

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051625\
 Data File : BF142435.D
 Acq On : 16 May 2025 19:11
 Operator : RC/JU
 Sample : Q2034-13
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 L3-WC-4

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-BF050525.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF142435.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.116	488	497	506	rBV	171070	259156	11.62%	1.508%
2	5.516	558	565	570	rBV	1575099	2130725	95.53%	12.397%
3	6.522	729	736	741	rBV	1646161	2230461	100.00%	12.977%
4	6.904	795	801	806	rBV	648682	824014	36.94%	4.794%
5	7.457	890	895	900	rBV	1010492	1303184	58.43%	7.582%
6	7.710	934	938	944	rBV	27528	36063	1.62%	0.210%
7	8.181	1013	1018	1033	rBV	811221	1040544	46.65%	6.054%
8	9.257	1195	1201	1206	rBV	1795422	2209009	99.04%	12.853%
9	9.933	1311	1316	1326	rBV	784711	1011117	45.33%	5.883%
10	10.645	1432	1437	1445	rBV	329967	413545	18.54%	2.406%
11	10.722	1445	1450	1461	rVV	887471	1108172	49.68%	6.448%
12	11.422	1564	1569	1577	rBV	630251	837039	37.53%	4.870%
13	11.945	1651	1658	1670	rBV3	105362	198199	8.89%	1.153%
14	12.033	1670	1673	1683	rVB2	15323	35240	1.58%	0.205%
15	12.633	1770	1775	1780	rBV	62858	83772	3.76%	0.487%
16	12.863	1804	1814	1819	rBV	65770	103622	4.65%	0.603%
17	13.010	1833	1839	1845	rVB	1213710	1582586	70.95%	9.208%
18	13.910	1987	1992	2002	rBV2	54735	106305	4.77%	0.619%
19	14.063	2011	2018	2030	rVB	605893	869380	38.98%	5.058%
20	14.533	2094	2098	2107	rBV5	18749	48160	2.16%	0.280%
21	14.827	2144	2148	2154	rBV4	25460	53097	2.38%	0.309%
22	15.115	2192	2197	2205	rBV	28831	63840	2.86%	0.371%
23	15.421	2246	2249	2254	rVB	16304	23714	1.06%	0.138%
24	15.486	2256	2260	2265	rVV	17865	26669	1.20%	0.155%
25	15.551	2265	2271	2284	rVB	360760	589613	26.43%	3.431%

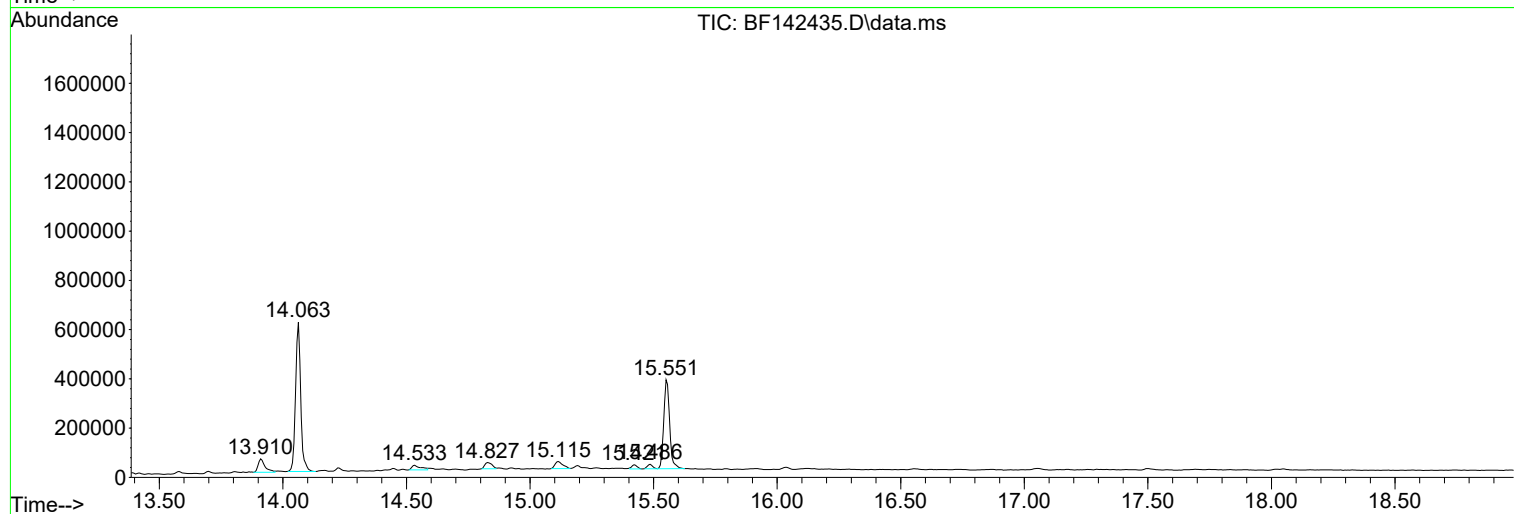
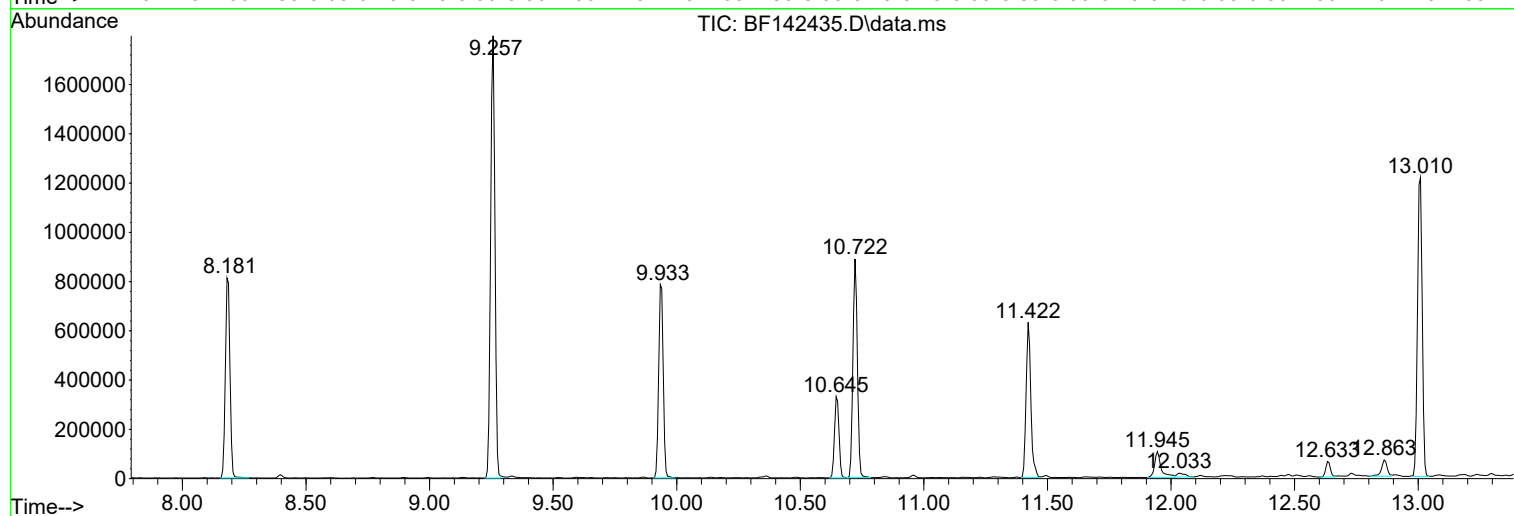
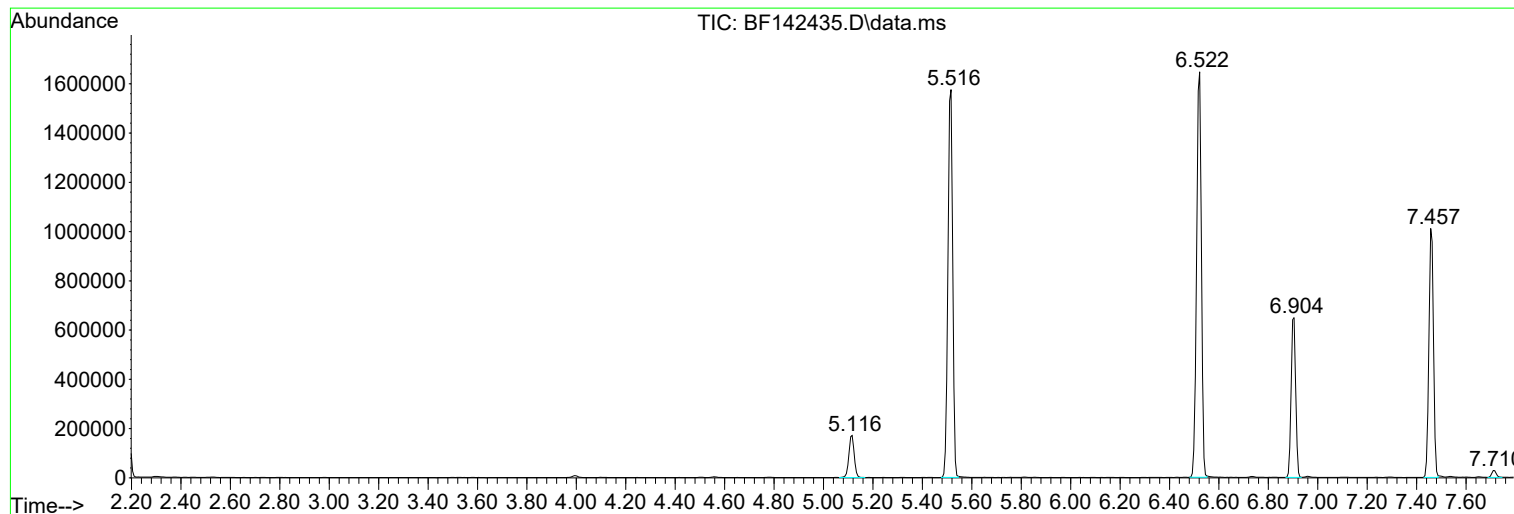
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Instrument :
BNA_F
ClientSampleId :
L3-WC-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF050525.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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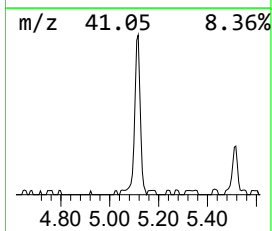
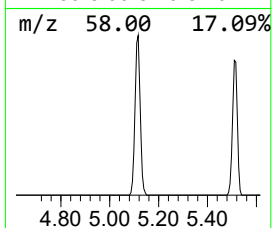
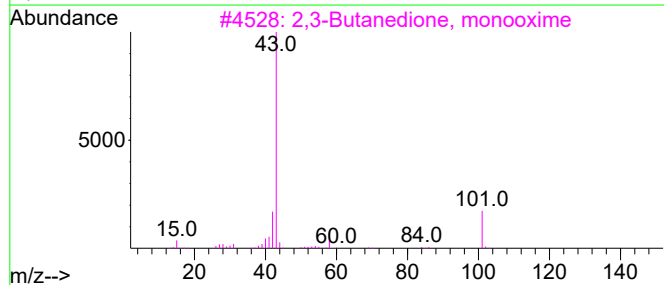
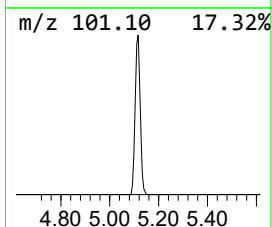
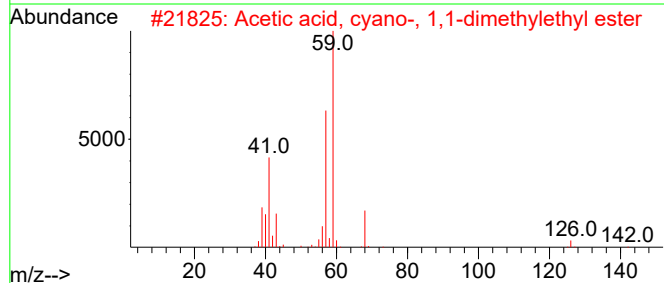
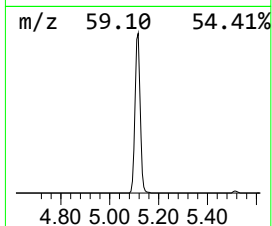
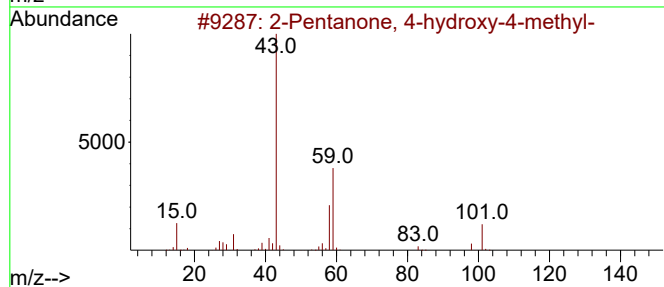
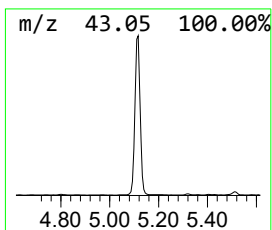
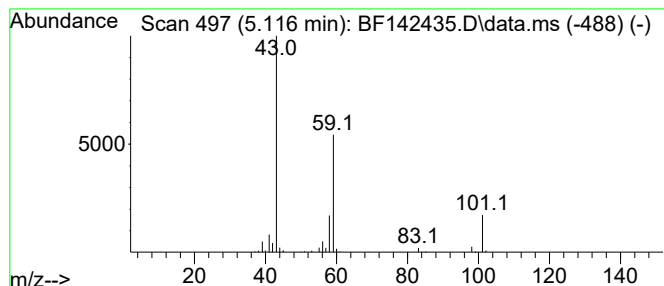
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF050525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.116	6.29 ng	259156	1,4-Dichlorobenzene-d4	6.904

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	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
5	1,2-Butanediol	90	C4H10O2	000584-03-2	9		



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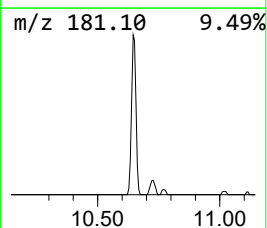
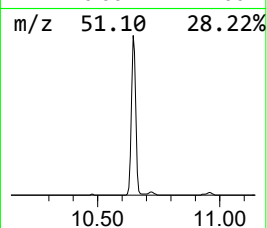
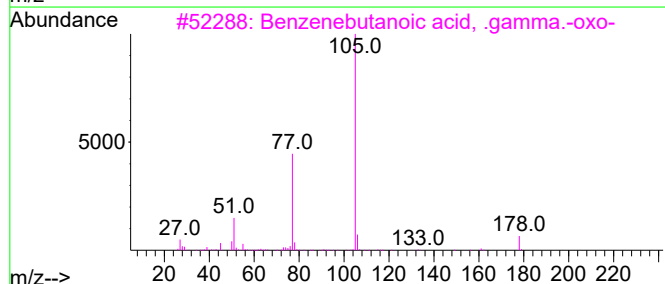
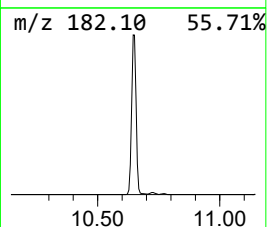
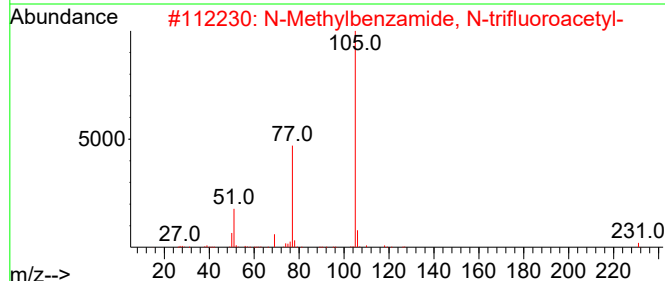
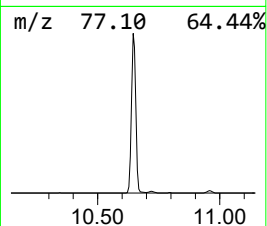
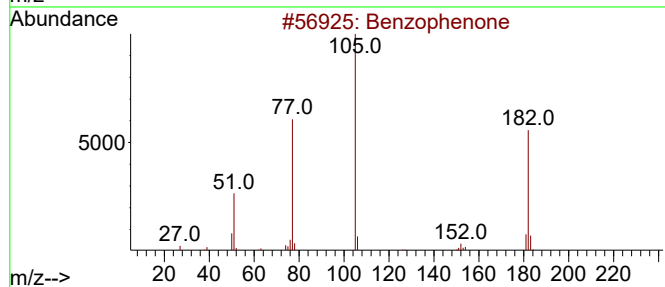
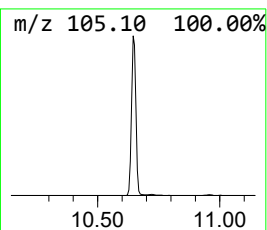
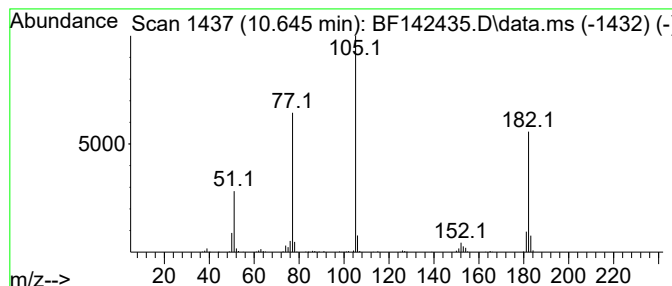
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.645	8.18 ng	413545	Acenaphthene-d10	9.933

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			N-Methylbenzamide, N-trifluoroac...	231	C10H8F3NO2	1000446-91-5	47
3			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
4			2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47
5			1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47



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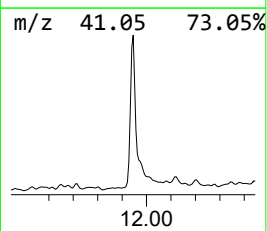
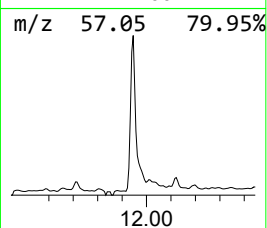
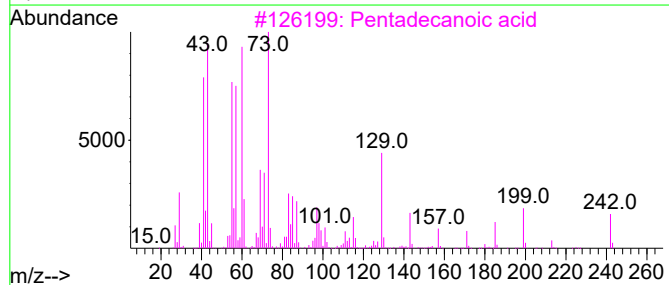
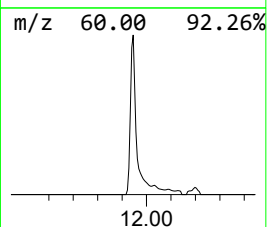
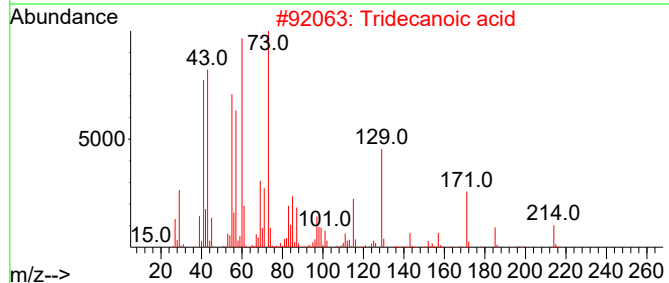
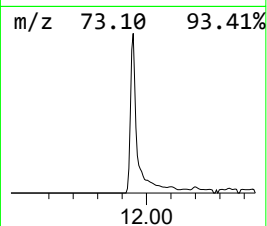
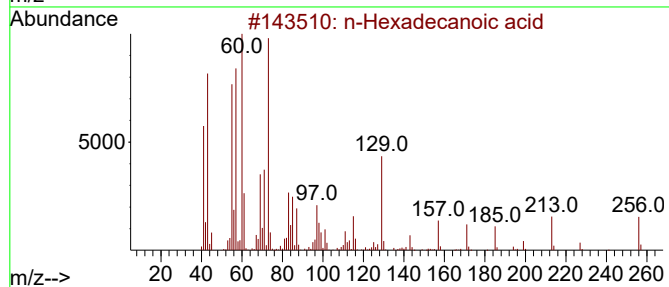
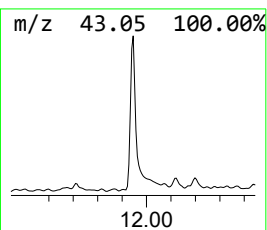
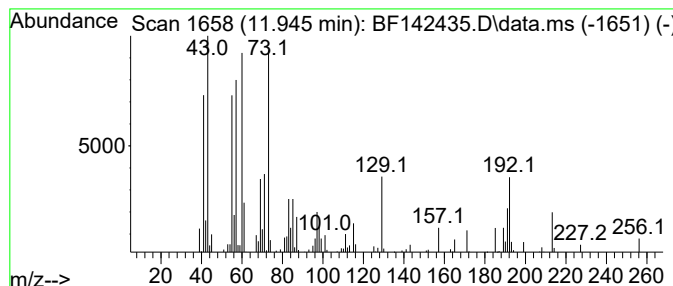
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	4.74 ng	198199	Phenanthrene-d10	11.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2			Tridecanoic acid	214	C13H26O2	000638-53-9	86
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	62
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	60



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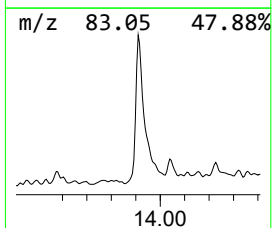
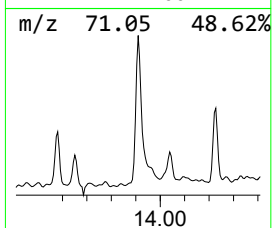
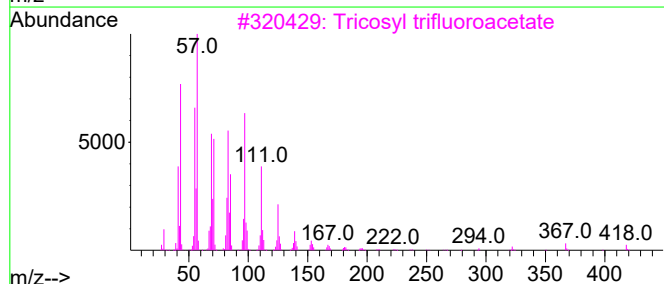
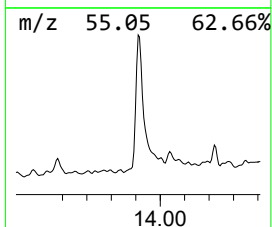
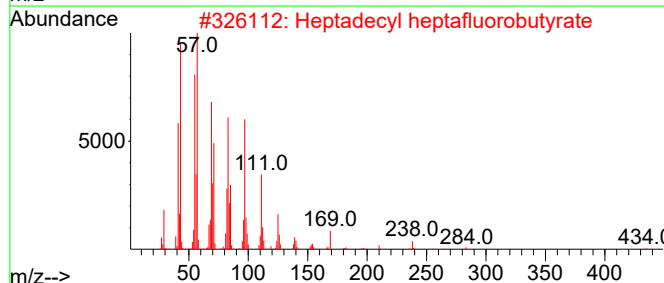
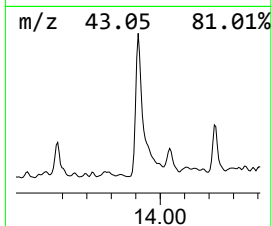
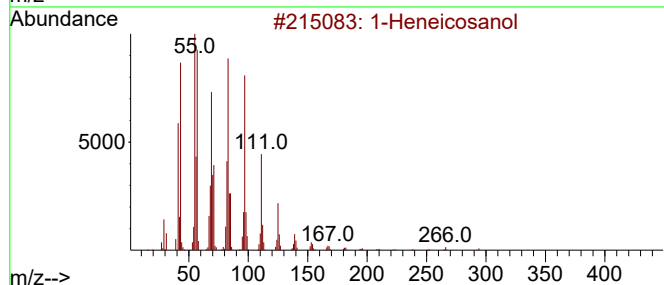
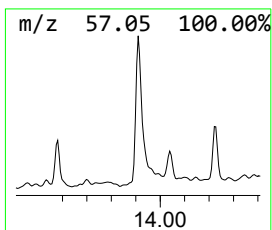
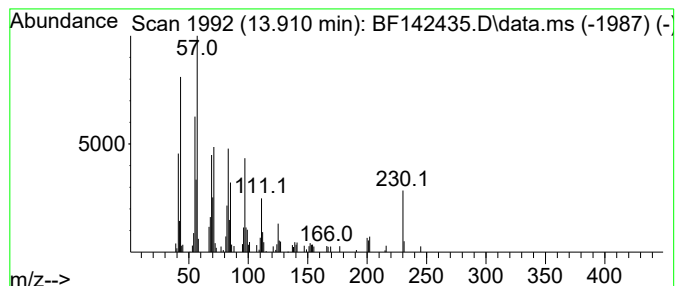
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L3-WC-4

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

Peak Number 5 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.910	2.45 ng	106305	Chrysene-d12	14.063	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	92
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
3	Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	90
4	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	90
5	Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	90



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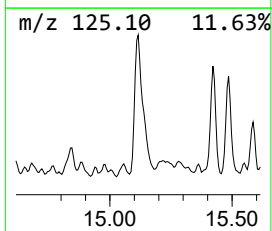
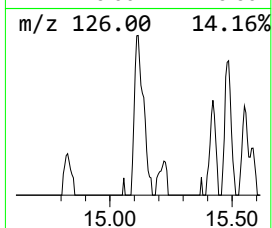
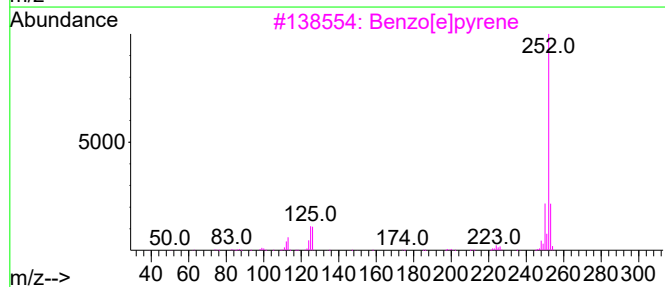
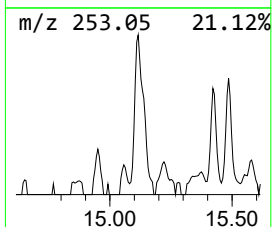
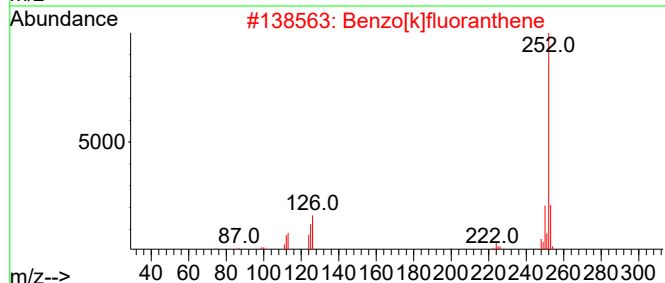
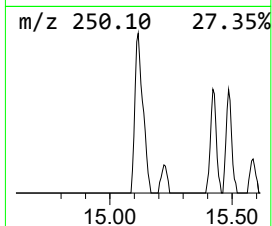
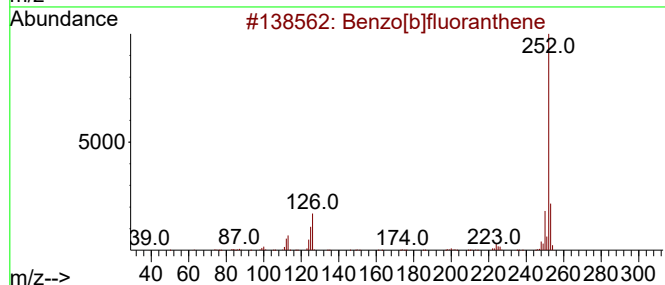
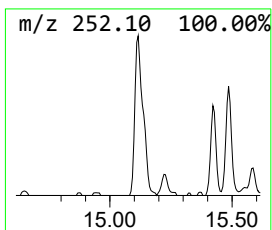
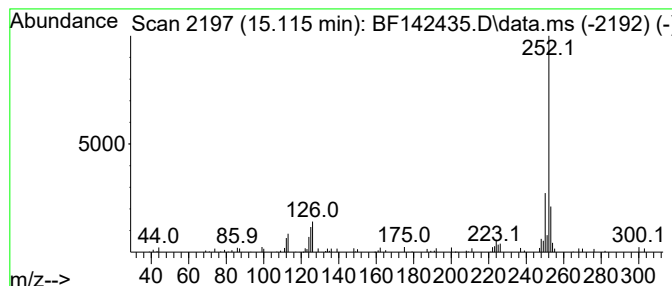
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.115	2.17 ng	63840	Perylene-d12	15.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3			Benzo[e]pyrene	252	C20H12	000192-97-2	94
4			Perylene	252	C20H12	000198-55-0	93
5			Benzo[a]pyrene	252	C20H12	000050-32-8	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.116	6.3	ng	259156	1	6.904	824014	20.0
Benzophenone	10.645	8.2	ng	413545	3	9.933	1011120	20.0
n-Hexadecanoic ...	11.945	4.7	ng	198199	4	11.422	837039	20.0
1-Heneicosanol	13.910	2.5	ng	106305	5	14.063	869380	20.0
Benzo[e]pyrene	15.115	2.2	ng	63840	6	15.551	589613	20.0