

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022625\
 Data File : BF141783.D
 Acq On : 26 Feb 2025 15:29
 Operator : RC/JU
 Sample : Q1422-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-154-SB01

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-BF020625.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141783.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.404	558	563	590	rBV	1065314	1576200	74.12%	10.037%
2	6.422	731	736	761	rBV	1079996	1638735	77.06%	10.435%
3	6.781	792	797	803	rBV	337053	425048	19.99%	2.707%
4	7.345	887	893	904	rBV	751256	1021262	48.02%	6.503%
5	7.604	933	937	953	rBV	21211	43645	2.05%	0.278%
6	8.063	1010	1015	1032	rBV	416039	573697	26.98%	3.653%
7	9.134	1192	1197	1208	rBV	1506567	1907486	89.69%	12.146%
8	9.810	1307	1312	1323	rBV	495610	671293	31.57%	4.275%
9	10.528	1429	1434	1442	rVB	131518	171742	8.08%	1.094%
10	10.604	1442	1447	1460	rBV	847082	1141750	53.69%	7.270%
11	11.298	1560	1565	1580	rBV	535385	698875	32.86%	4.450%
12	11.827	1648	1655	1677	rBV	441718	835820	39.30%	5.322%
13	12.539	1770	1776	1781	rBV6	20894	50167	2.36%	0.319%
14	12.610	1782	1788	1801	rVV	205653	413008	19.42%	2.630%
15	12.886	1829	1835	1840	rBV	1590082	2126650	100.00%	13.542%
16	13.063	1856	1865	1879	rVB	494523	1031348	48.50%	6.567%
17	13.421	1921	1926	1933	rBV3	17594	39553	1.86%	0.252%
18	13.786	1982	1988	2003	rBV2	42252	119371	5.61%	0.760%
19	13.939	2006	2014	2021	rVV	444902	621442	29.22%	3.957%
20	14.010	2022	2026	2034	rVB	18333	31159	1.47%	0.198%
21	14.504	2106	2110	2115	rVB	18259	26555	1.25%	0.169%
22	14.986	2187	2192	2196	rBV2	16360	30690	1.44%	0.195%
23	15.386	2255	2260	2271	rVB	300180	508970	23.93%	3.241%

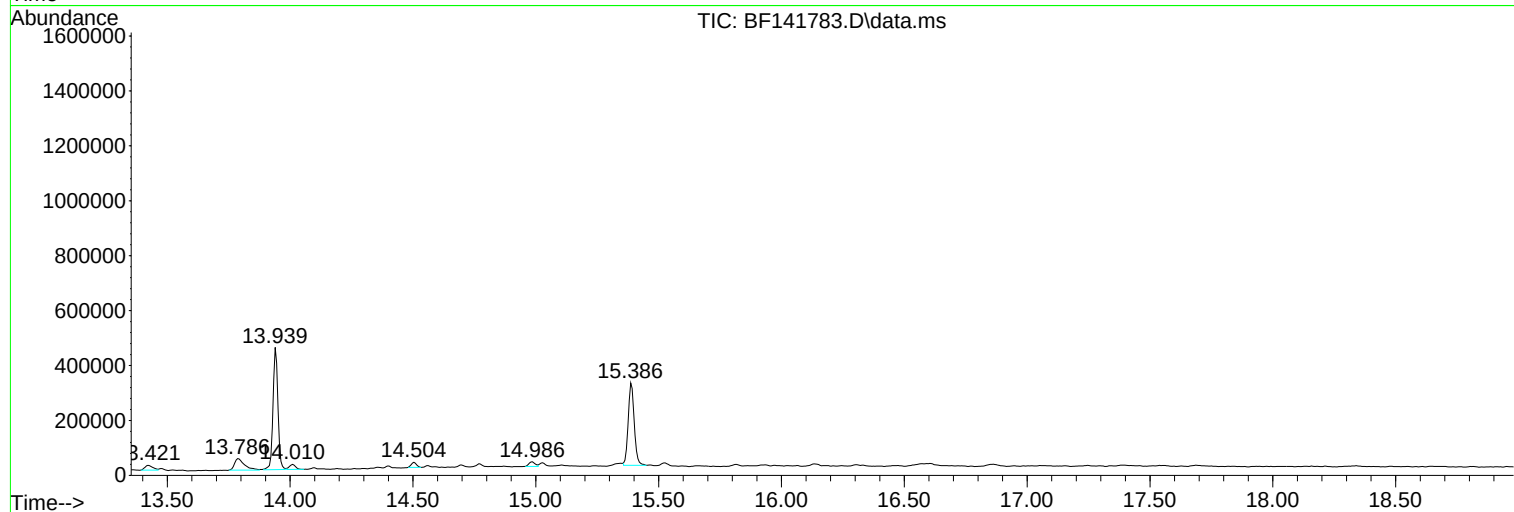
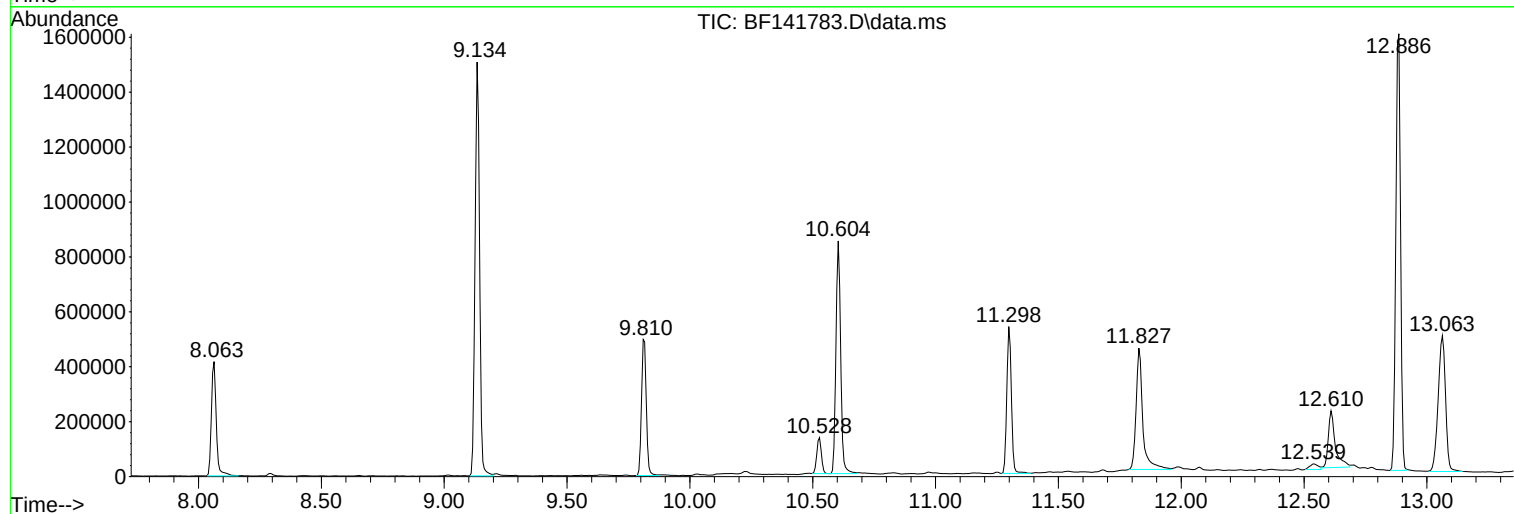
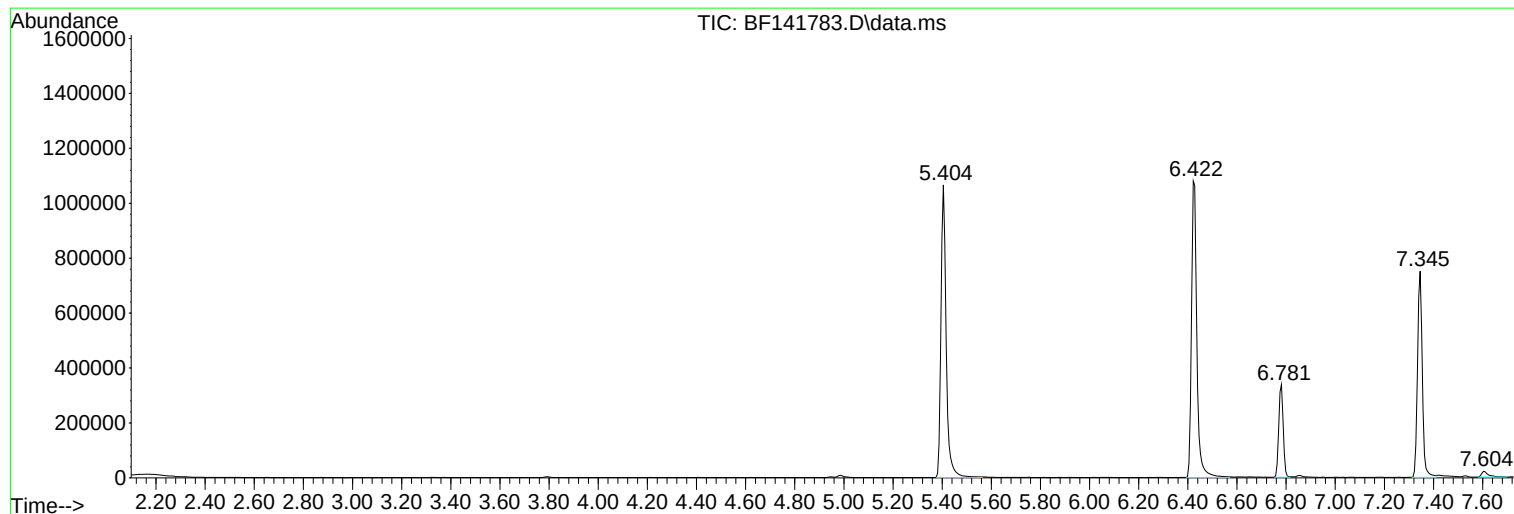
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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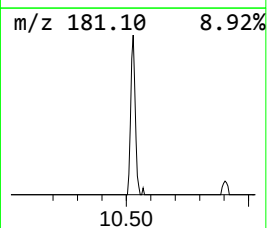
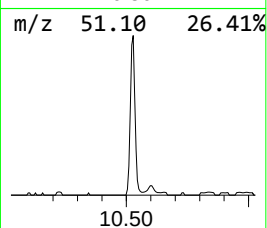
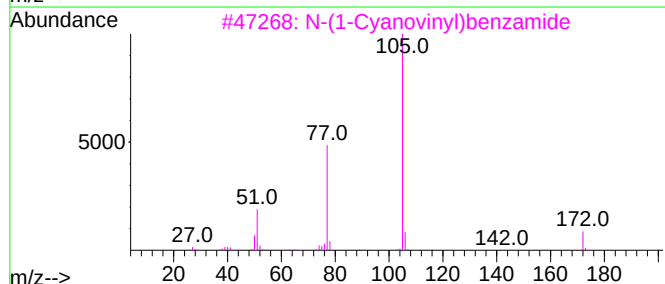
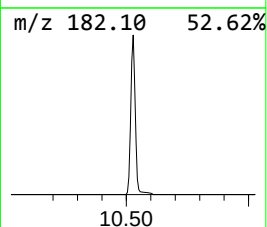
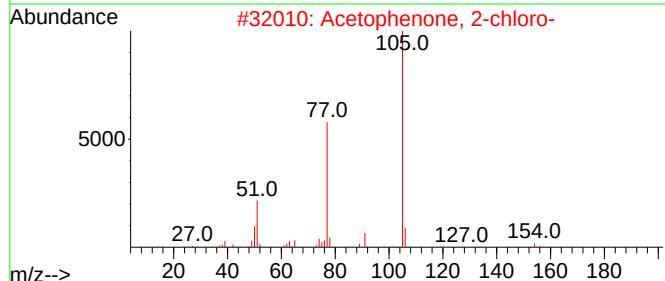
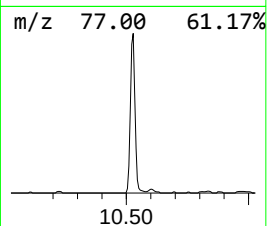
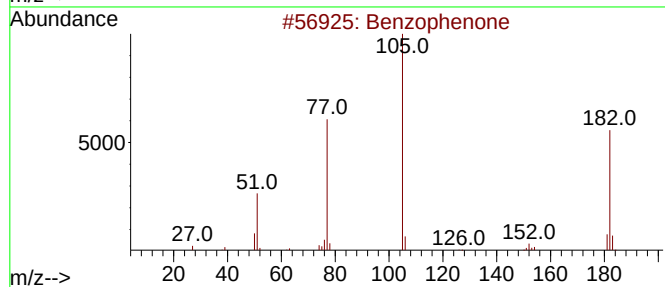
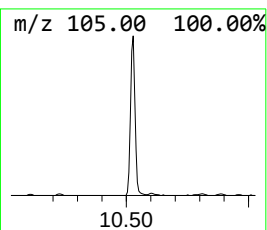
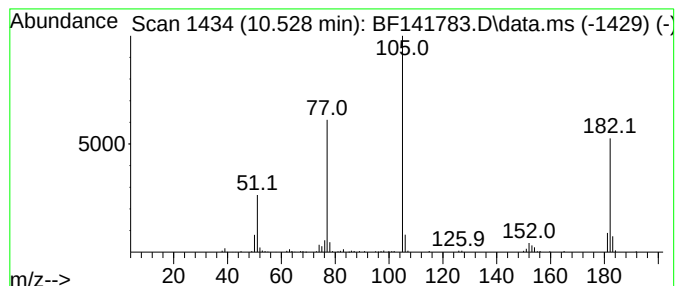
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.528	5.12 ng	171742	Acenaphthene-d10	9.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	53
3			N-(1-Cyanovinyl)benzamide	172	C10H8N2O	1000186-19-1	47
4			2-Benzoyl-4-bromo-5-methyl-1,2-o...	281	C11H8BrNO3	1000444-92-3	47
5			1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47



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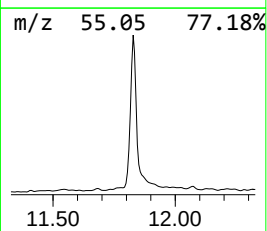
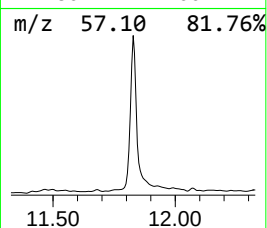
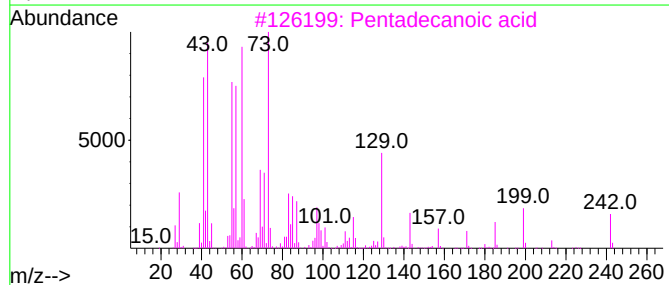
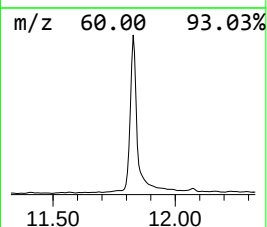
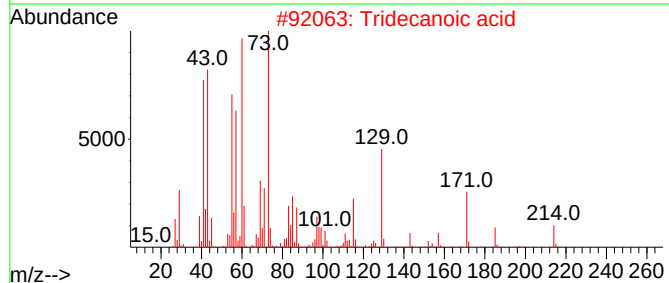
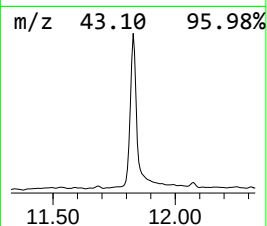
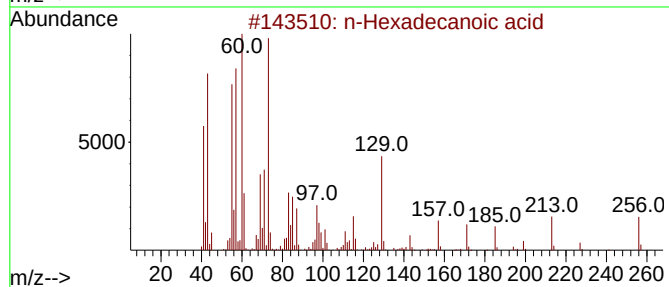
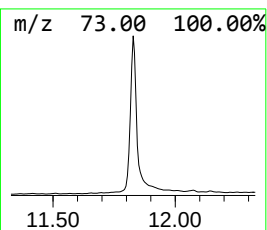
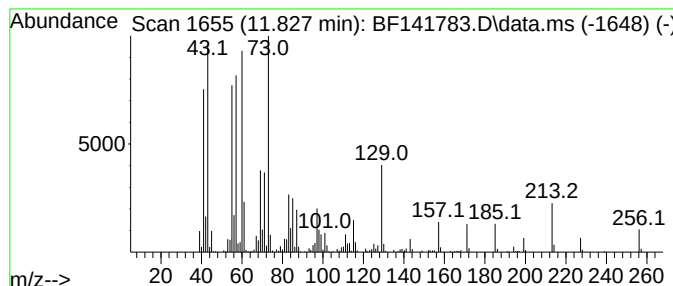
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.828	23.92 ng	835820	Phenanthrene-d10	11.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	91
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5			n-Decanoic acid	172	C10H20O2	000334-48-5	70



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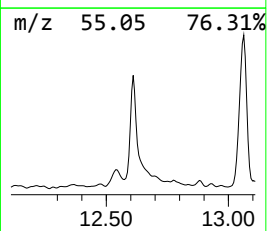
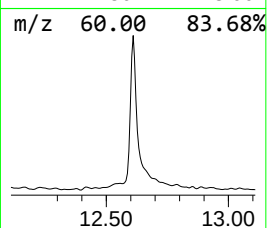
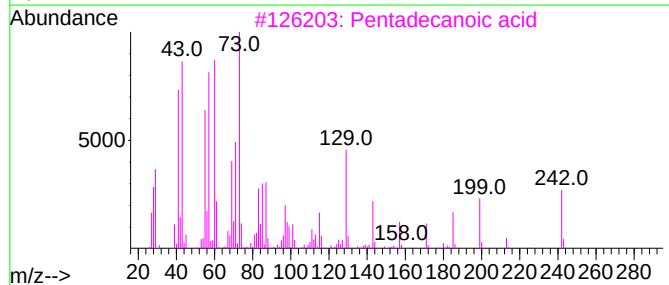
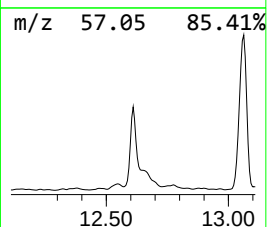
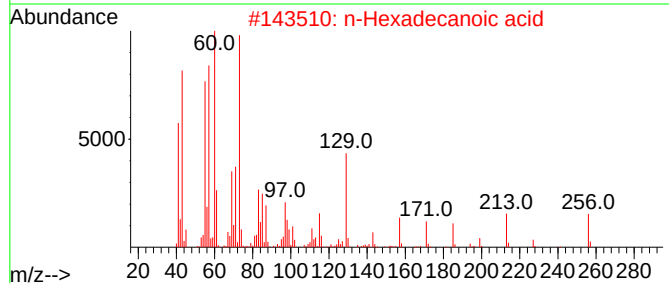
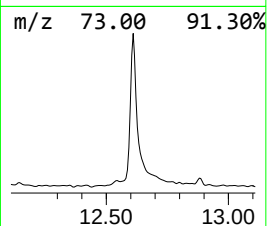
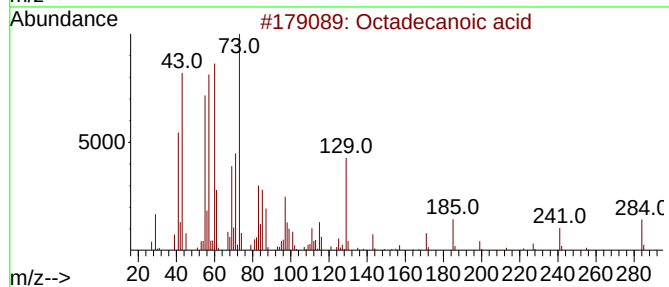
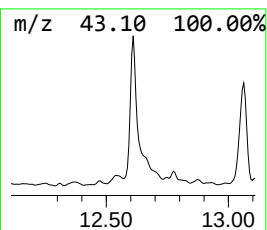
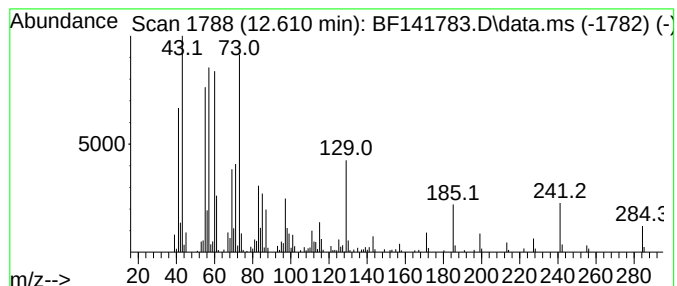
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Peak Number 3 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	11.82 ng	413008	Phenanthrene-d10	11.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			n-Decanoic acid	172	C10H20O2	000334-48-5	70



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ALS Vial : 10 Sample Multiplier: 1

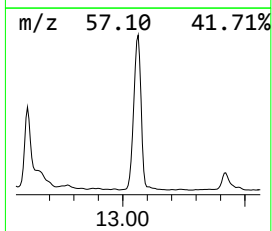
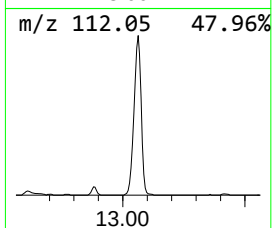
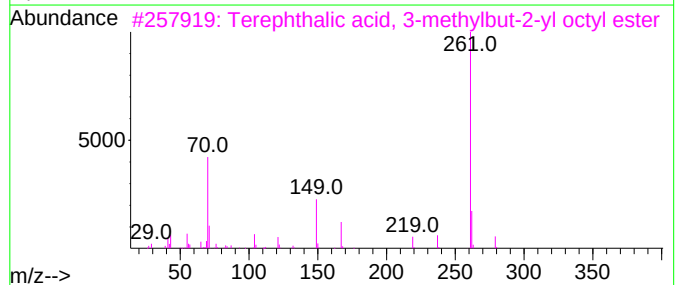
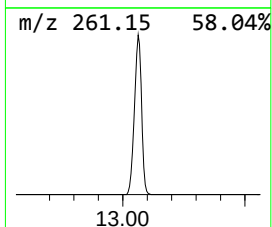
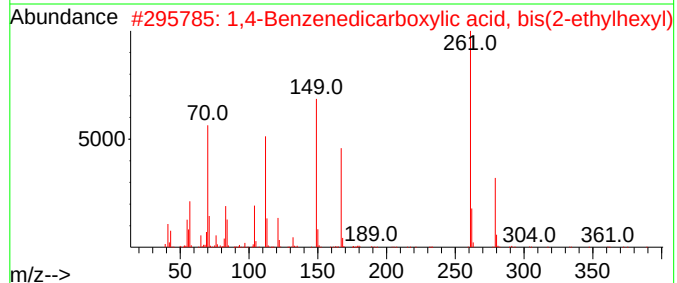
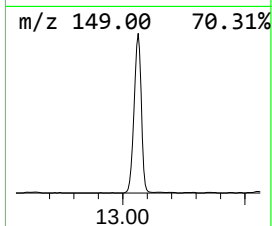
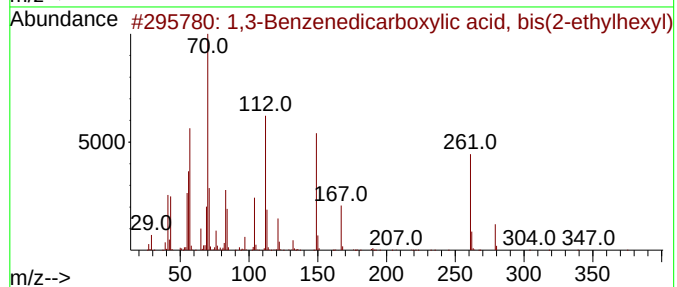
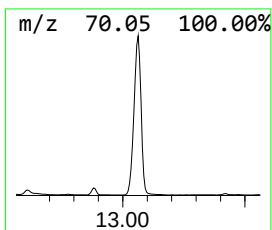
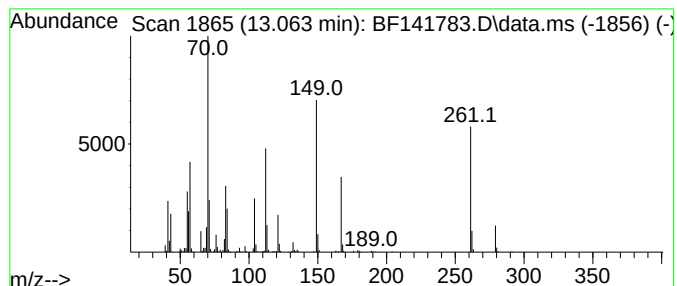
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Peak Number 4 1,3-Benzenedicarboxylic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.063	33.19 ng	1031350	Chrysene-d12	13.939	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	87
2	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	72
3	Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	53
4	Terephthalic acid, monochloride,...	296	C16H21ClO3	1000324-22-9	47
5	Terephthalic acid, 3-hexyl octyl...	362	C22H34O4	1000356-23-1	35



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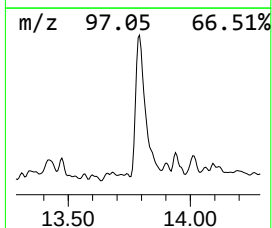
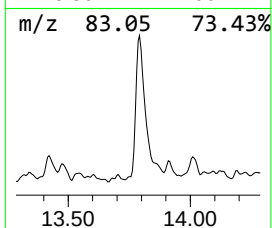
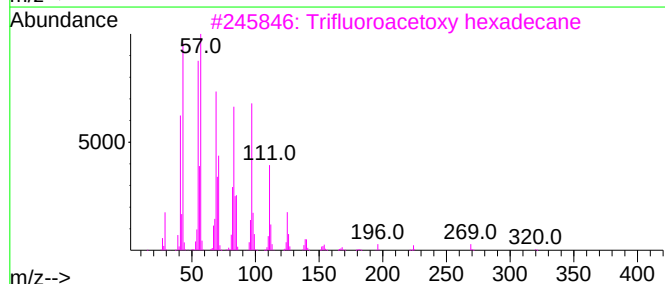
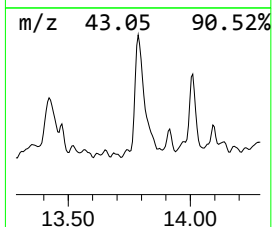
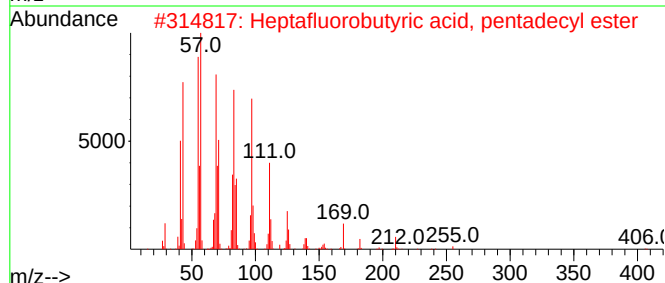
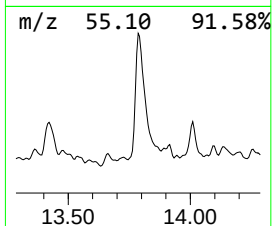
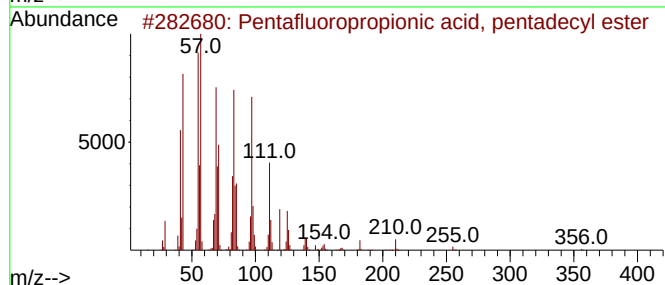
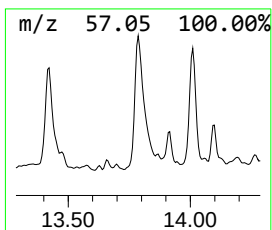
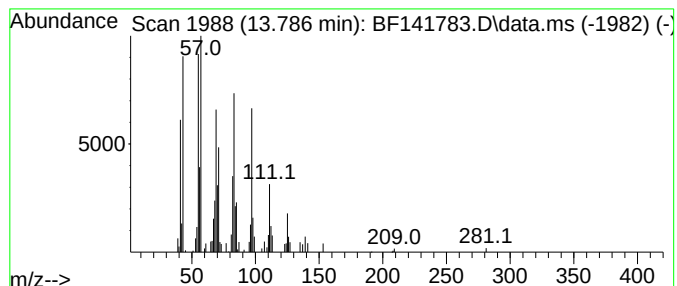
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Peak Number 5 Pentafluoropropionic acid, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.786	3.84 ng	119371	Chrysene-d12	13.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	95
2			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	95
3			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
5			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	10.528	5.1	ng	171742	3	9.810	671293	20.0
n-Hexadecanoic ...	11.828	23.9	ng	835820	4	11.298	698875	20.0
Octadecanoic acid	12.610	11.8	ng	413008	4	11.298	698875	20.0
1,3-Benzenedica...	13.063	33.2	ng	1031350	5	13.939	621442	20.0
Pentafluoroprop...	13.786	3.8	ng	119371	5	13.939	621442	20.0