

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081318\
 Data File : BF108125.D
 Acq On : 13 Aug 2018 19:21
 Operator : JU/SJ
 Sample : J4442-41
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1-5

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	3.643	220	222	229	rBV	55631	96991	1.94%	0.226%
2	5.266	493	498	508	rVB	755371	989083	19.82%	2.306%
3	5.648	558	563	567	rBV	4566619	4782972	95.85%	11.153%
4	6.642	726	732	735	rBV	4842242	4989946	100.00%	11.635%
5	6.666	735	736	744	rVB	64804	56568	1.13%	0.132%
6	6.795	753	758	761	rBV	5333693	4877299	97.74%	11.373%
7	7.025	793	797	801	rVB	1568595	1234430	24.74%	2.878%
8	7.183	819	824	827	rBV	4274724	3663757	73.42%	8.543%
9	7.583	887	892	895	rBV	3150945	3036374	60.85%	7.080%
10	8.307	1011	1015	1019	rBV	1757145	1422420	28.51%	3.317%
11	9.383	1193	1198	1201	rBV	5453477	4608965	92.37%	10.747%
12	9.460	1207	1211	1214	rBV3	53208	55670	1.12%	0.130%
13	9.772	1261	1264	1267	rBV	497679	385437	7.72%	0.899%
14	10.072	1309	1315	1319	rBV2	1536359	1409952	28.26%	3.288%
15	10.483	1382	1385	1388	rVB	65908	57891	1.16%	0.135%
16	10.707	1419	1423	1425	rVB	52280	52767	1.06%	0.123%
17	10.871	1446	1451	1454	rBV	2723390	2686199	53.83%	6.263%
18	10.971	1463	1468	1471	rVB	118399	148110	2.97%	0.345%
19	11.442	1545	1548	1551	rBV2	59030	59100	1.18%	0.138%
20	11.571	1566	1570	1574	rBV	1598040	1394620	27.95%	3.252%
21	13.165	1836	1841	1844	rBV	4656381	4343787	87.05%	10.129%
22	14.054	1989	1992	2004	rBV2	195668	397274	7.96%	0.926%
23	14.230	2019	2022	2026	rBV	1237510	1193578	23.92%	2.783%
24	15.806	2286	2290	2296	rVB	638170	943471	18.91%	2.200%

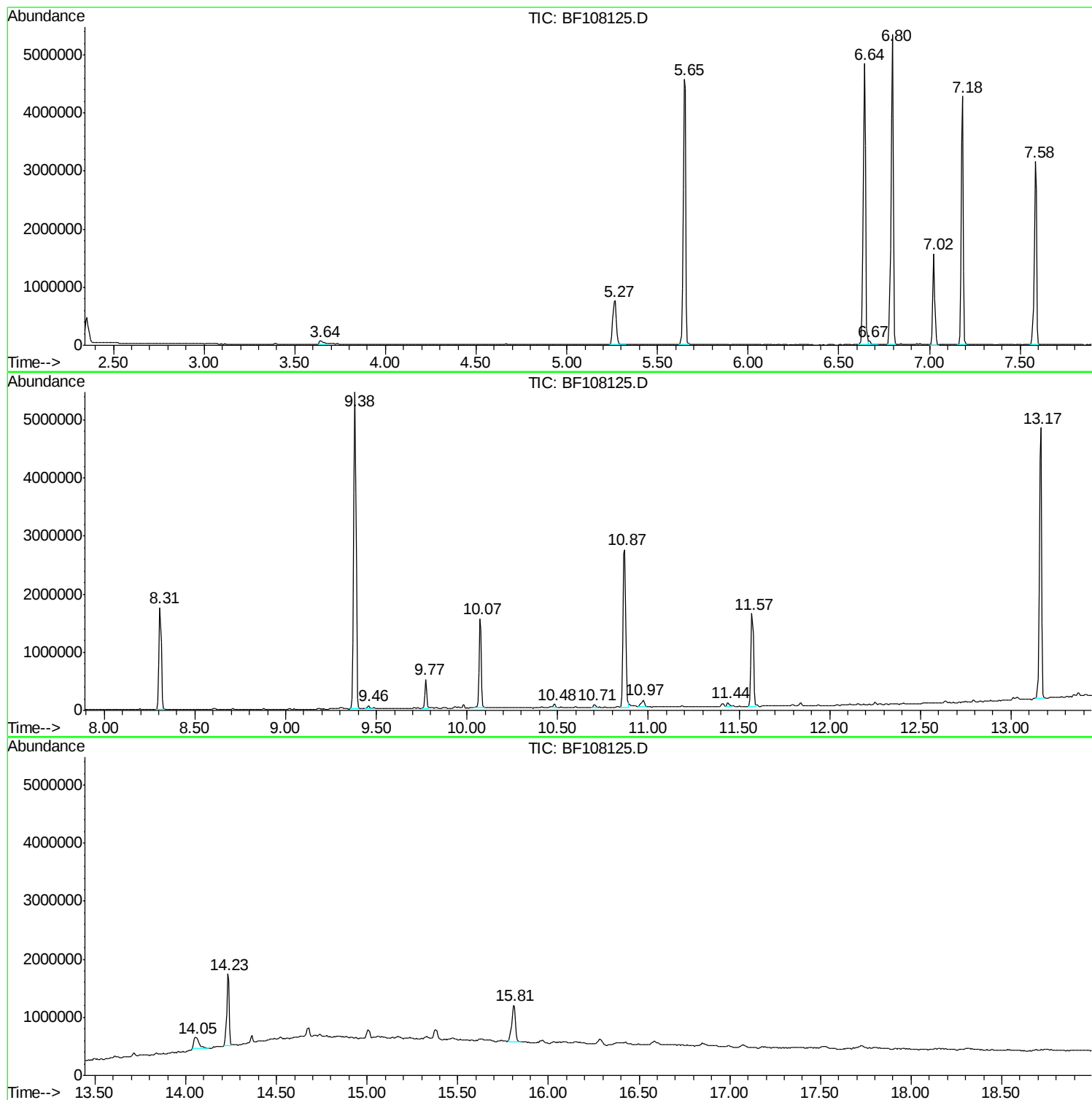
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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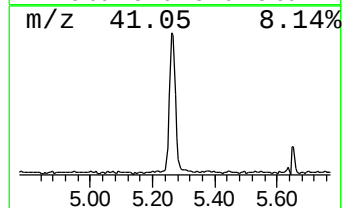
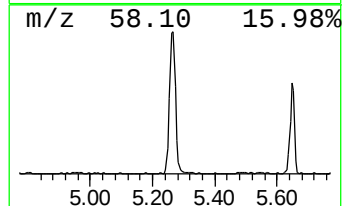
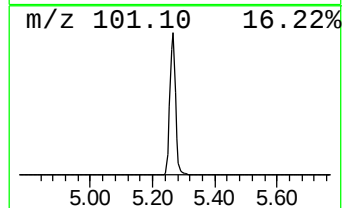
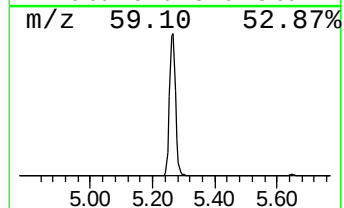
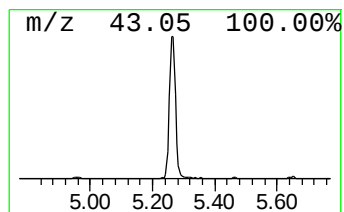
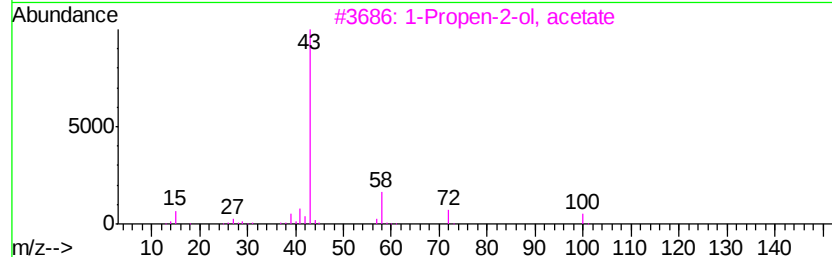
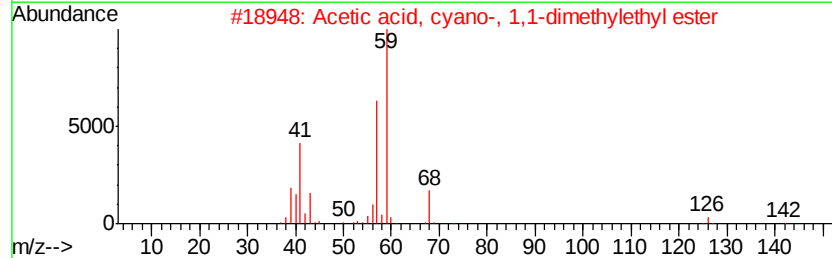
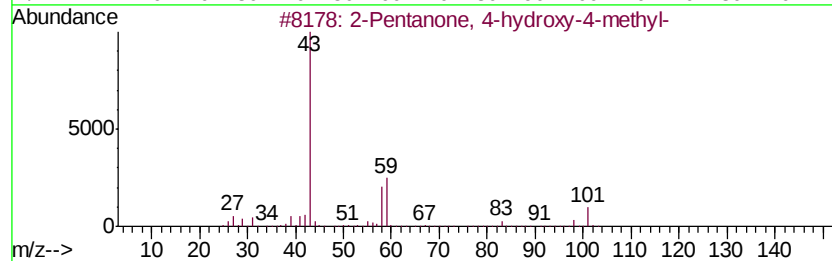
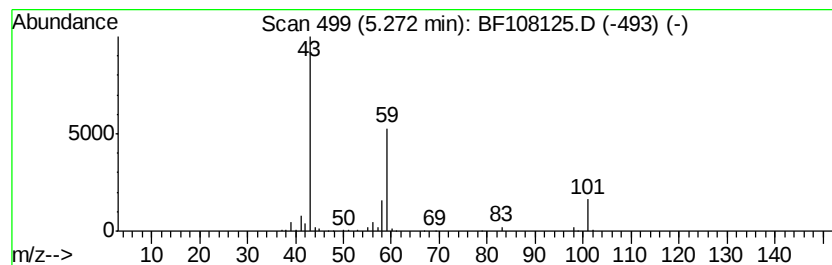
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.27	16.02 ng	989083	1,4-Dichlorobenzene-d4	7.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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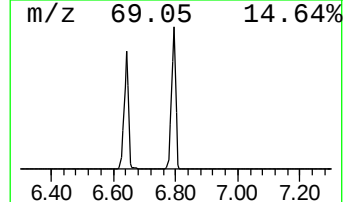
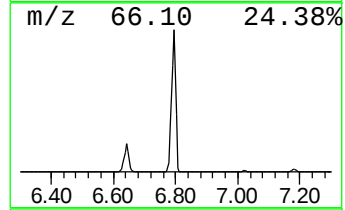
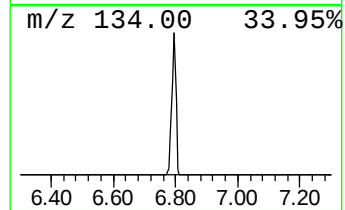
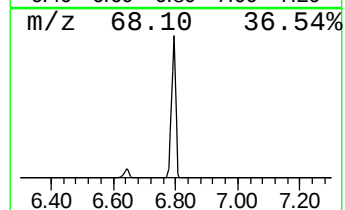
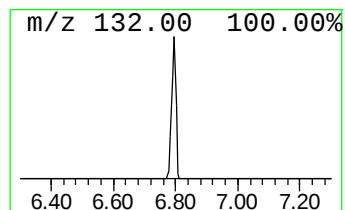
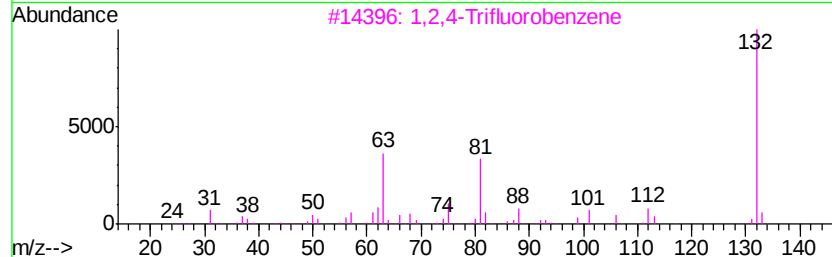
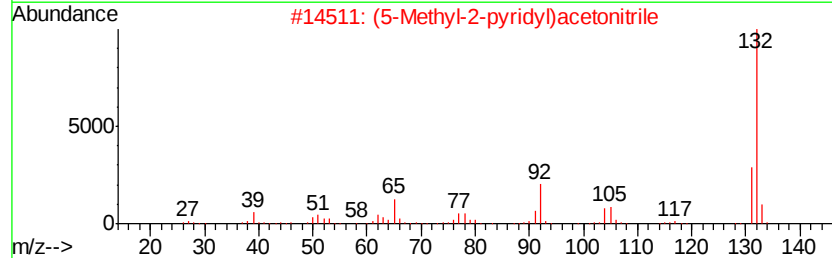
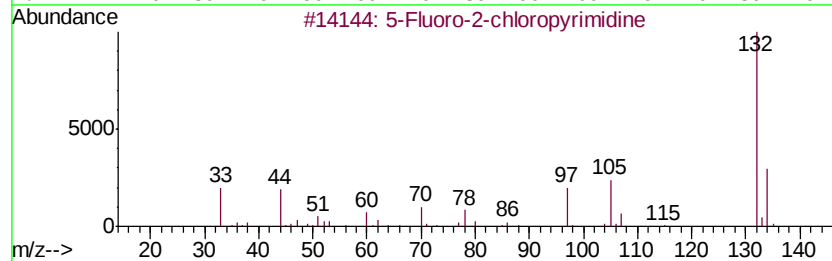
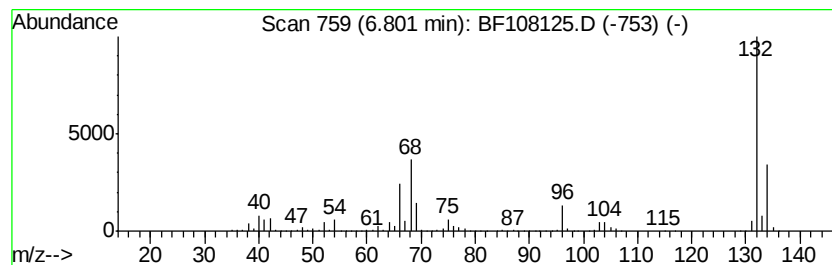
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Peak Number 2 unknown6.80 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	79.02 ng	4877300	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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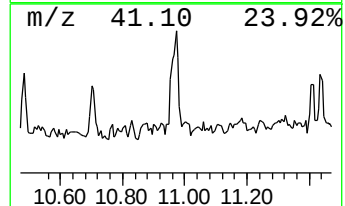
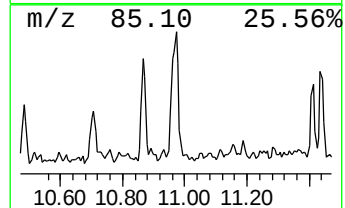
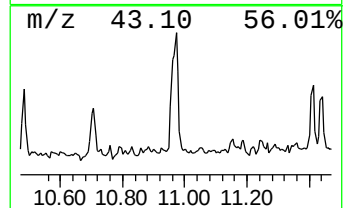
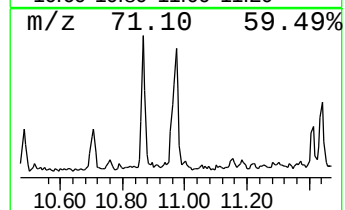
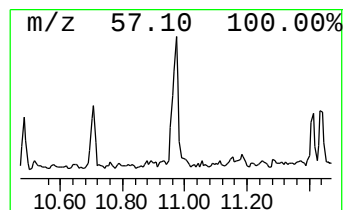
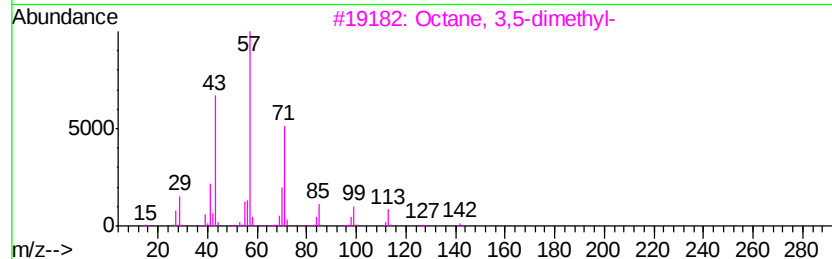
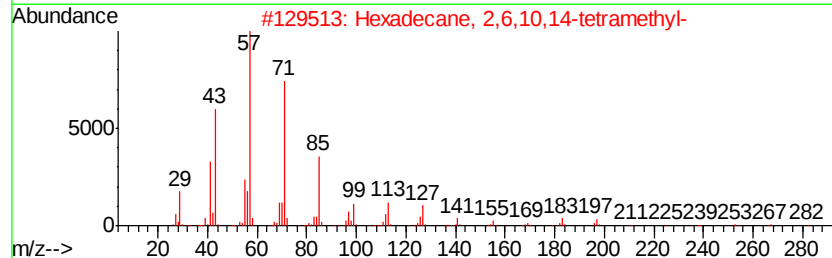
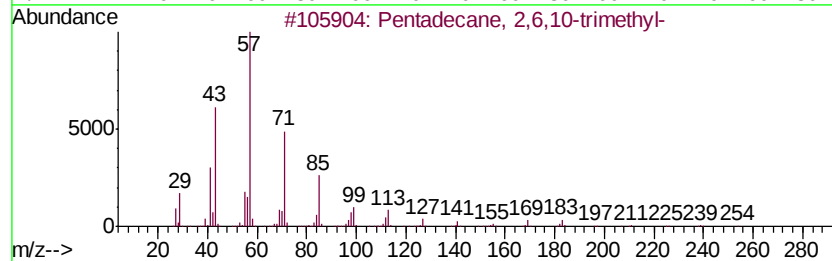
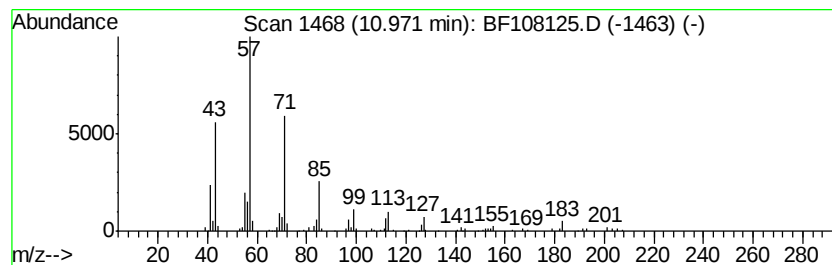
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Peak Number 4 Pentadecane, 2,6,10-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	2.12 ng	148110	Phenanthrene-d10	11.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	80
3		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	68
4		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	64
5		Heptadecane, 7-methyl-	254	C18H38	020959-33-5	64



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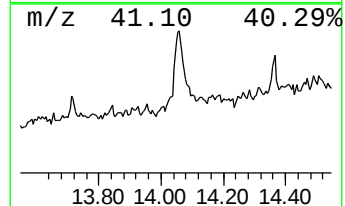
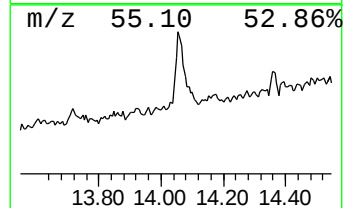
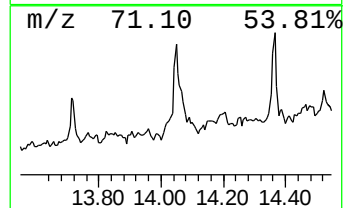
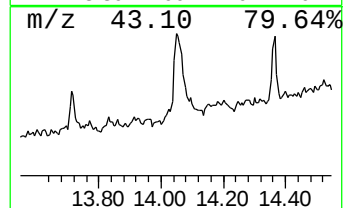
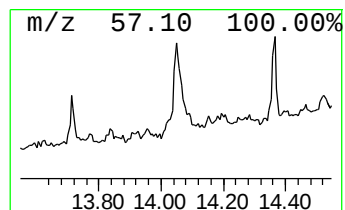
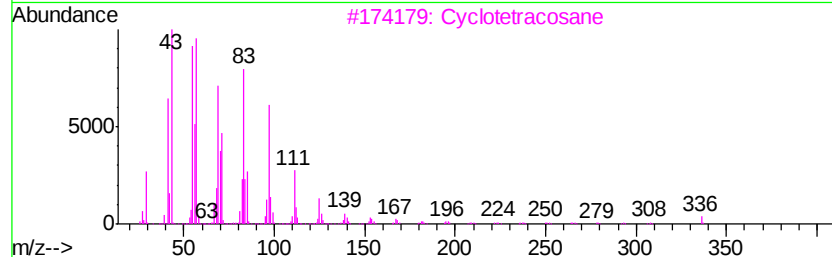
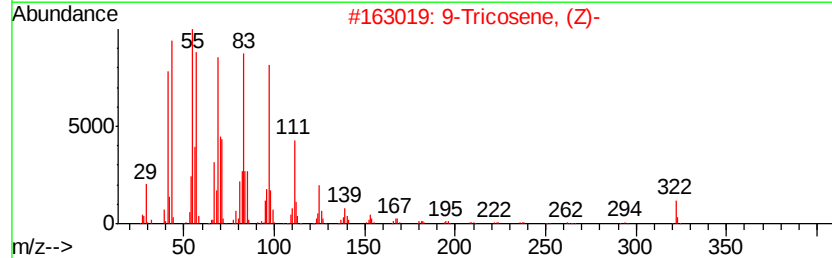
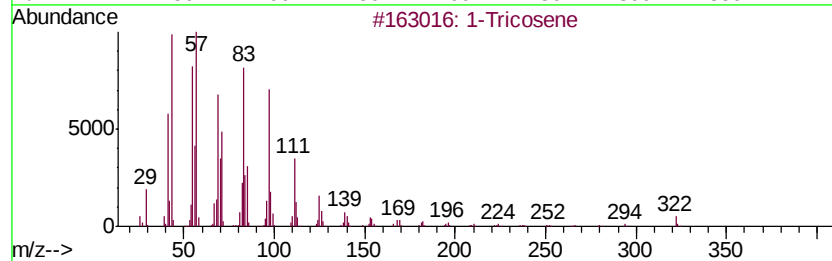
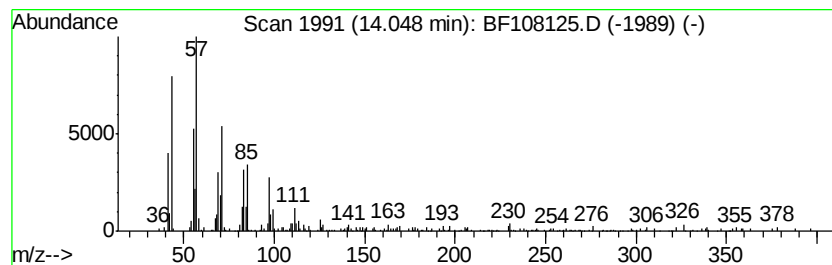
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Peak Number 5 1-Tricosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.05	6.66 ng	397274	Chrysene-d12	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Tricosene		322 C23H46	018835-32-0 97
2	9-Tricosene, (Z)-		322 C23H46	027519-02-4 96
3	Cyclotetracosane		336 C24H48	000297-03-0 94
4	3-Chloropropionic acid, heptadec...		346 C20H39ClO2	1000283-05-1 90
5	Cycloeicosane		280 C20H40	000296-56-0 90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.27	16.0	ng	989083	1	7.02	1234430	20.0
unknown6.80	6.80	79.0	ng	4877300	1	7.02	1234430	20.0
Pentadecane, 2,6,...	10.97	2.1	ng	148110	4	11.57	1394620	20.0
1-Tricosene	14.05	6.7	ng	397274	5	14.23	1193580	20.0