

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
 Data File : BP019806.D
 Acq On : 21 Feb 2024 14:05
 Operator : MA/JU
 Sample : P1523-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20240763-COMPOSITE-39N-N-3-01

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-BP013124.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP019806.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	24	rVB	2138599	2611411	12.63%	5.474%
2	4.069	160	167	175	rVB	150487	247101	1.20%	0.518%
3	4.640	259	264	269	rBV	246858	353647	1.71%	0.741%
4	5.016	322	328	337	rBV	284380	414783	2.01%	0.870%
5	5.469	399	405	415	rBV	891990	1340813	6.49%	2.811%
6	7.069	669	677	684	rBV	798386	1330925	6.44%	2.790%
7	7.893	811	817	824	rVB	251842	398771	1.93%	0.836%
8	9.063	1009	1016	1027	rBV	525016	879238	4.25%	1.843%
9	10.693	1283	1293	1302	rBV	294523	532199	2.57%	1.116%
10	13.157	1705	1712	1720	rBV	1126801	1716886	8.31%	3.599%
11	13.428	1753	1758	1764	rBV	174985	279287	1.35%	0.585%
12	14.528	1939	1945	1958	rVB	459157	685396	3.32%	1.437%
13	16.022	2187	2199	2221	rVB	1325467	2930491	14.18%	6.143%
14	16.416	2238	2266	2282	rBV2	4215786	20668488	100.00%	43.328%
15	16.686	2307	2312	2320	rVB2	155059	271457	1.31%	0.569%
16	17.286	2408	2414	2420	rVB	590080	851952	4.12%	1.786%
17	18.192	2556	2568	2581	rBV	1376985	2347484	11.36%	4.921%
18	18.457	2608	2613	2616	rBV	639395	872217	4.22%	1.828%
19	18.498	2616	2620	2630	rVB	667879	1022169	4.95%	2.143%
20	19.422	2770	2777	2789	rBV	1149927	1653262	8.00%	3.466%
21	19.851	2845	2850	2855	rBV	2345527	2857063	13.82%	5.989%
22	20.204	2905	2910	2918	rVB	207458	349363	1.69%	0.732%
23	21.098	3059	3062	3066	rBV	190629	212104	1.03%	0.445%
24	21.380	3105	3110	3120	rVV	857096	1210249	5.86%	2.537%
25	21.933	3200	3204	3213	rVB2	231046	381737	1.85%	0.800%
26	23.792	3514	3520	3533	rVB	600396	1283683	6.21%	2.691%

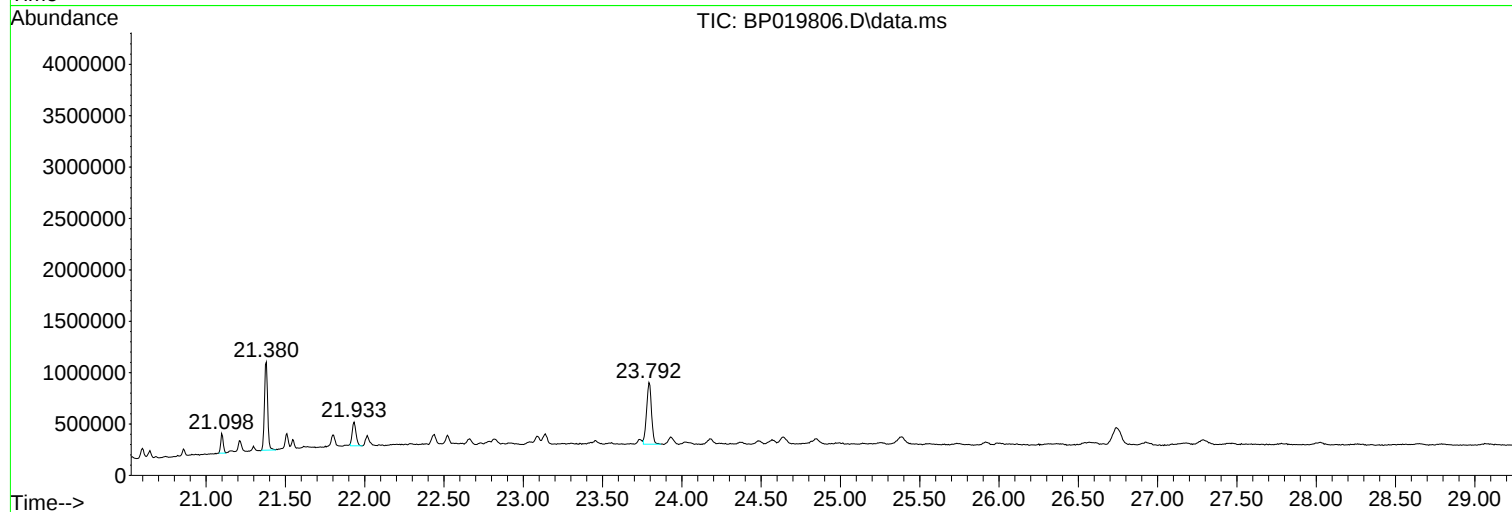
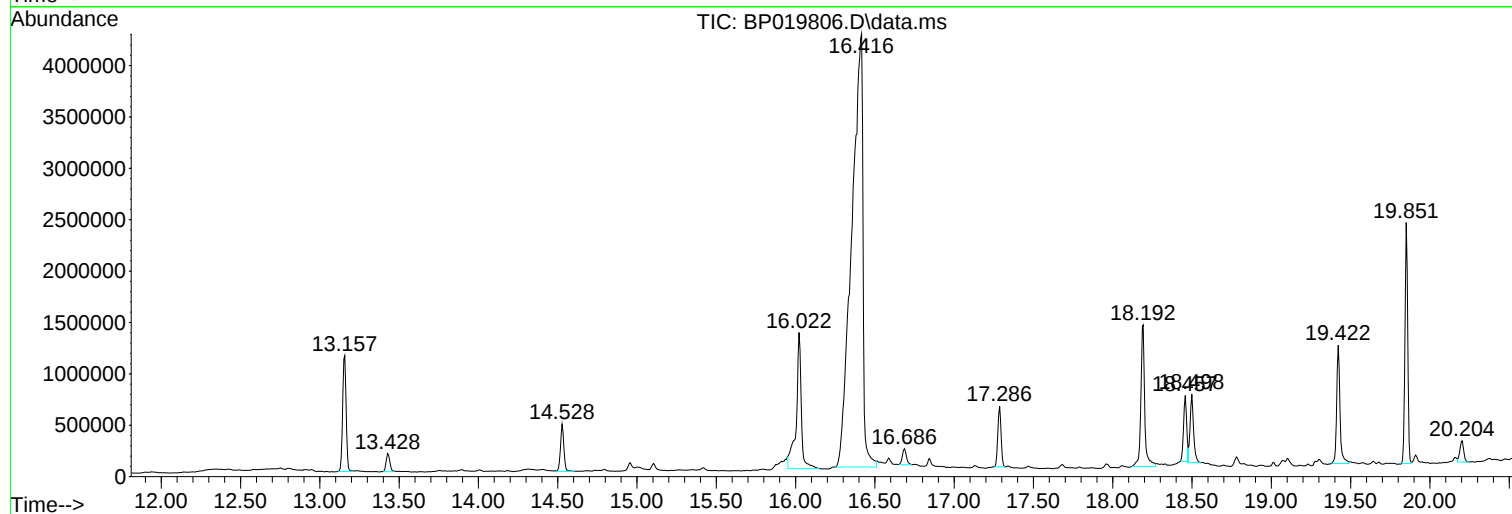
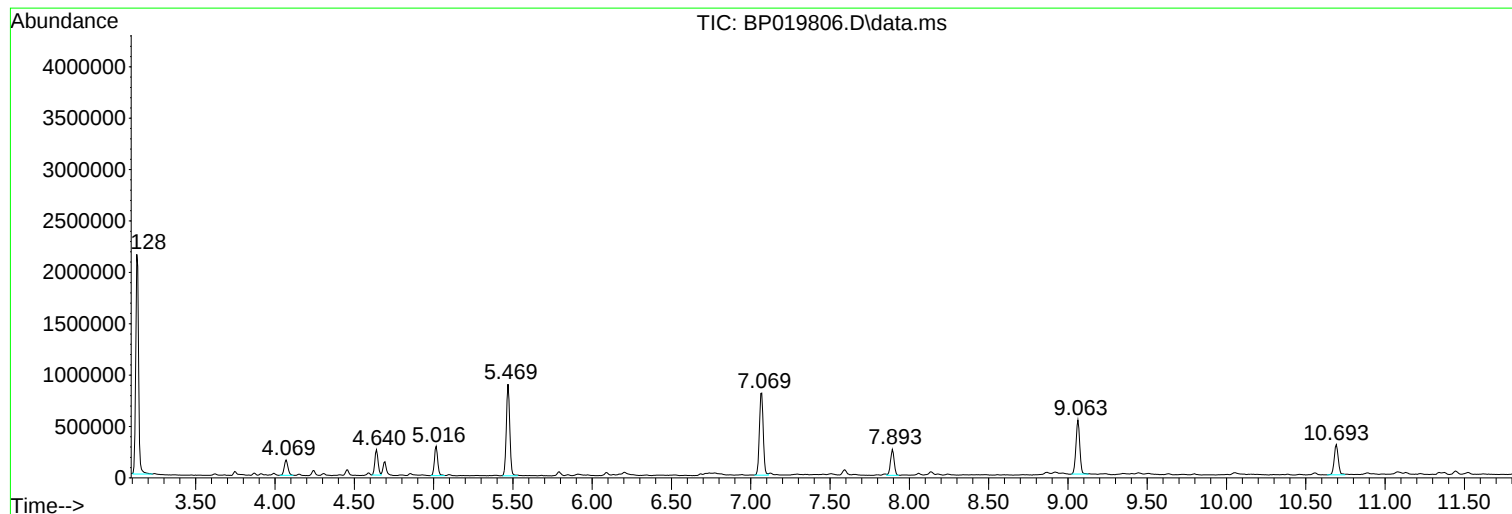
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.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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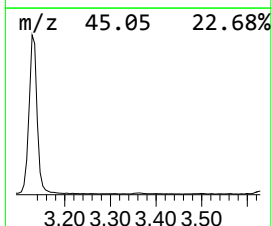
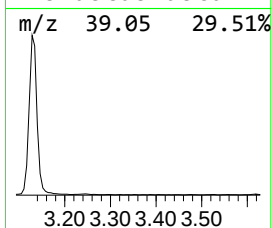
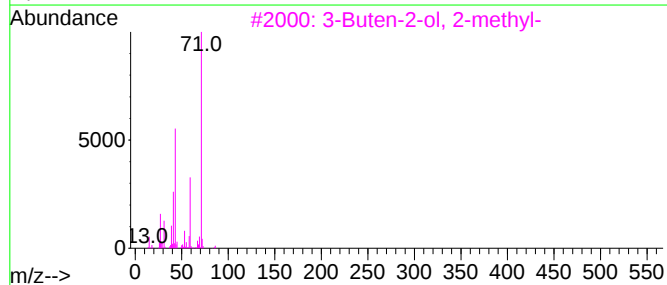
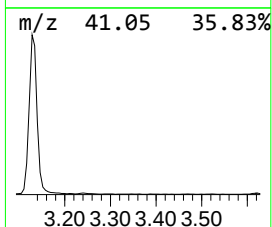
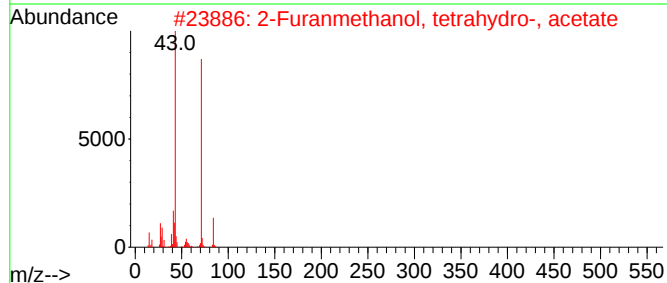
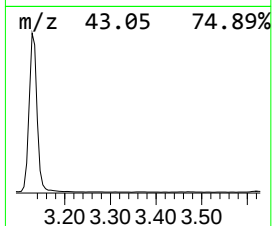
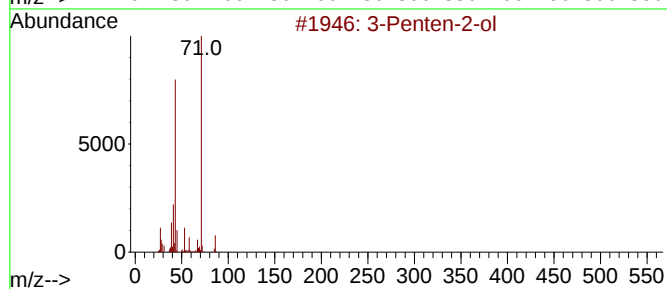
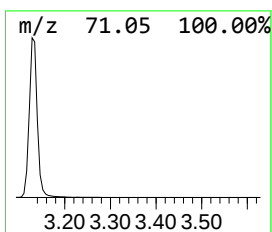
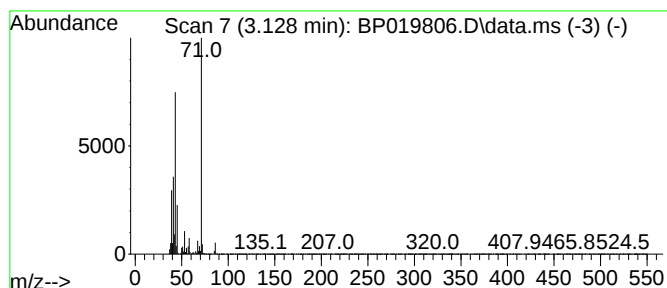
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	130.97 ng	2611410	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-ol	86	C5H10O	001569-50-2	80
2			2-Furanmethanol, tetrahydro-, ac...	144	C7H12O3	000637-64-9	53
3			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	53
4			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	50
5			Butanoic acid, anhydride	158	C8H14O3	000106-31-0	50



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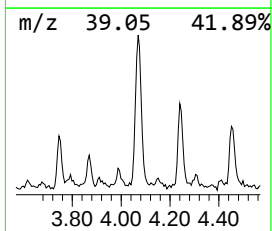
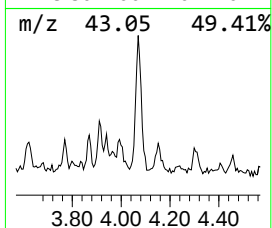
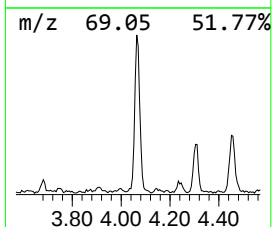
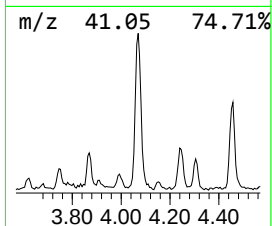
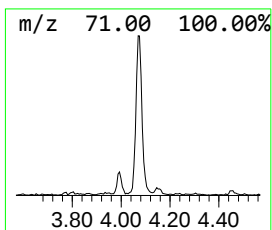
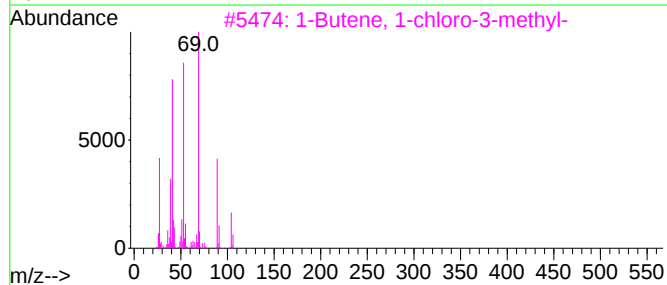
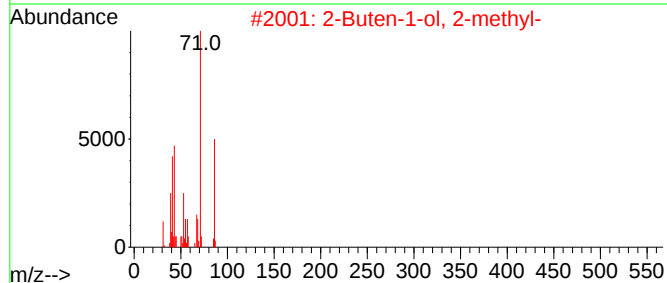
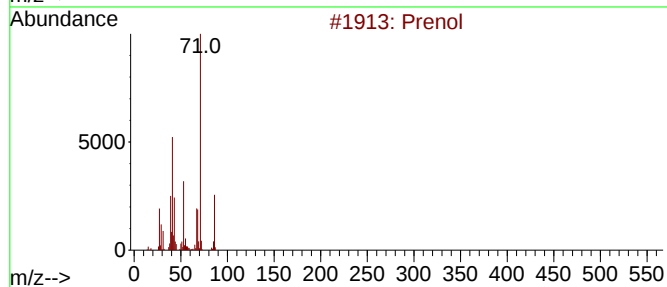
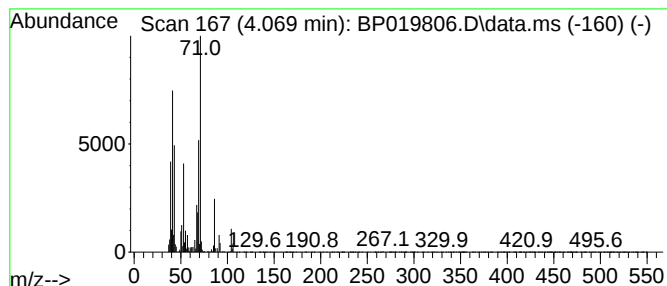
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Prenol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.069	12.39 ng	247101	1,4-Dichlorobenzene-d4	7.893

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Prenol	86	C5H10O	000556-82-1	81
2	2-Buten-1-ol, 2-methyl-	86	C5H10O	004675-87-0	74
3	1-Butene, 1-chloro-3-methyl-	104	C5H9Cl	023010-00-6	38
4	2-Butene, 1-methoxy-, (E)-	86	C5H10O	010034-14-7	27
5	2-Butyne, 1-methoxy-	84	C5H8O	002768-41-4	25



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Instrument :
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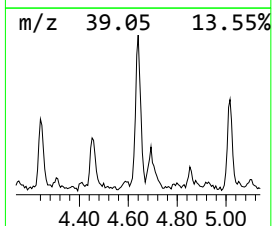
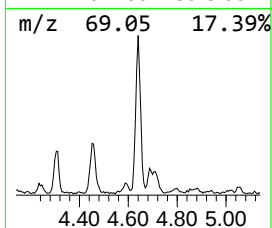
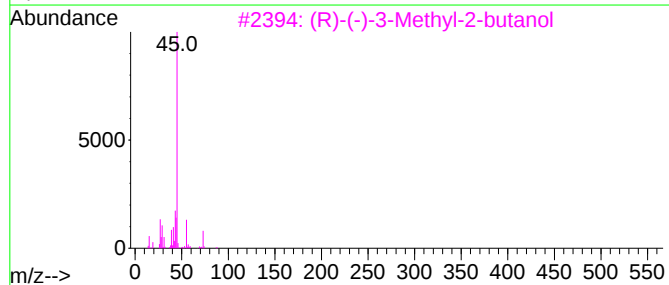
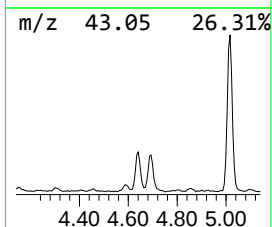
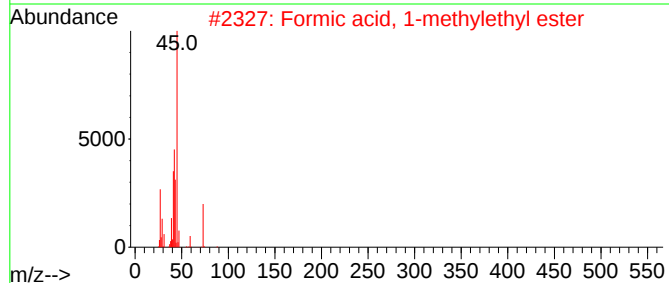
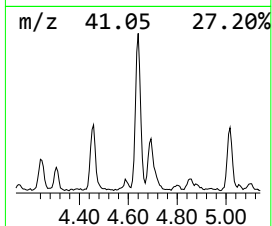
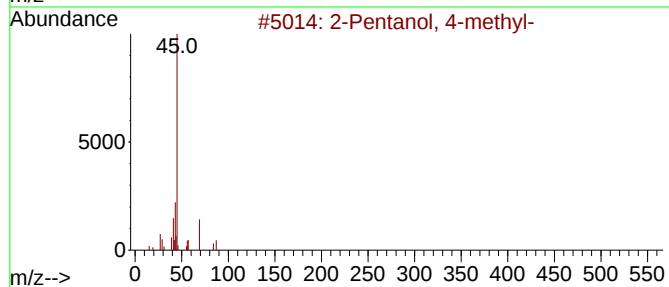
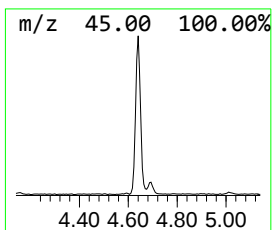
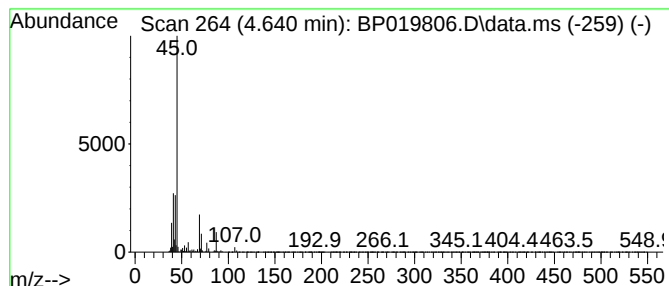
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanol, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.640	17.74 ng	353647	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	72
2			Formic acid, 1-methylethyl ester	88	C4H8O2	000625-55-8	50
3			(R)-(-)-3-Methyl-2-butanol	88	C5H12O	001572-93-6	50
4			2-Hexanol	102	C6H14O	000626-93-7	40
5			Propanenitrile, 3-methoxy-	85	C4H7NO	000110-67-8	40



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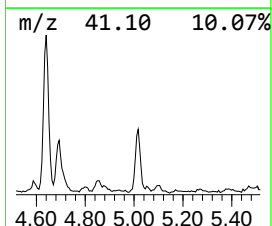
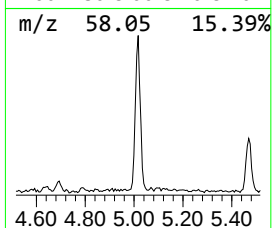
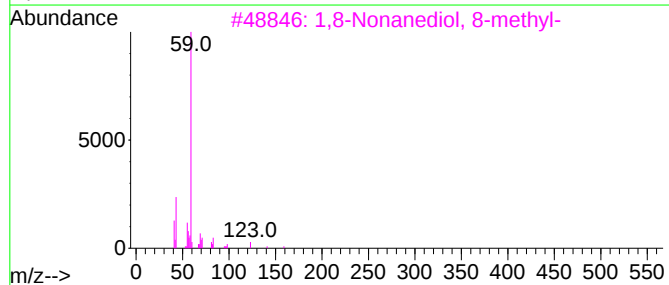
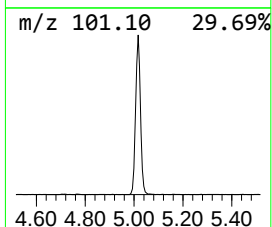
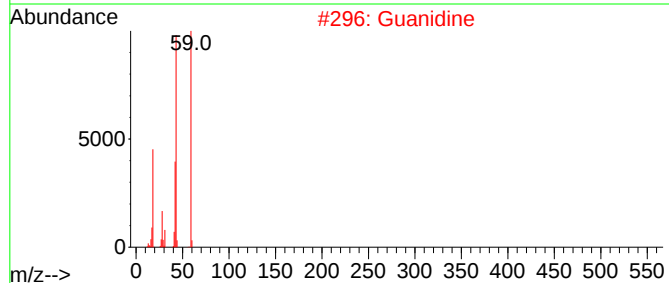
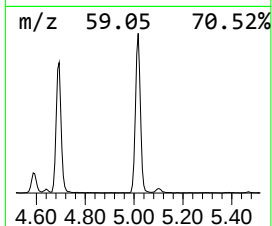
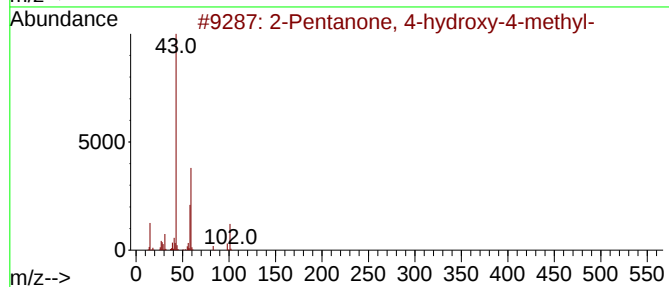
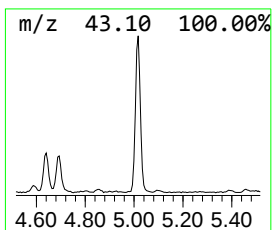
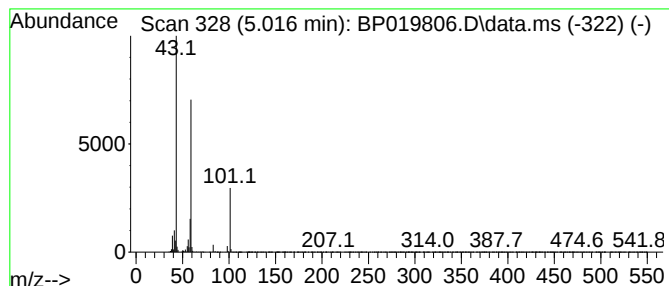
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.016	20.80 ng	414783	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Guanidine	59	CH5N3	000113-00-8	43
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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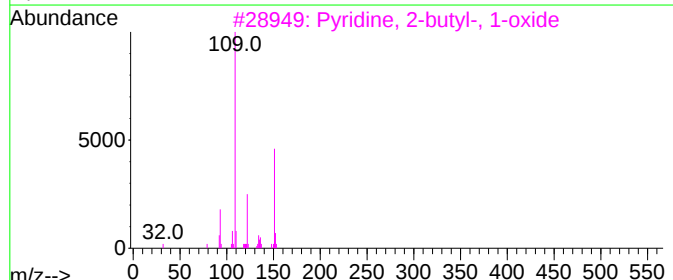
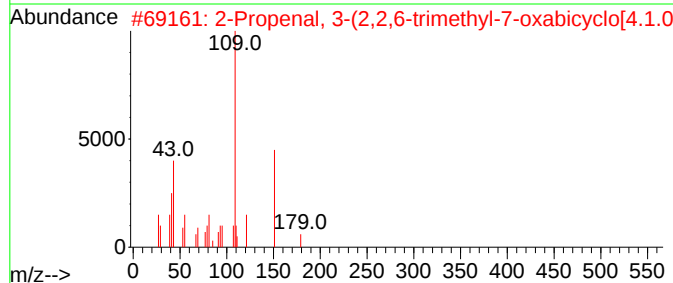
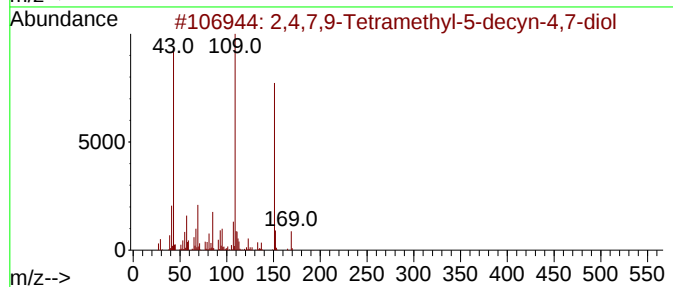
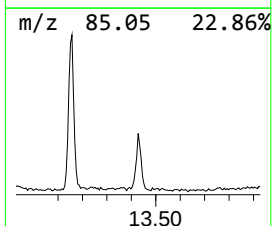
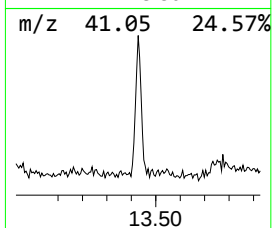
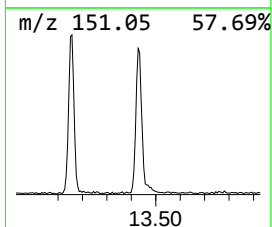
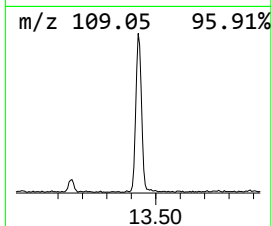
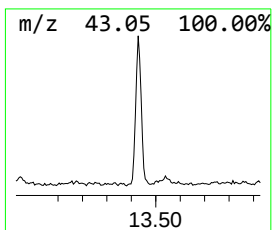
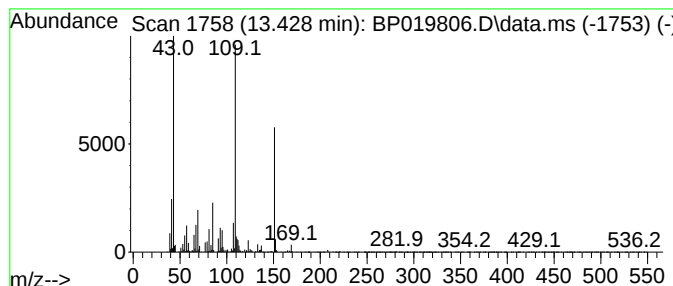
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Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.428	8.15 ng	279287	Acenaphthene-d10	14.528	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	87
2	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	59
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	53
4	Acetaminophen	151	C8H9NO2	000103-90-2	50
5	Tiocarlide	400	C23H32N2O2S	000910-86-1	50



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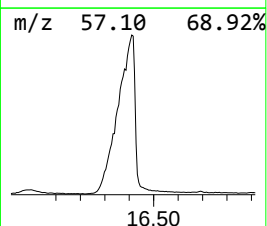
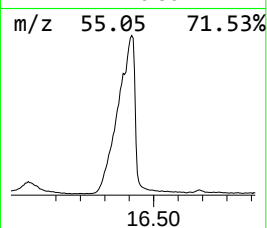
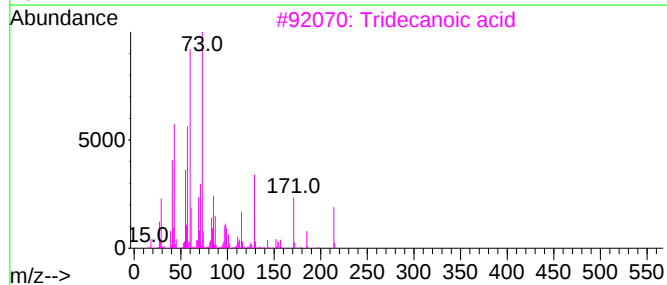
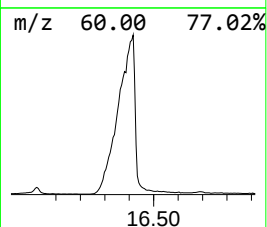
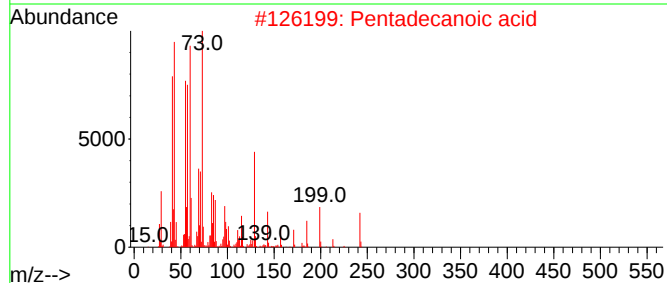
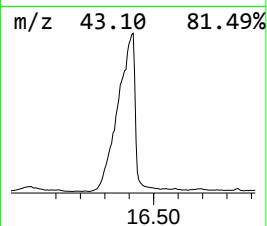
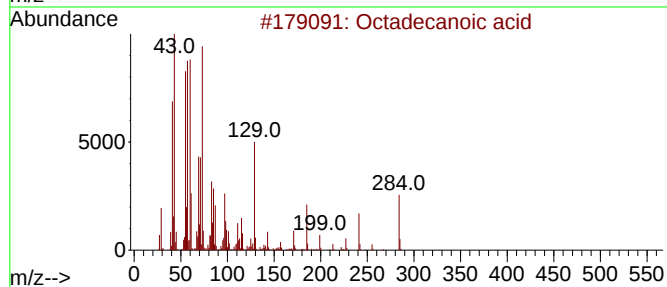
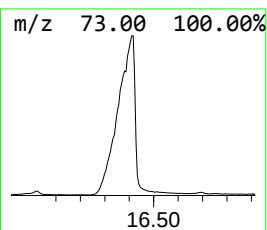
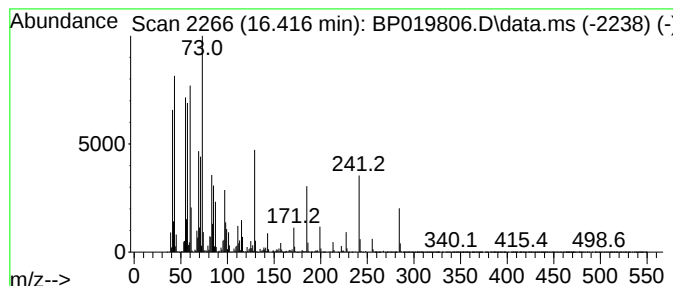
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.416	485.20 ng	20668500	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Tridecanoic acid	214	C13H26O2	000638-53-9	86
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	83
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

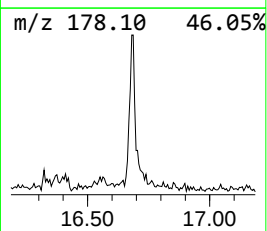
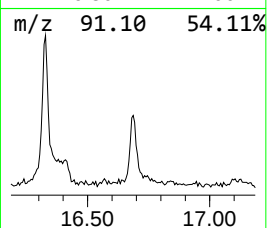
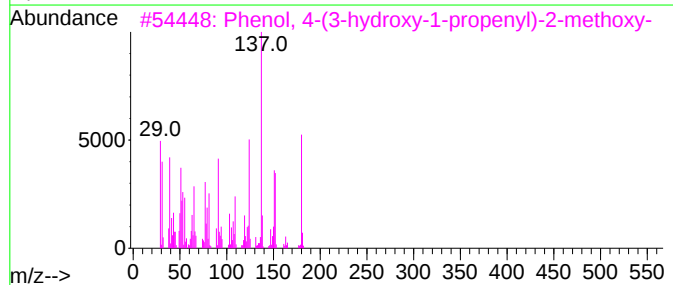
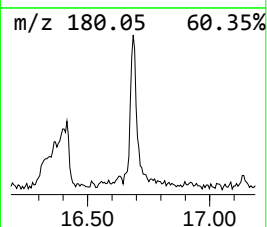
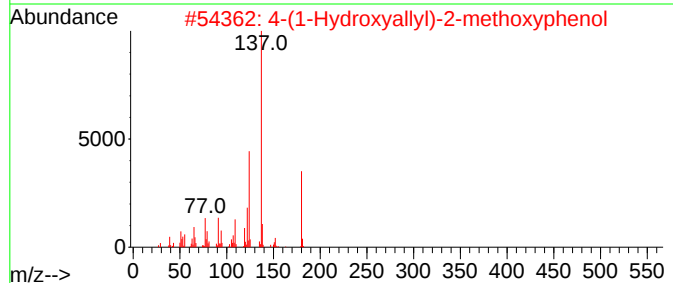
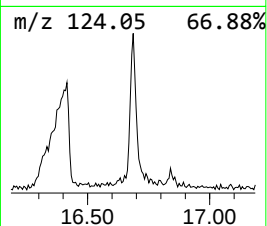
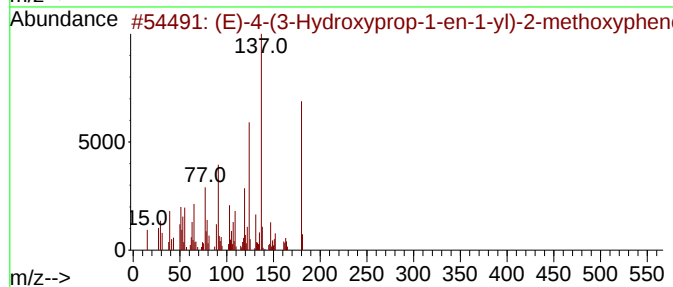
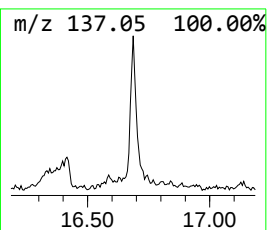
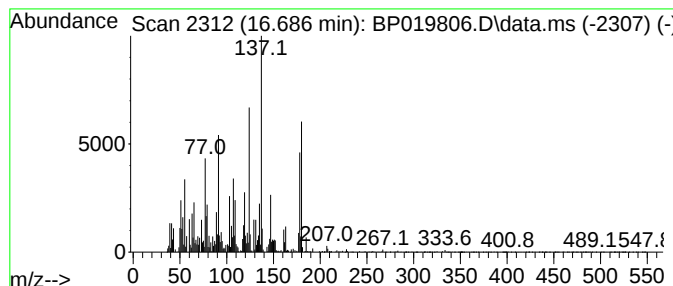
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 (E)-4-(3-Hydroxyprop-1-en-1-yl)-2-methoxyphenol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.686	6.37 ng	271457	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-4-(3-Hydroxyprop-1-en-1-yl)-2-methoxyphenol	180	C10H12O3	032811-40-8	86		
2	4-(1-Hydroxyallyl)-2-methoxyphenol	180	C10H12O3	112465-50-6	53		
3	Phenol, 4-(3-hydroxy-1-propenyl)-2-methoxy-	180	C10H12O3	000458-35-5	52		
4	3,7-Benzofurandiyl, 2,3-dihydro-1,4-benzodioxole-	180	C10H12O3	017781-15-6	43		
5	1-(Prop-2-ynyloxy)-3,4-methylenebis(2-methoxyphenyl)	178	C10H10O3	019202-22-3	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

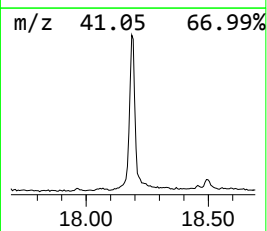
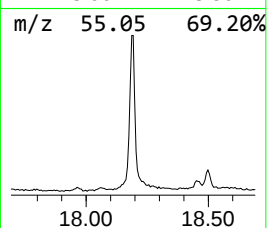
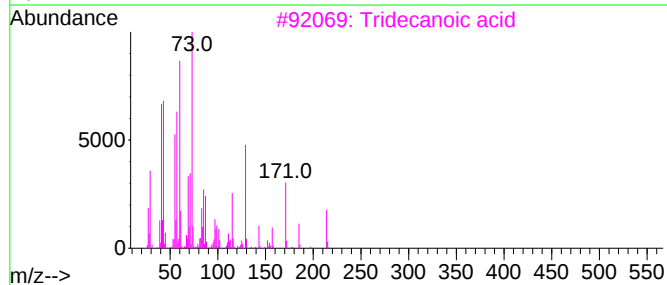
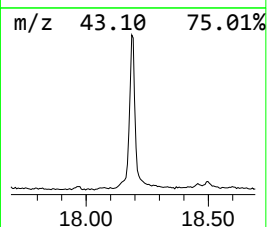
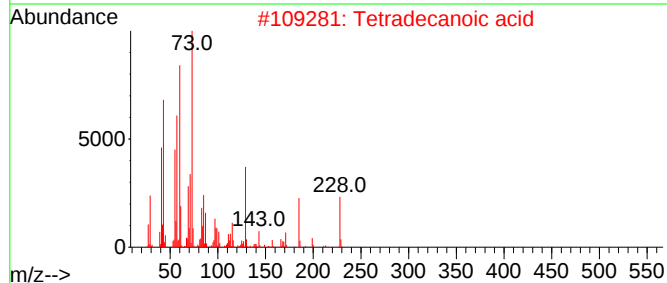
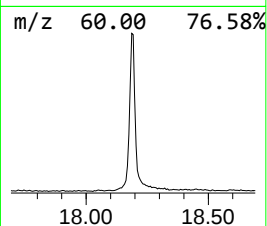
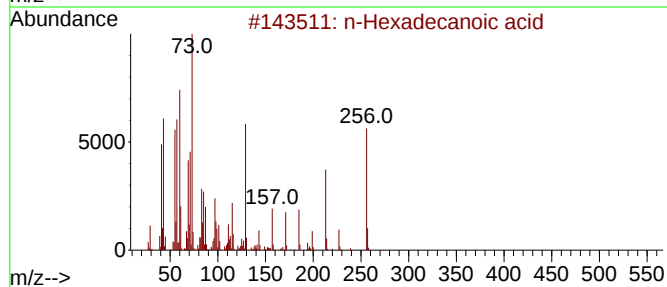
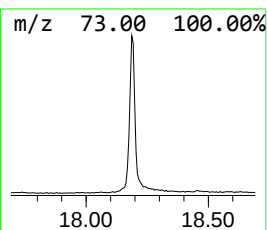
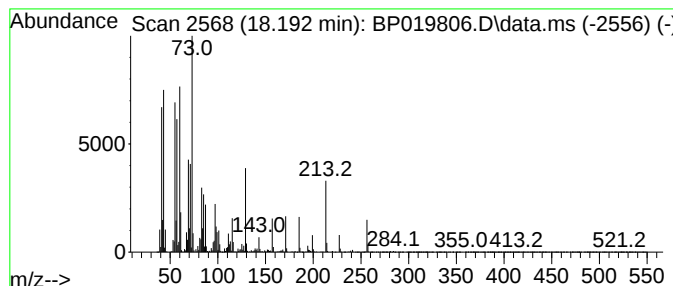
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	55.11 ng	2347480	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	95
3			Tridecanoic acid	214	C13H26O2	000638-53-9	91
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			Octadecanoic acid	284	C18H36O2	000057-11-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

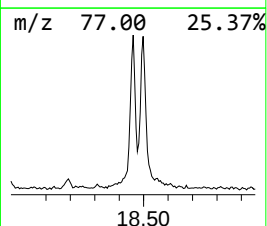
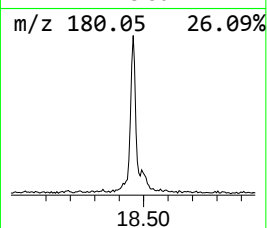
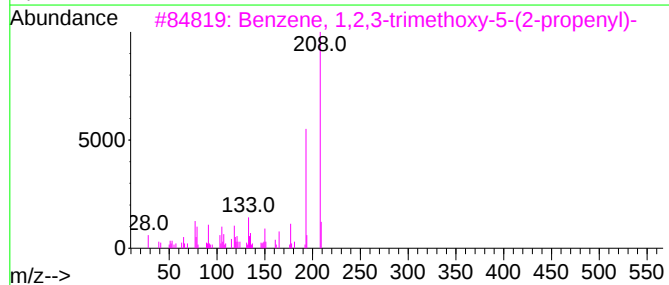
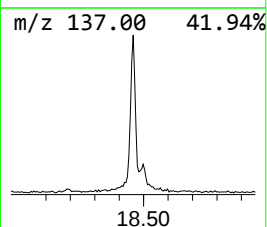
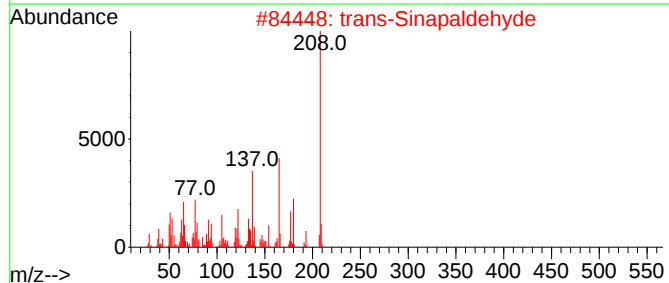
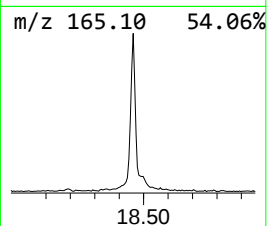
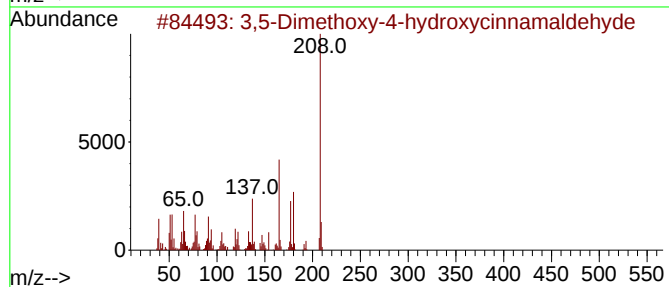
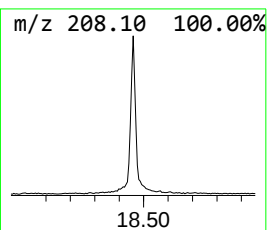
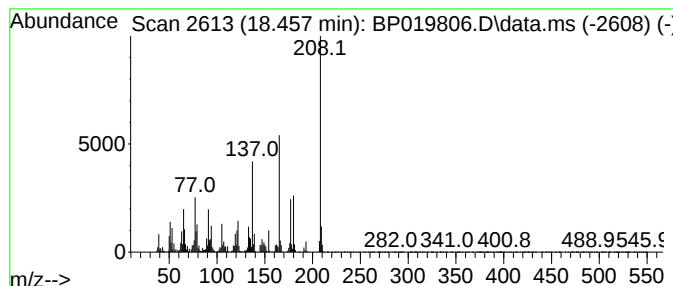
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.457	20.48 ng	872217	Phenanthrene-d10	17.286	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethoxy-4-hydroxycinnamal...	208	C11H12O4	087345-53-7	96
2	trans-Sinapaldehyde	208	C11H12O4	004206-58-0	93
3	Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	46
4	2-Propenoic acid, 3-(2,3-dimetho...	208	C11H12O4	007345-82-6	43
5	trans-2,5-Dimethoxycinnamic acid	208	C11H12O4	038489-74-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

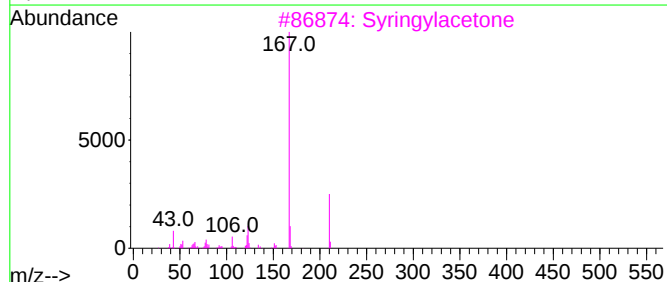
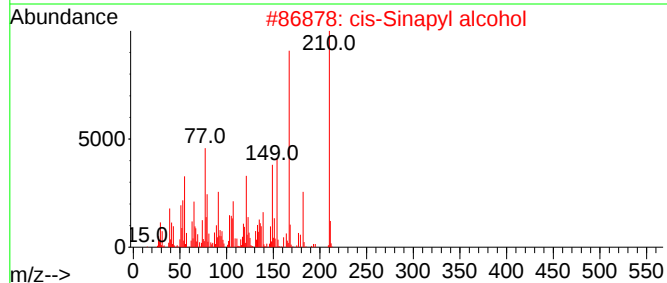
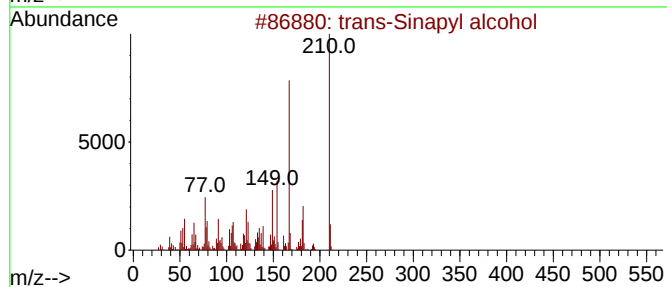
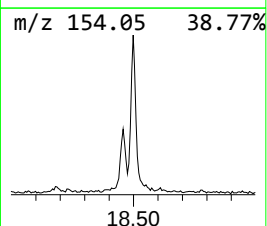
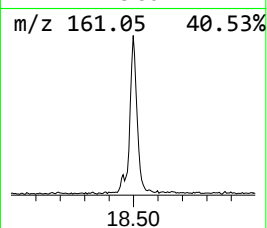
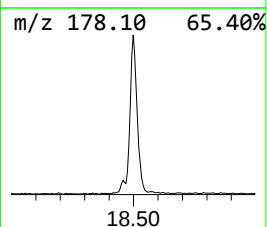
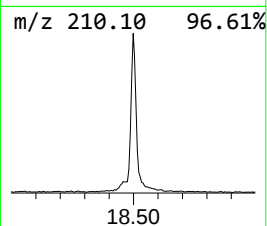
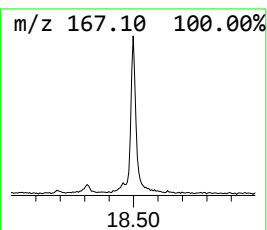
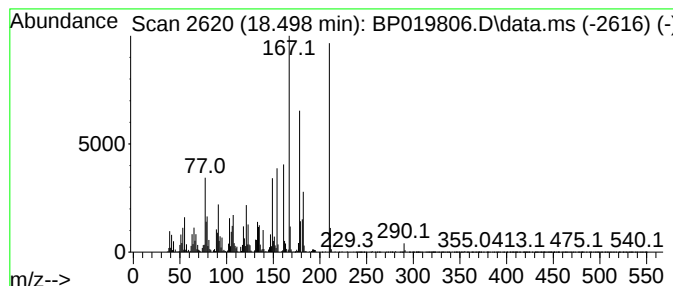
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 trans-Sinapyl alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.498	24.00 ng	1022170	Phenanthrene-d10	17.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		trans-Sinapyl alcohol	210	C11H14O4	020675-96-1	89
2		cis-Sinapyl alcohol	210	C11H14O4	104330-63-4	86
3		Syringylacetone	210	C11H14O4	019037-58-2	50
4		2-Pentanone, 1-(2,4,6-trihydroxy...	210	C11H14O4	1000116-22-3	38
5		4-Ethyl-2,6-dimethoxyphenol	182	C10H14O3	014059-92-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

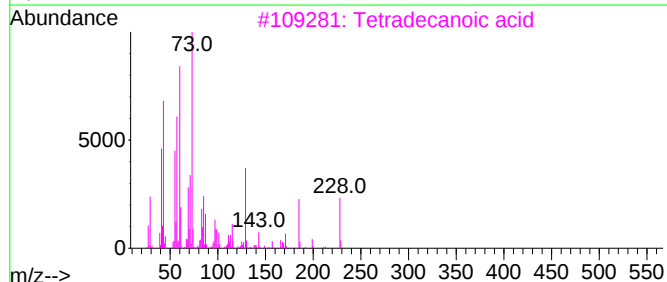
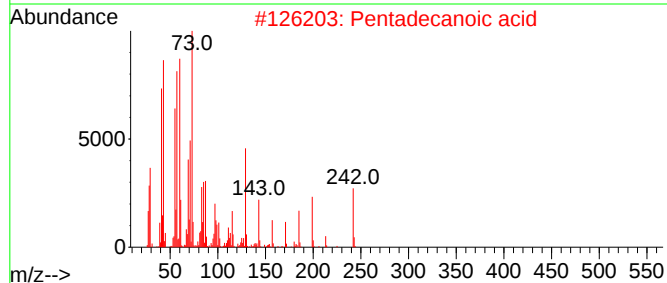
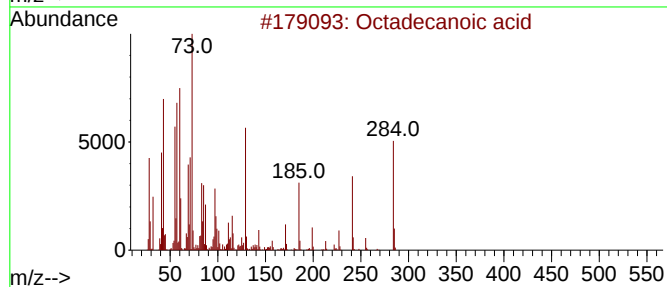
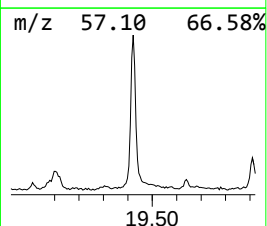
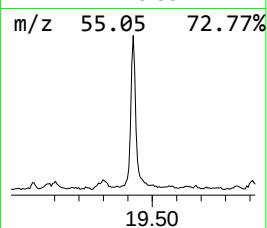
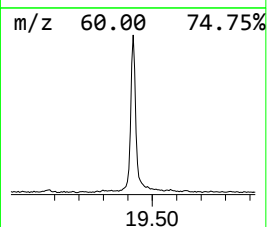
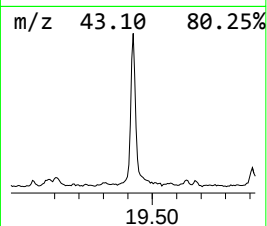
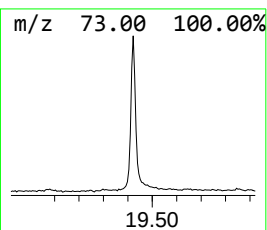
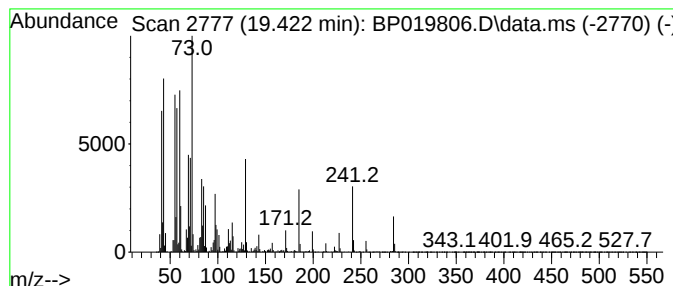
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.422	27.32 ng	1653260	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64
5			n-Decanoic acid	172	C10H20O2	000334-48-5	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

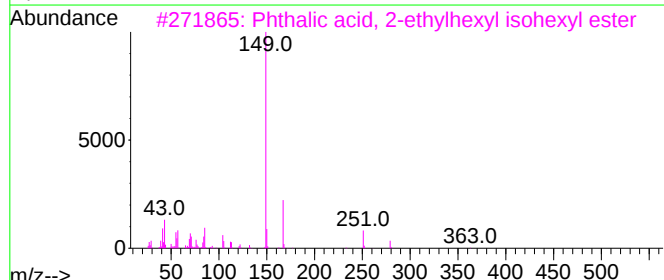
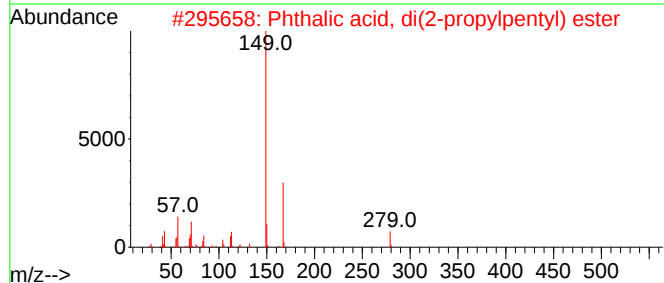
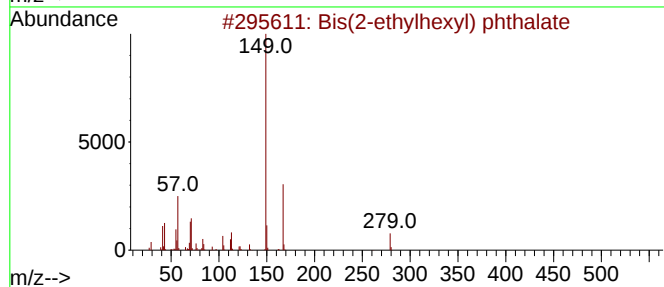
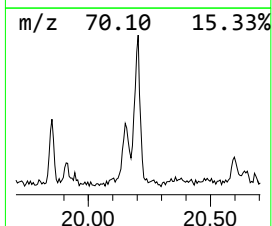
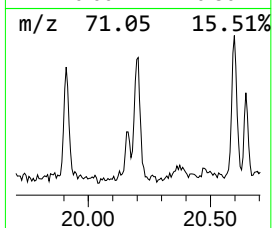
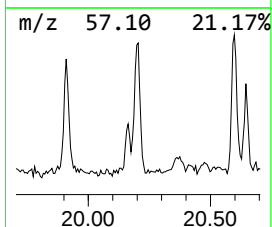
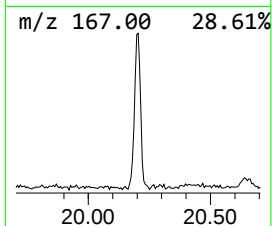
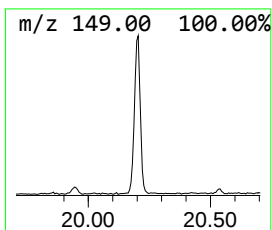
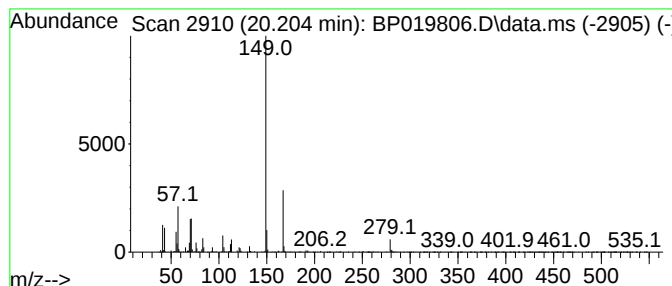
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Phthalic acid, di(2-propylp... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.204	5.77 ng	349363	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2			Phthalic acid, di(2-propylpentyl...	390	C24H38O4	1000377-93-5	90
3			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	90
4			Phthalic acid, di(oct-3-yl) ester	390	C24H38O4	1000377-72-3	86
5			Phthalic acid, 2-ethylhexyl hexy...	362	C22H34O4	1000309-02-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

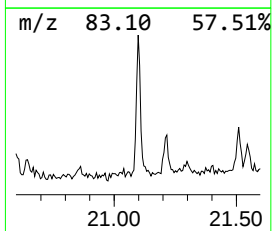
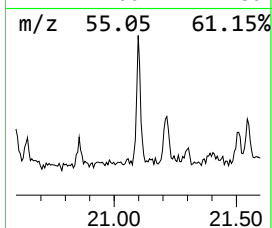
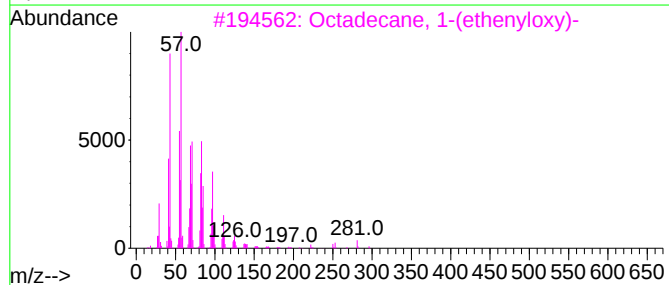
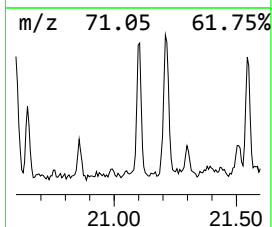
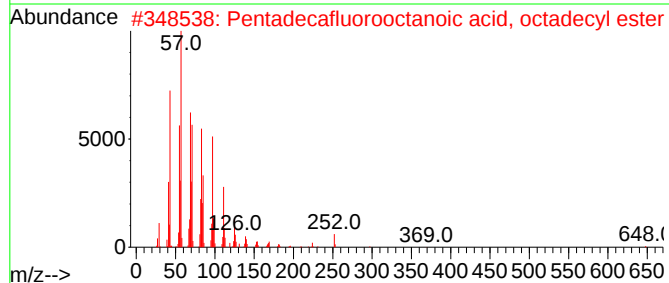
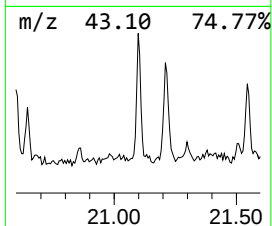
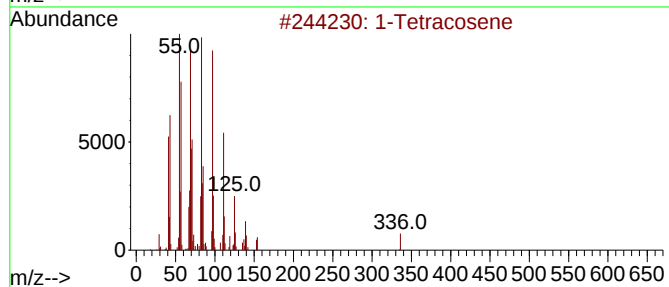
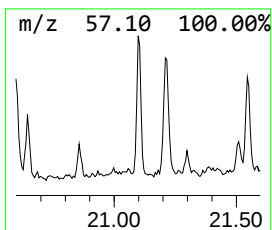
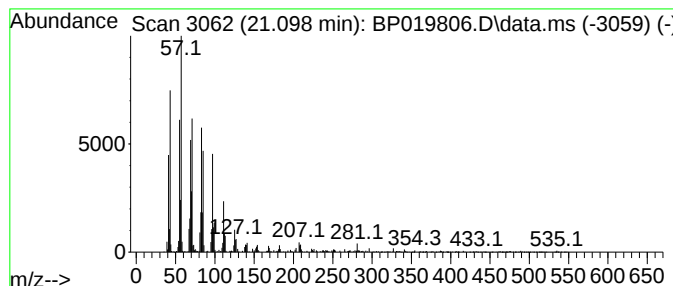
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Tetracosene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	3.51 ng	212104	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tetracosene	336	C24H48	010192-32-2	96
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
3			Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	90
4			7-Hexadecene, (Z)-	224	C16H32	035507-09-6	89
5			Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

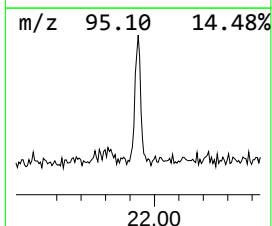
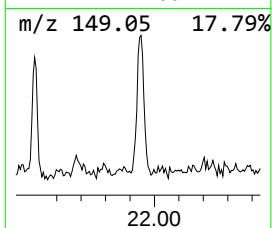
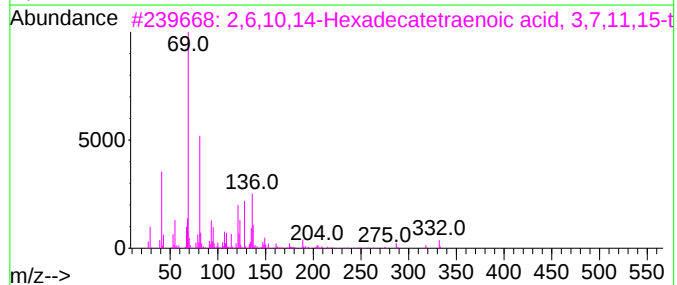
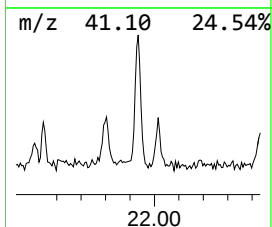
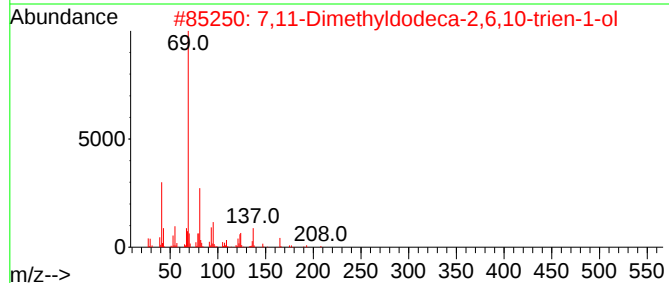
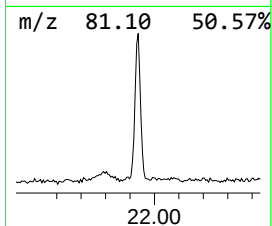
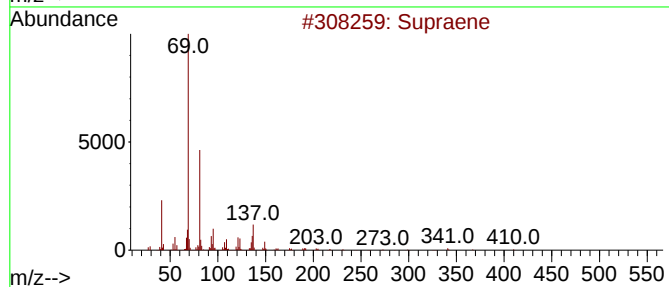
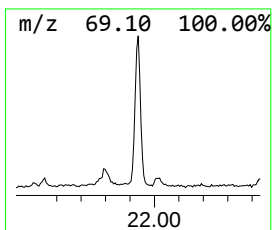
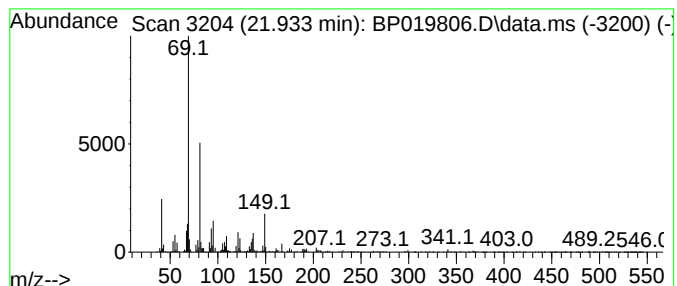
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Supraene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.933	6.31 ng	381737	Chrysene-d12	21.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Supraene	410	C30H50	007683-64-9	72
2		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
3		2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	59
4		3-Ethyl-1,5-octadiene	138	C10H18	1000114-87-7	53
5		4-Methyl-1,5-Heptadiene	110	C8H14	000998-94-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019806.D
Acq On : 21 Feb 2024 14:05
Operator : MA/JU
Sample : P1523-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240763-COMPOSITE-39N-N-3-01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-ol	3.128	131.0	ng	2611410	1	7.893	398771	20.0
Prenol	4.069	12.4	ng	247101	1	7.893	398771	20.0
2-Pentanol, 4-m...	4.640	17.7	ng	353647	1	7.893	398771	20.0
2-Pentanone, 4-...	5.016	20.8	ng	414783	1	7.893	398771	20.0
2,4,7,9-Tetrame...	13.428	8.2	ng	279287	3	14.528	685396	20.0
Octadecanoic acid	16.416	485.2	ng	20668500	4	17.286	851952	20.0
(E)-4-(3-Hydrox...	16.686	6.4	ng	271457	4	17.286	851952	20.0
n-Hexadecanoic ...	18.192	55.1	ng	2347480	4	17.286	851952	20.0
3,5-Dimethoxy-4...	18.457	20.5	ng	872217	4	17.286	851952	20.0
trans-Sinapyl a...	18.498	24.0	ng	1022170	4	17.286	851952	20.0
Pentadecanoic acid	19.422	27.3	ng	1653260	5	21.380	1210250	20.0
Phthalic acid, ...	20.204	5.8	ng	349363	5	21.380	1210250	20.0
1-Tetracosene	21.098	3.5	ng	212104	5	21.380	1210250	20.0
Supraene	21.933	6.3	ng	381737	5	21.380	1210250	20.0