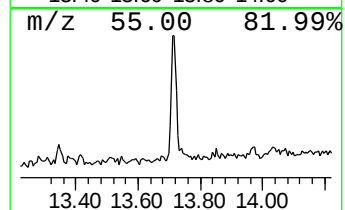
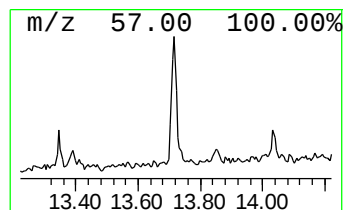
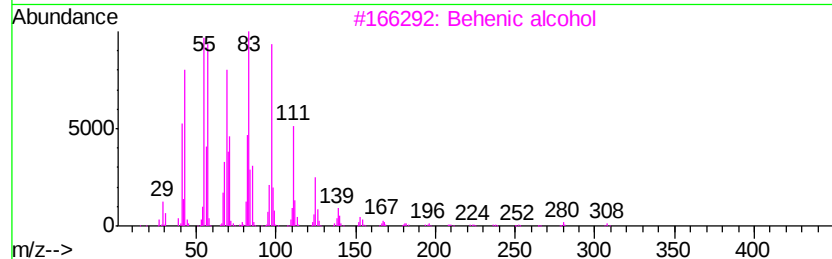
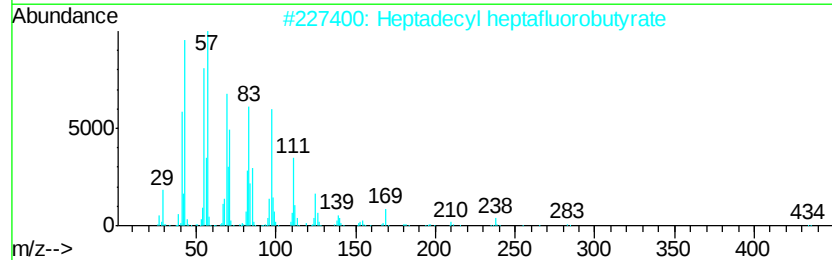
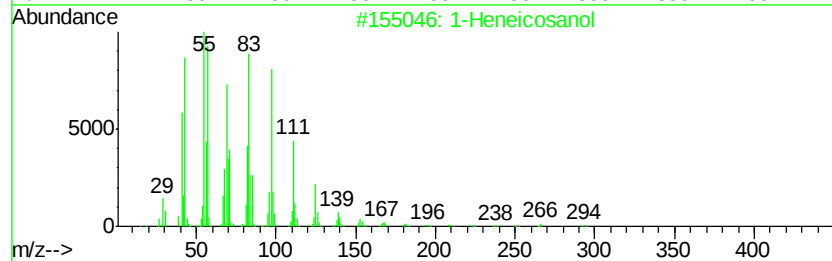
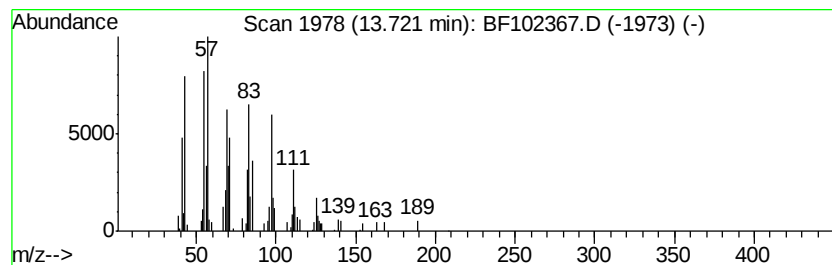
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Peak Number 5 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	4.29 ng	163387	Chrysene-d12	13.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	95
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	95
3		Behenic alcohol	326	C22H46O	000661-19-8	95
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
5		Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.92	3.7	ng	144052	1	6.72	786072	20.0
unknown6.49	6.49	73.8	ng	2901220	1	6.72	786072	20.0
Benzophenone	10.46	2.4	ng	131539	3	9.75	1084410	20.0
1-Heneicosanol	13.72	4.3	ng	163387	5	13.87	761739	20.0