

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092722\
 Data File : BP011816.D
 Acq On : 27 Sep 2022 15:25
 Operator : CG/JU
 Sample : N4837-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 251563-3

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_P\Methods\8270E-BP092122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP011816.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.046	6	10	32	rVB	7553156	11934089	48.86%	10.165%
2	5.469	415	422	440	rBV	2659531	4175549	17.10%	3.556%
3	7.069	686	694	711	rBV	1499329	2470781	10.12%	2.104%
4	7.910	829	837	845	rBV	1317397	2098007	8.59%	1.787%
5	9.075	1028	1035	1056	rBV	3926988	6776785	27.75%	5.772%
6	10.499	1271	1277	1301	rBV	294809	666036	2.73%	0.567%
7	10.722	1307	1315	1328	rBV	1829363	3089379	12.65%	2.631%
8	13.181	1725	1733	1750	rBV	10501197	15971528	65.39%	13.603%
9	14.551	1959	1966	1984	rBV	2863102	4318147	17.68%	3.678%
10	16.045	2214	2220	2241	rBV	8673627	12682636	51.93%	10.802%
11	17.310	2428	2435	2446	rBV	3775577	5332871	21.83%	4.542%
12	18.192	2580	2585	2595	rBV	521630	1061848	4.35%	0.904%
13	19.427	2790	2795	2817	rBV	1359980	2991558	12.25%	2.548%
14	19.863	2864	2869	2885	rBV	18547467	24424308	100.00%	20.803%
15	21.098	3076	3079	3088	rBV2	271964	495117	2.03%	0.422%
16	21.392	3124	3129	3136	rBV	4606991	5836786	23.90%	4.971%
17	21.516	3145	3150	3166	rBV	4996945	7159413	29.31%	6.098%
18	23.833	3538	3544	3555	rVB2	2831833	5923126	24.25%	5.045%

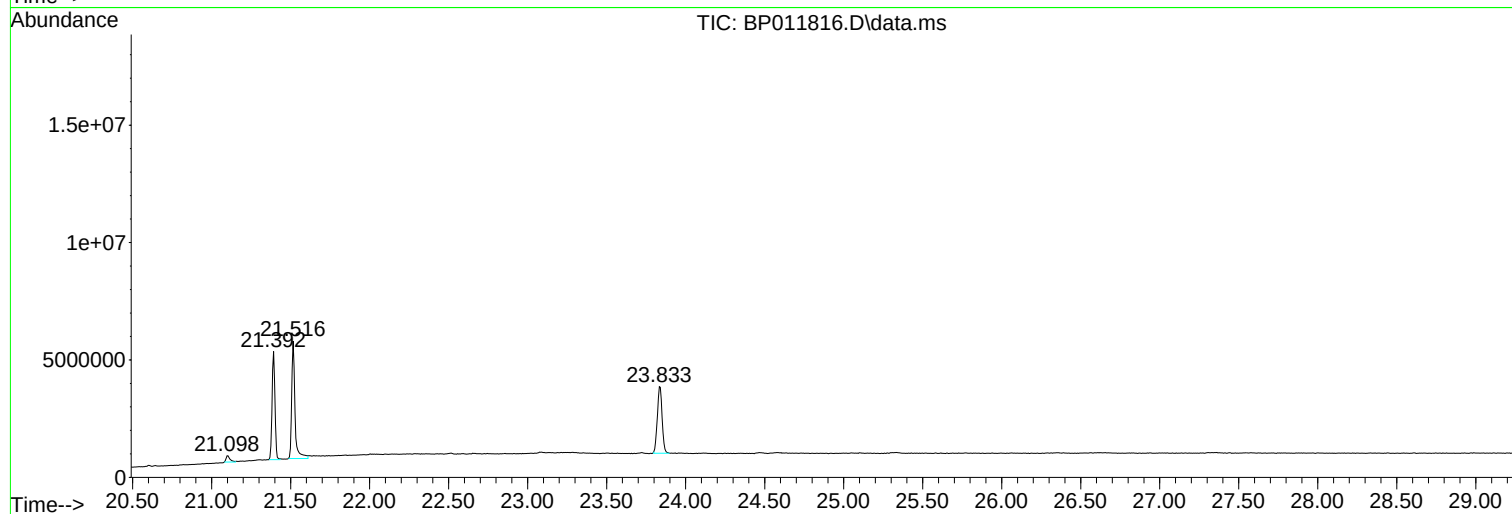
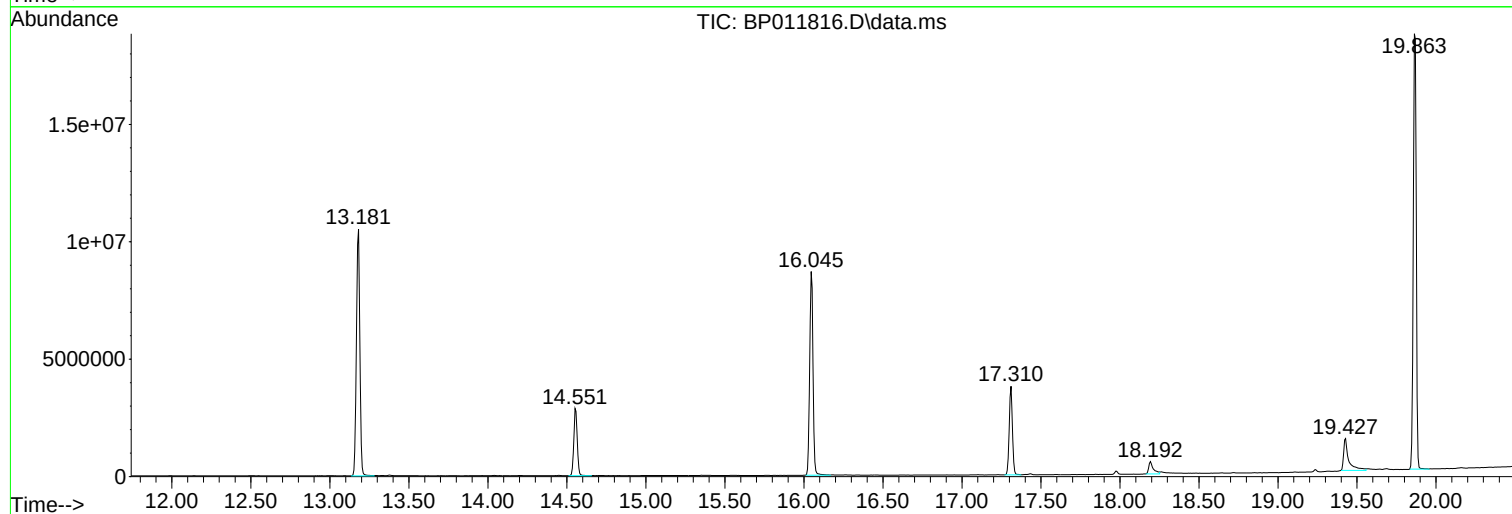
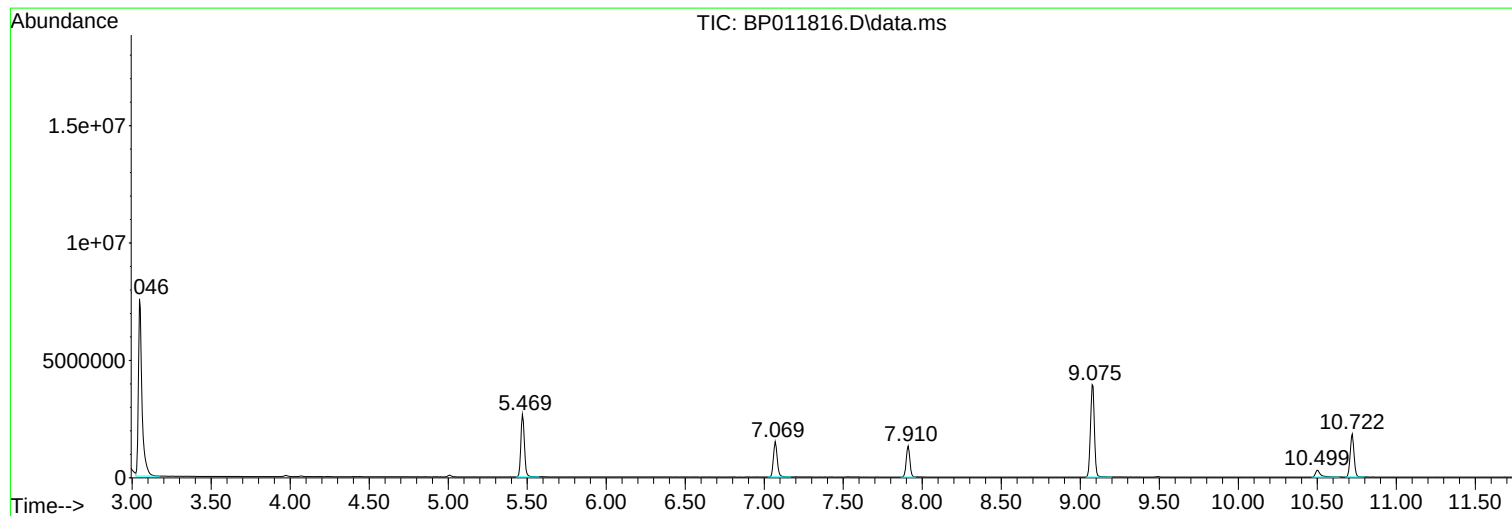
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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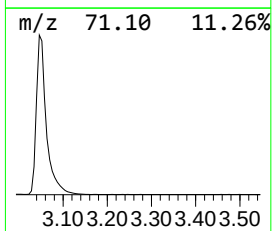
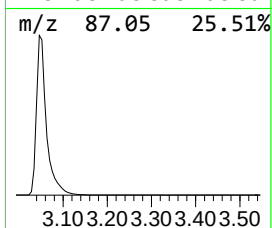
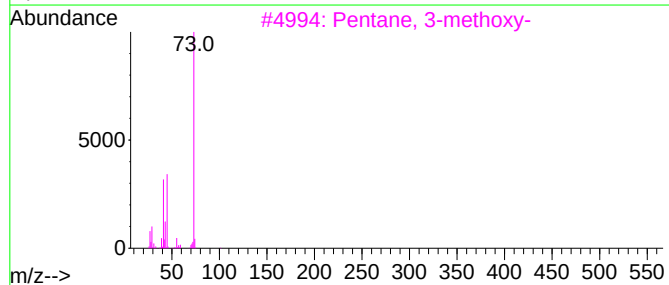
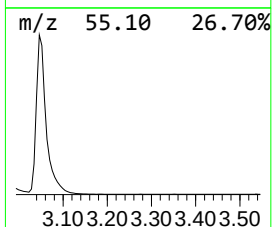
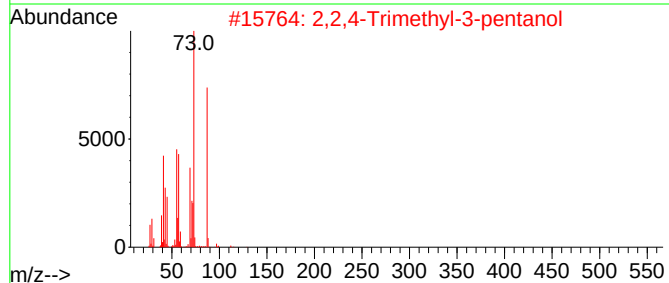
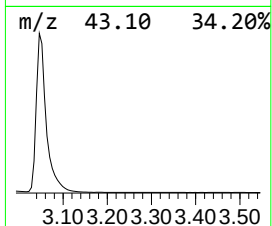
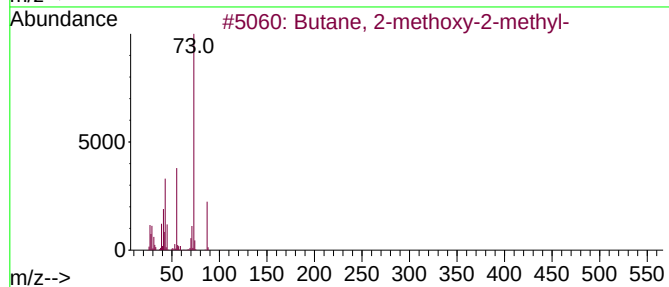
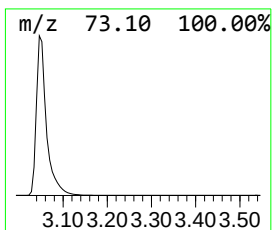
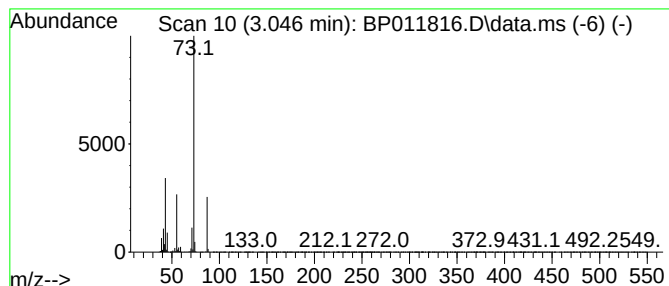
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.046	113.77 ng	11934100	1,4-Dichlorobenzene-d4	7.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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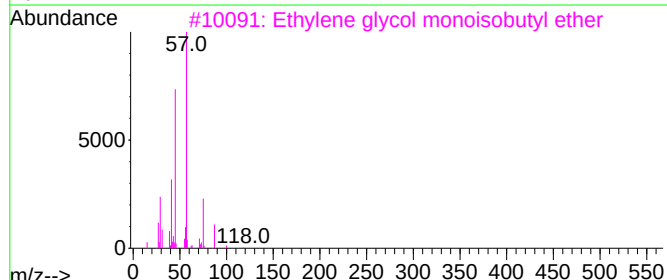
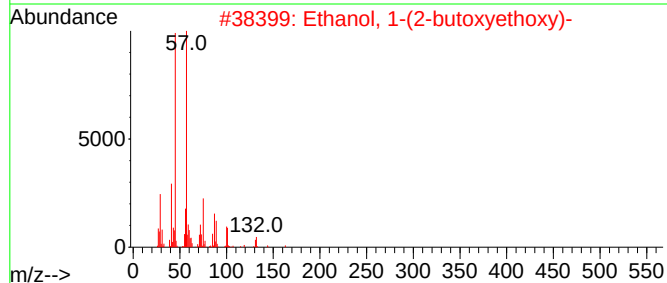
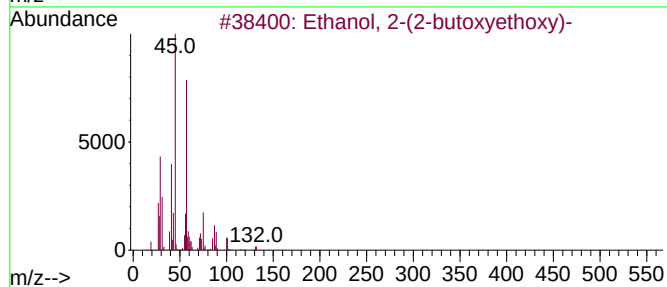
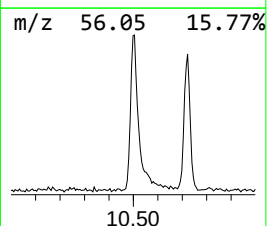
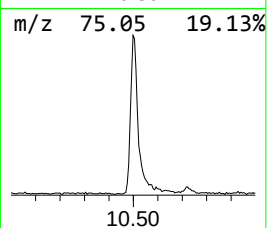
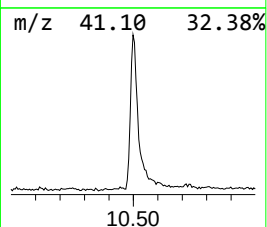
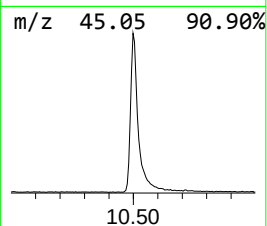
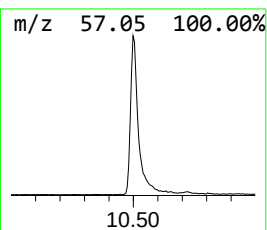
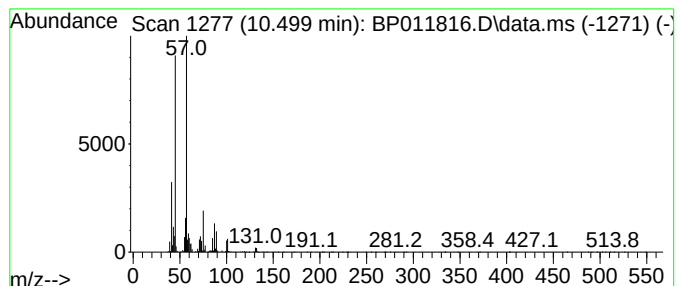
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.499	4.31 ng	666036	Naphthalene-d8	10.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50
5			Di-sec-Butyl ether	130	C8H18O	006863-58-7	50



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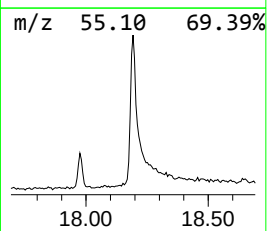
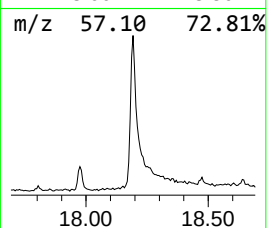
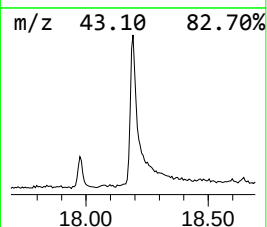
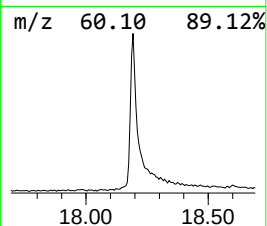
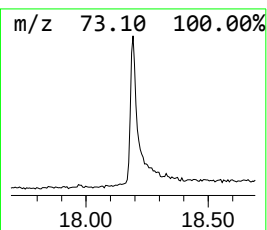
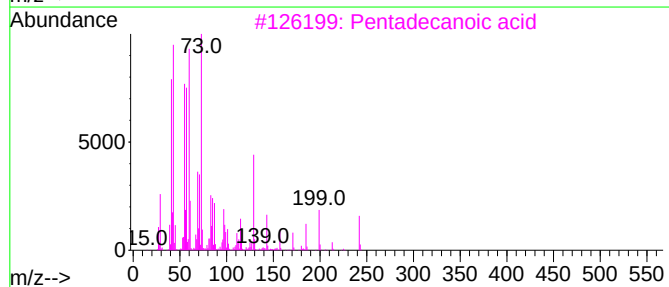
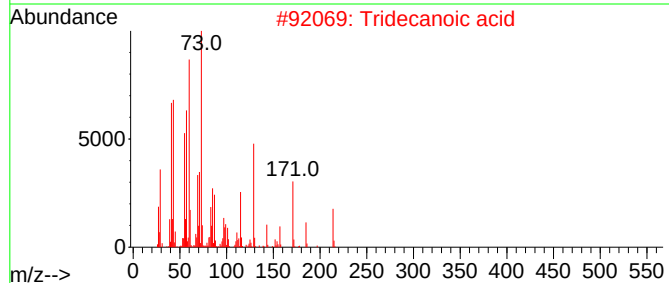
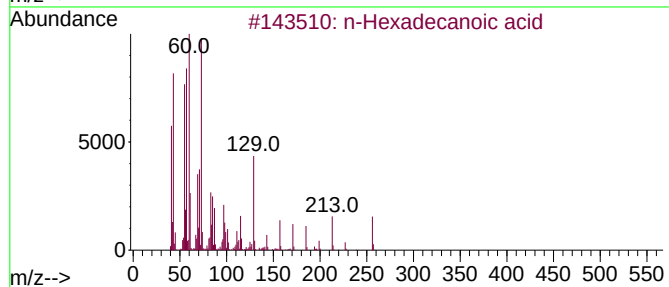
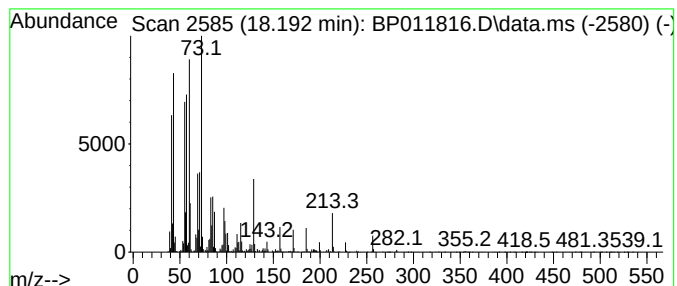
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.192	3.98 ng	1061850	Phenanthrene-d10		17.310	
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	94
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	90
4	Octadecanoic acid		284	C18H36O2	000057-11-4	87
5	Tetradecanoic acid		228	C14H28O2	000544-63-8	72



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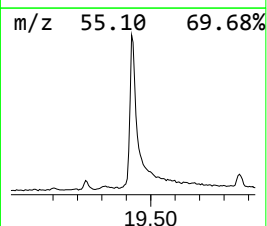
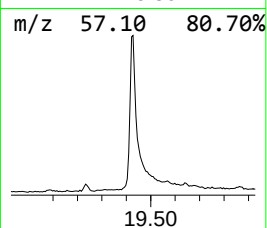
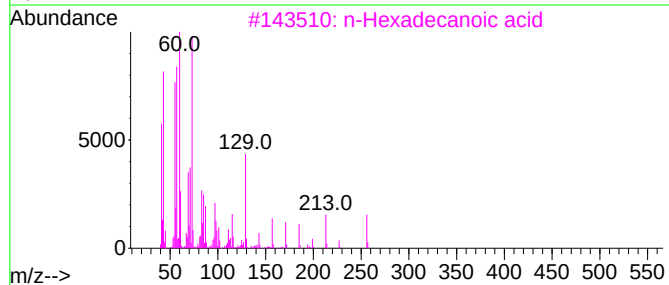
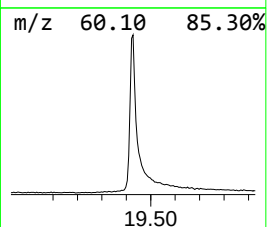
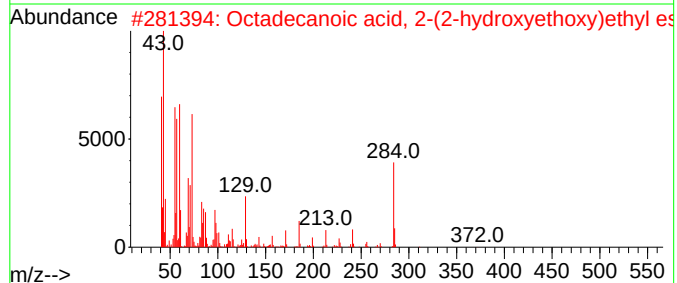
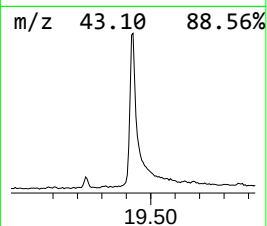
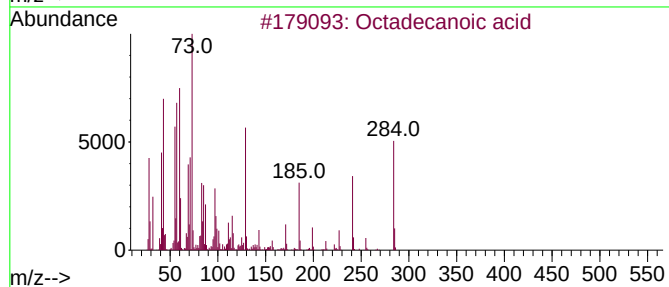
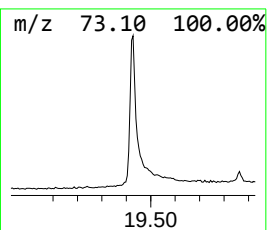
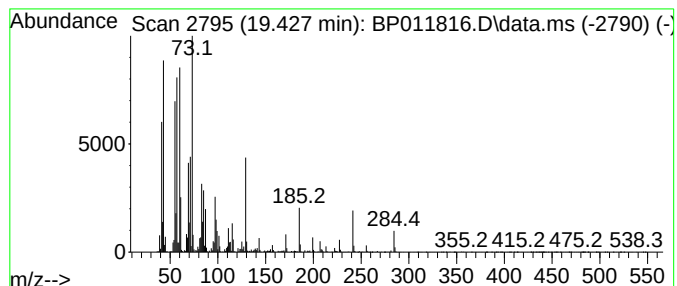
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Peak Number 4 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.427	10.25 ng	2991560	Chrysene-d12	21.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	87
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	76



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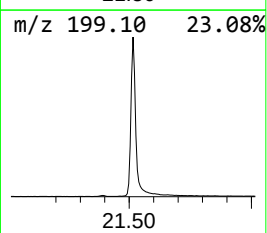
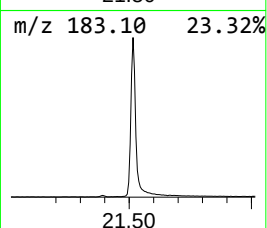
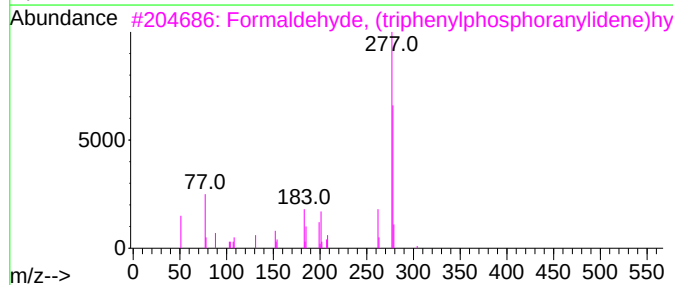
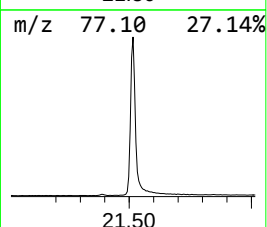
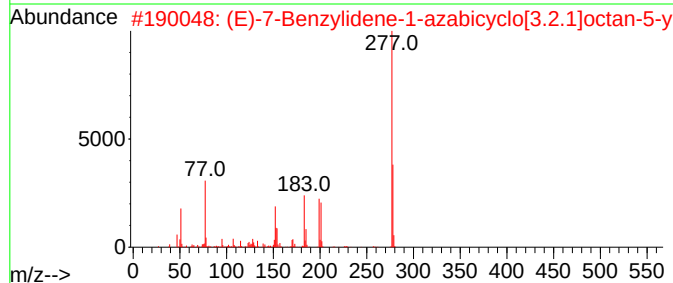
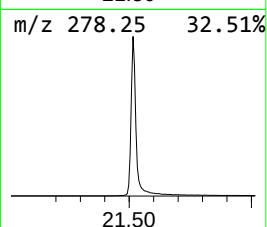
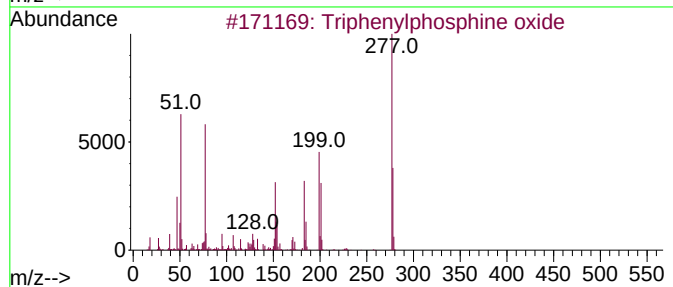
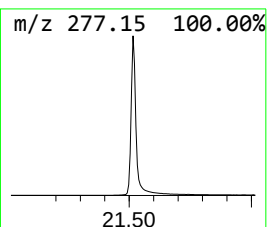
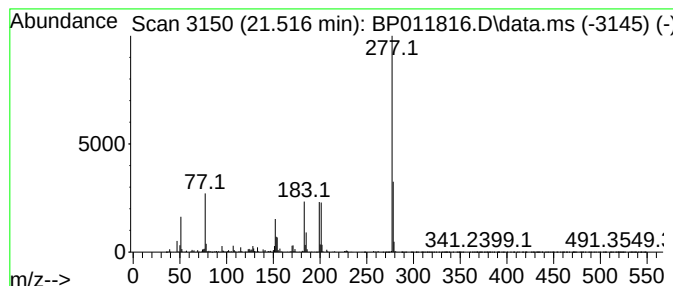
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Peak Number 5 Triphenylphosphine oxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.516	24.53 ng	7159410	Chrysene-d12	21.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	91
2			(E)-7-Benzylidene-1-azabicyclo[3.1.0]hex-5-ene	293	C15H19N03S	1000483-83-4	91
3			Formaldehyde, (triphenylphosphoranylidene)hy	304	C19H17N2P	015990-54-2	64
4			4-Nitro-4'-methyldiphenylsulfone	277	C13H11N04S	004094-37-5	43
5			Vanadium, cycloheptatrienyl-(pen...)	277	C17H22V	1000155-20-5	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.046	113.8	ng	11934100	1	7.910	2098010	20.0
Ethanol, 2-(2-b...	10.499	4.3	ng	666036	2	10.722	3089380	20.0
n-Hexadecanoic ...	18.192	4.0	ng	1061850	4	17.310	5332870	20.0
Octadecanoic acid	19.427	10.3	ng	2991560	5	21.392	5836790	20.0
Triphenylphosph...	21.516	24.5	ng	7159410	5	21.392	5836790	20.0