

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121824\
 Data File : BF140906.D
 Acq On : 18 Dec 2024 15:28
 Operator : RC/JU
 Sample : P5277-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-3A

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-BF121624.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF140906.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.063	500	505	520	rVB	36472	57486	4.42%	0.719%
2	5.487	572	577	604	rBV	380023	621686	47.81%	7.778%
3	6.498	744	749	779	rBV	448231	730514	56.18%	9.139%
4	6.845	803	808	816	rBV	311104	389574	29.96%	4.874%
5	7.410	899	904	923	rBV	353000	484364	37.25%	6.060%
6	8.128	1021	1026	1043	rBV	396673	509923	39.22%	6.379%
7	9.198	1203	1208	1218	rBV	760973	972636	74.80%	12.168%
8	9.880	1319	1324	1345	rVB	415817	553151	42.54%	6.920%
9	10.598	1441	1446	1452	rBV	48414	65399	5.03%	0.818%
10	10.680	1455	1460	1474	rBV	301312	449794	34.59%	5.627%
11	11.374	1573	1578	1586	rBV	605479	773702	59.50%	9.679%
12	12.957	1841	1847	1858	rBV	1036430	1300278	100.00%	16.267%
13	13.433	1922	1928	1937	rBV	65412	92256	7.10%	1.154%
14	13.515	1937	1942	1951	rVB5	8528	14449	1.11%	0.181%
15	13.851	1994	1999	2017	rVB2	29238	82085	6.31%	1.027%
16	14.027	2024	2029	2039	rBV	369689	479520	36.88%	5.999%
17	15.127	2213	2216	2222	rVB4	12052	16521	1.27%	0.207%
18	15.527	2278	2284	2301	rVB	238370	399904	30.76%	5.003%

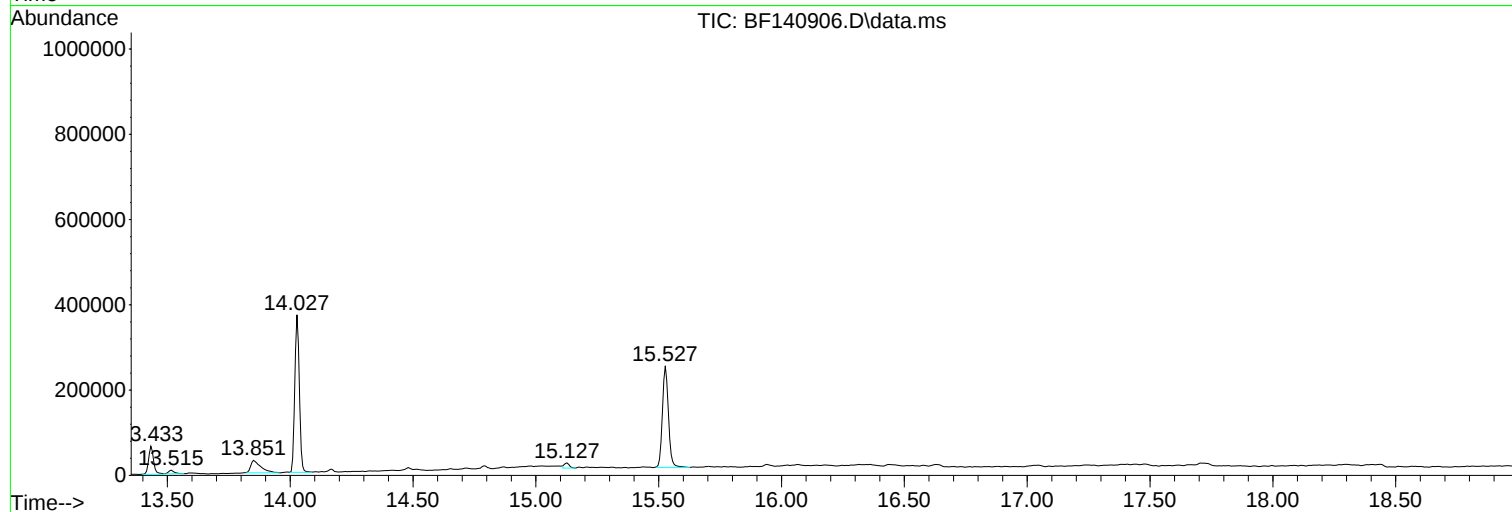
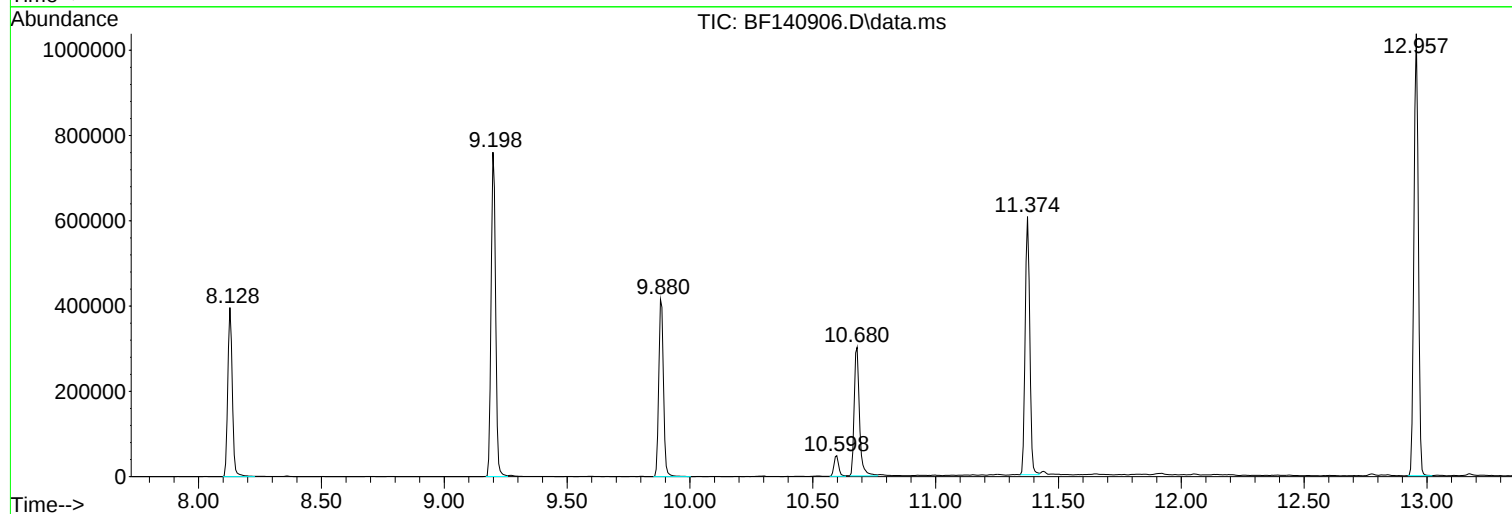
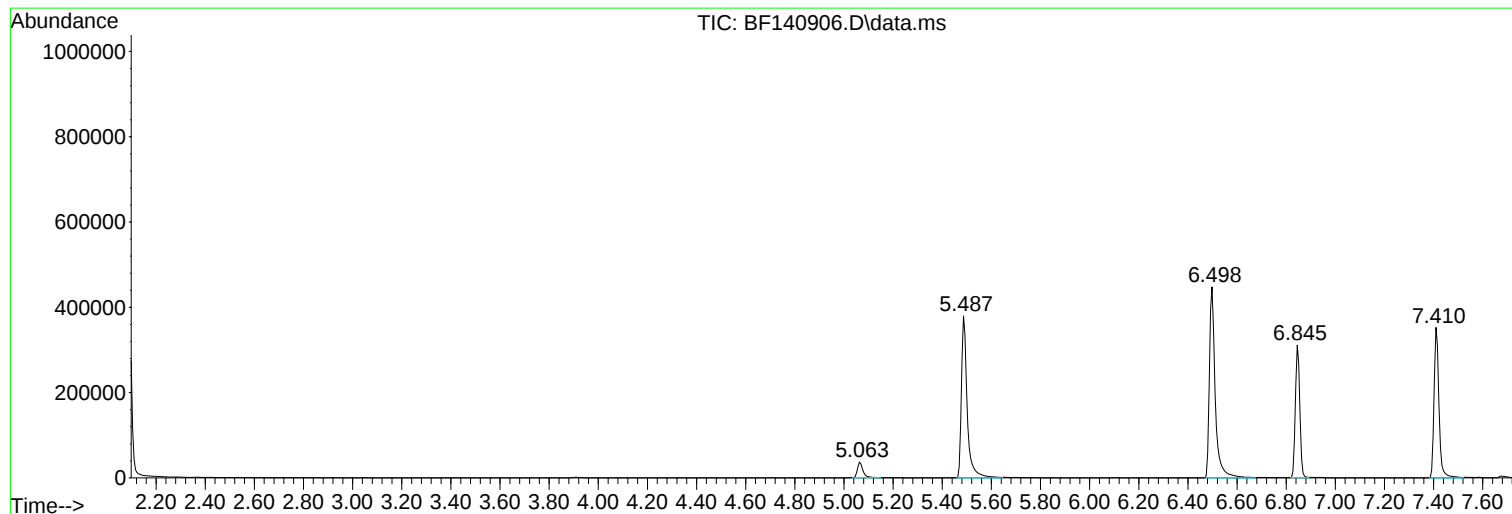
Sum of corrected areas: 7993242

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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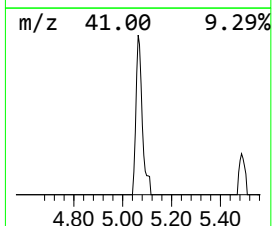
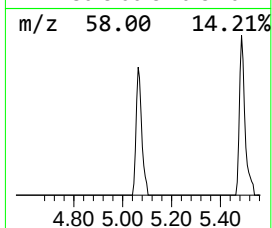
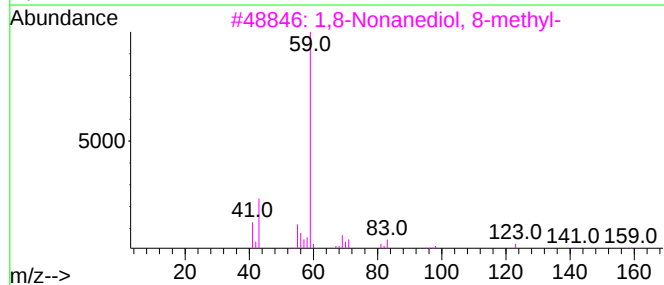
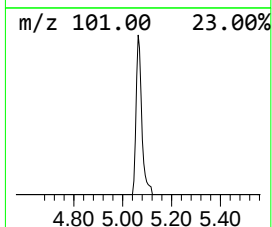
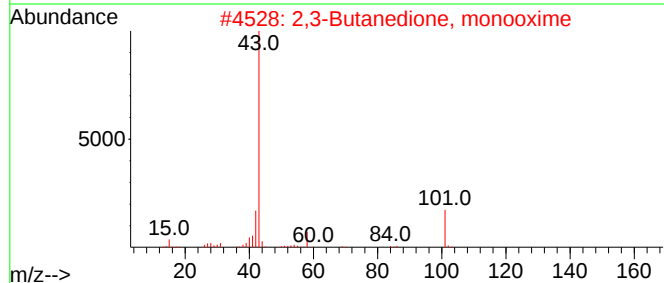
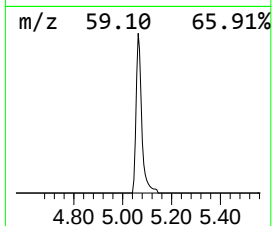
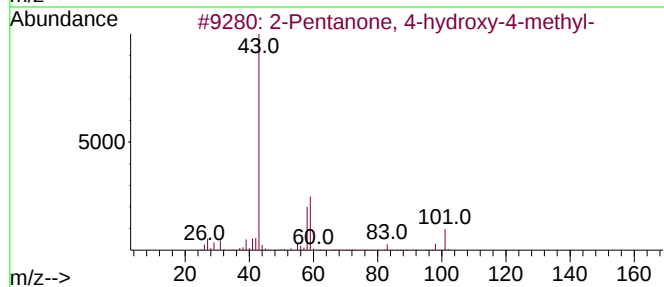
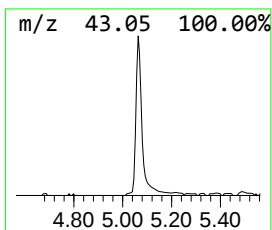
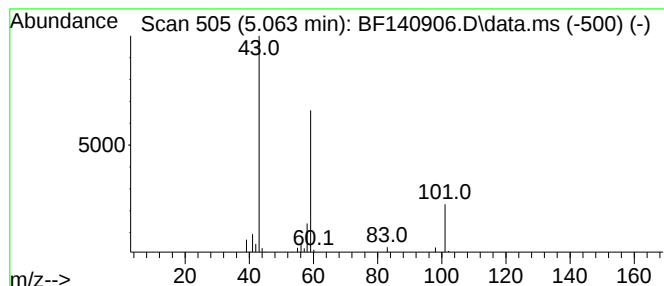
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.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.
5.063	2.95 ng	57486	1,4-Dichlorobenzene-d4	6.845

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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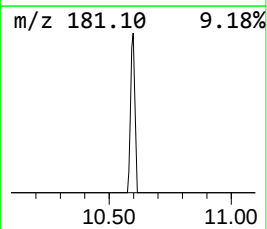
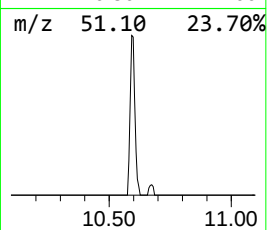
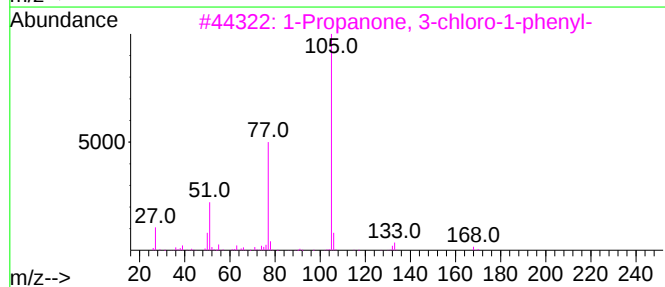
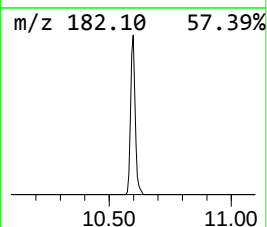
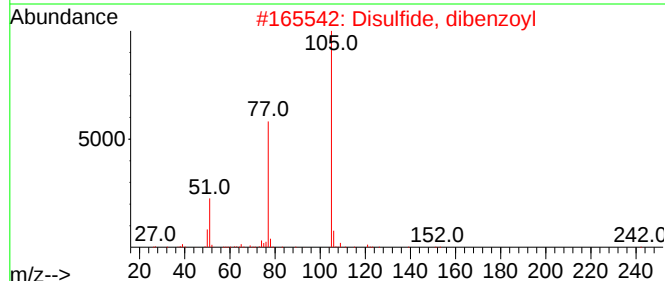
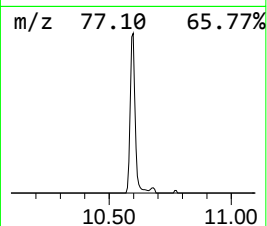
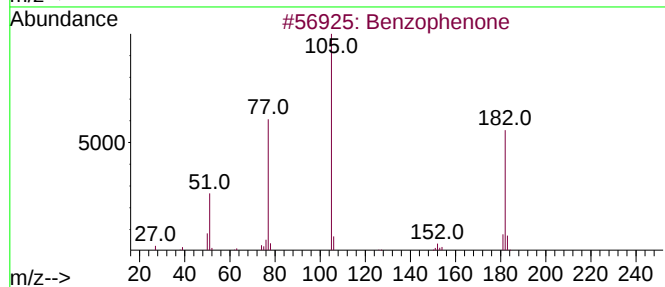
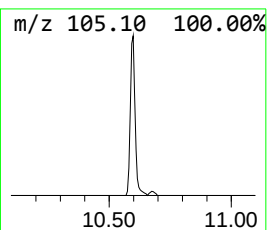
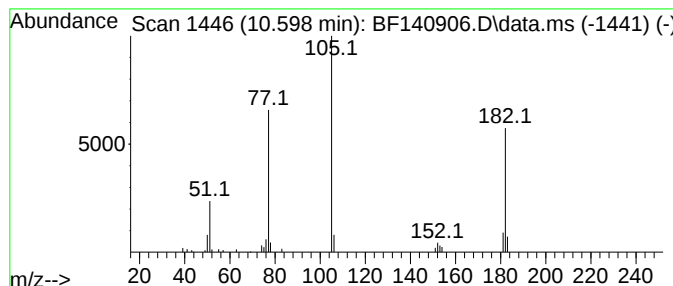
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.598	2.36 ng	65399	Acenaphthene-d10	9.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Disulfide, dibenzoyl	274	C14H10O2S2	000644-32-6	47
3			1-Propanone, 3-chloro-1-phenyl-	168	C9H9ClO	000936-59-4	47
4			.alpha.,.alpha.-Dichloroacetophe...	188	C8H6Cl2O	002648-61-5	47
5			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	47



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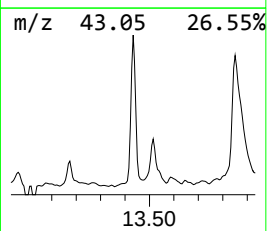
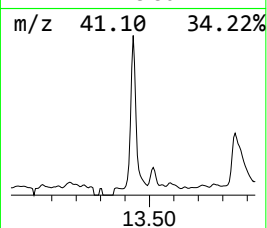
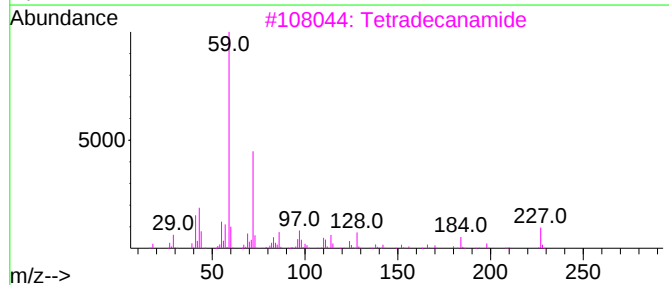
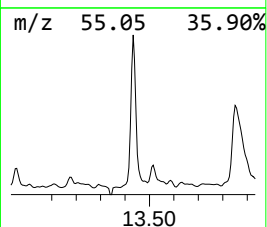
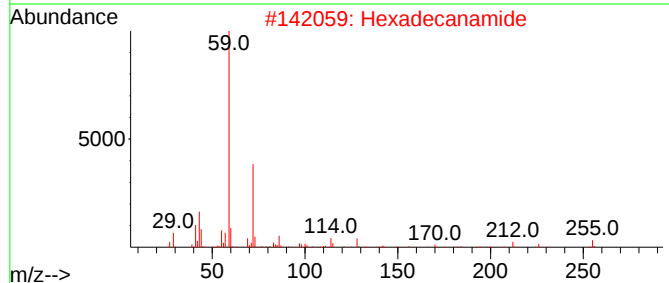
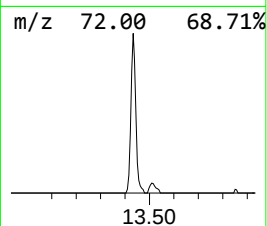
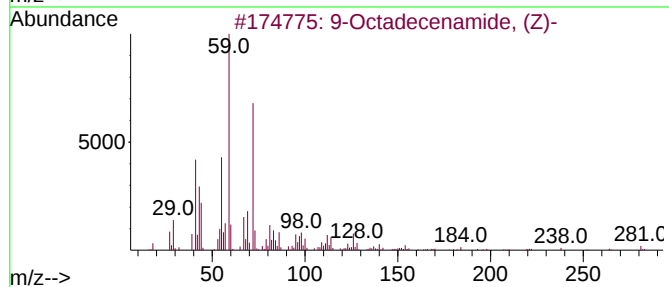
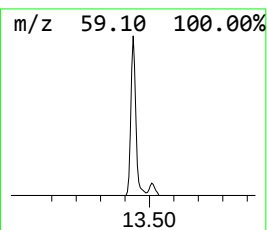
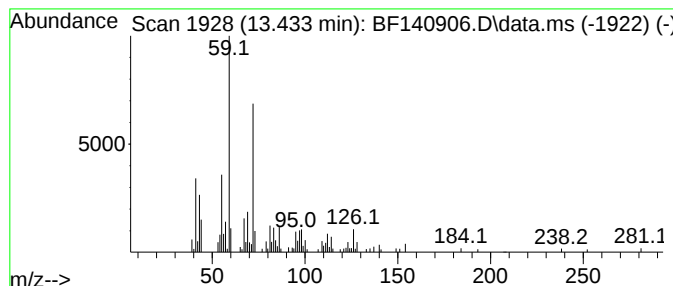
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Peak Number 3 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.433	3.85 ng	92256	Chrysene-d12	14.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	98
2		Hexadecanamide	255	C16H33NO	000629-54-9	64
3		Tetradecanamide	227	C14H29NO	000638-58-4	64
4		Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	59
5		Octanamide	143	C8H17NO	000629-01-6	53



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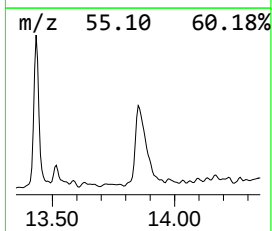
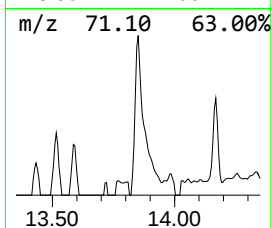
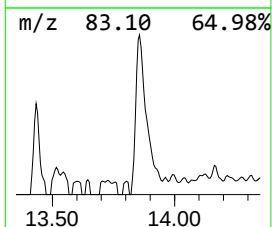
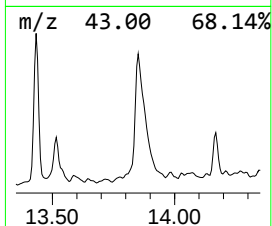
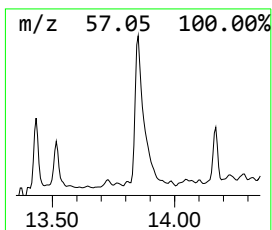
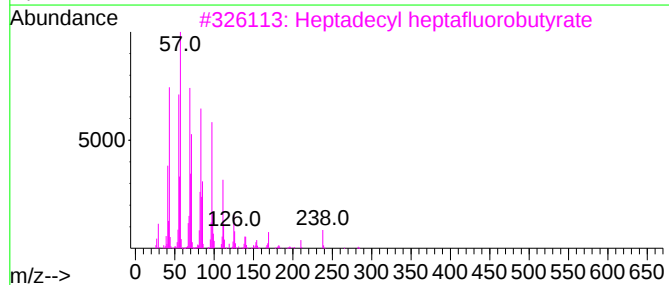
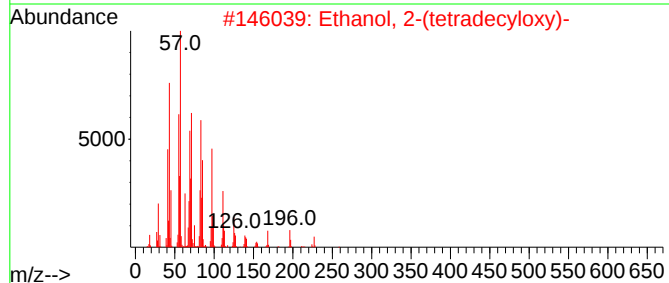
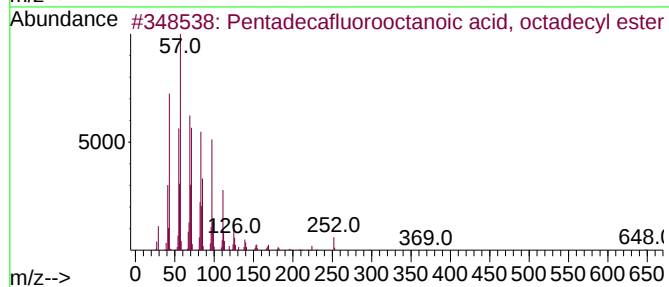
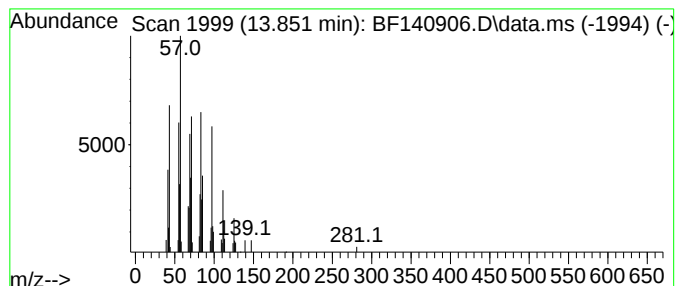
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Peak Number 4 Pentadecafluorooctanoic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.851	3.42 ng	82085	Chrysene-d12	14.027	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	95
2	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
4	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.063	3.0	ng	57486	1	6.845	389574	20.0
Benzophenone	10.598	2.4	ng	65399	3	9.880	553151	20.0
9-Octadecenamid...	13.433	3.9	ng	92256	5	14.027	479520	20.0
Pentadecafluoro...	13.851	3.4	ng	82085	5	14.027	479520	20.0