

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
 Data File : BF139941.D
 Acq On : 22 Oct 2024 21:20
 Operator : RC/JU
 Sample : P4385-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1

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-BF101824.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139941.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.187	7	16	30	rVB	2027002	3279149	99.73%	11.868%
2	5.110	505	513	522	rVB	150442	225704	6.86%	0.817%
3	5.504	573	580	585	rBV	2151430	2836162	86.26%	10.265%
4	6.504	744	750	763	rVB	2145012	2968756	90.29%	10.745%
5	6.887	810	815	820	rBV	669818	832153	25.31%	3.012%
6	7.445	904	910	915	rBV	1501687	1928927	58.67%	6.981%
7	7.698	949	953	958	rBV	86622	102736	3.12%	0.372%
8	8.169	1028	1033	1038	rBV	874857	1077443	32.77%	3.899%
9	8.381	1063	1069	1077	rVB	37634	50672	1.54%	0.183%
10	9.245	1210	1216	1221	rBV	2646168	3288006	100.00%	11.900%
11	9.922	1326	1331	1336	rBV	847517	1098305	33.40%	3.975%
12	10.633	1447	1452	1459	rBV	677761	882492	26.84%	3.194%
13	10.710	1460	1465	1475	rVV	1005821	1265856	38.50%	4.581%
14	10.945	1501	1505	1510	rVB	126416	152533	4.64%	0.552%
15	11.410	1579	1584	1591	rBV	722727	897512	27.30%	3.248%
16	11.933	1669	1673	1688	rVB2	199905	301586	9.17%	1.092%
17	12.433	1754	1758	1765	rVB	28848	40637	1.24%	0.147%
18	12.722	1802	1807	1815	rBV2	30349	66942	2.04%	0.242%
19	12.992	1848	1853	1859	rBV	1897464	2522693	76.72%	9.130%
20	13.227	1886	1893	1896	rBV3	22722	48968	1.49%	0.177%
21	13.898	2002	2007	2022	rBV	183889	362160	11.01%	1.311%
22	14.045	2027	2032	2041	rVB	648172	860025	26.16%	3.113%
23	14.527	2110	2114	2126	rBV4	36680	93159	2.83%	0.337%
24	14.980	2186	2191	2198	rVB	97141	139426	4.24%	0.505%
25	15.174	2220	2224	2226	rBV	71840	102232	3.11%	0.370%
26	15.204	2226	2229	2250	rVB2	101414	285476	8.68%	1.033%
27	15.527	2278	2284	2296	rVB	550334	899329	27.35%	3.255%
28	15.774	2321	2326	2336	rVB	58995	102481	3.12%	0.371%
29	16.015	2361	2367	2372	rBV	160794	285667	8.69%	1.034%
30	16.074	2372	2377	2387	rVV2	49582	137826	4.19%	0.499%
31	16.833	2502	2506	2513	rVB2	32447	61370	1.87%	0.222%
32	17.221	2567	2572	2597	rVB2	36964	171859	5.23%	0.622%
33	17.639	2636	2643	2660	rBV4	78691	262088	7.97%	0.949%

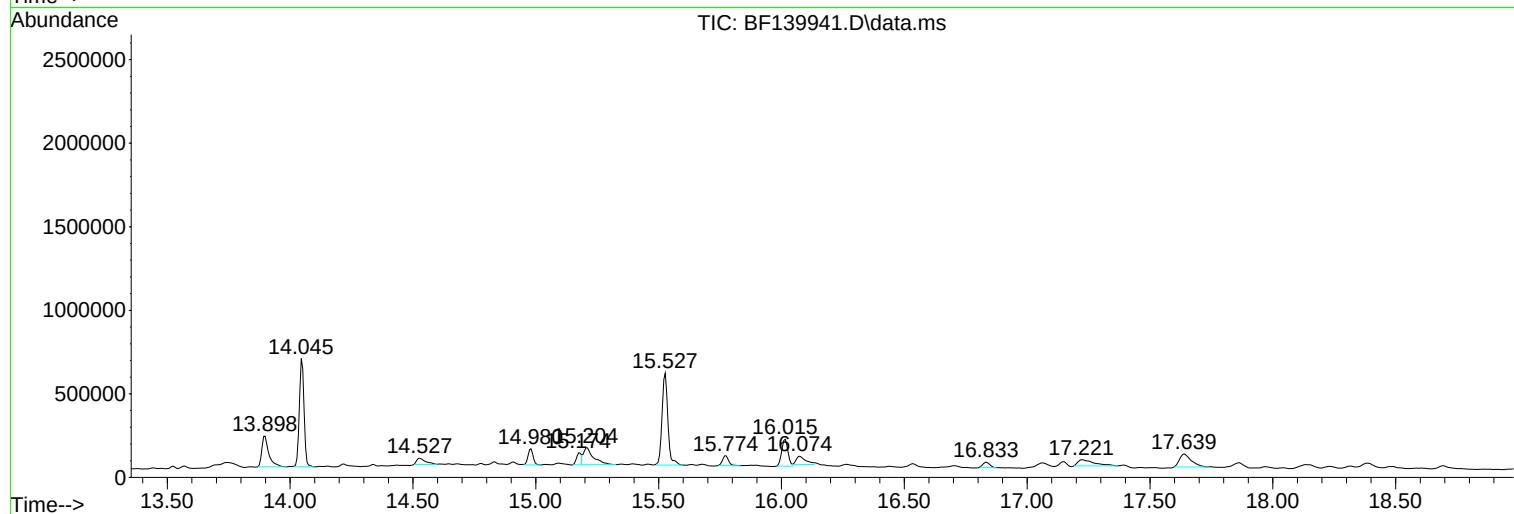
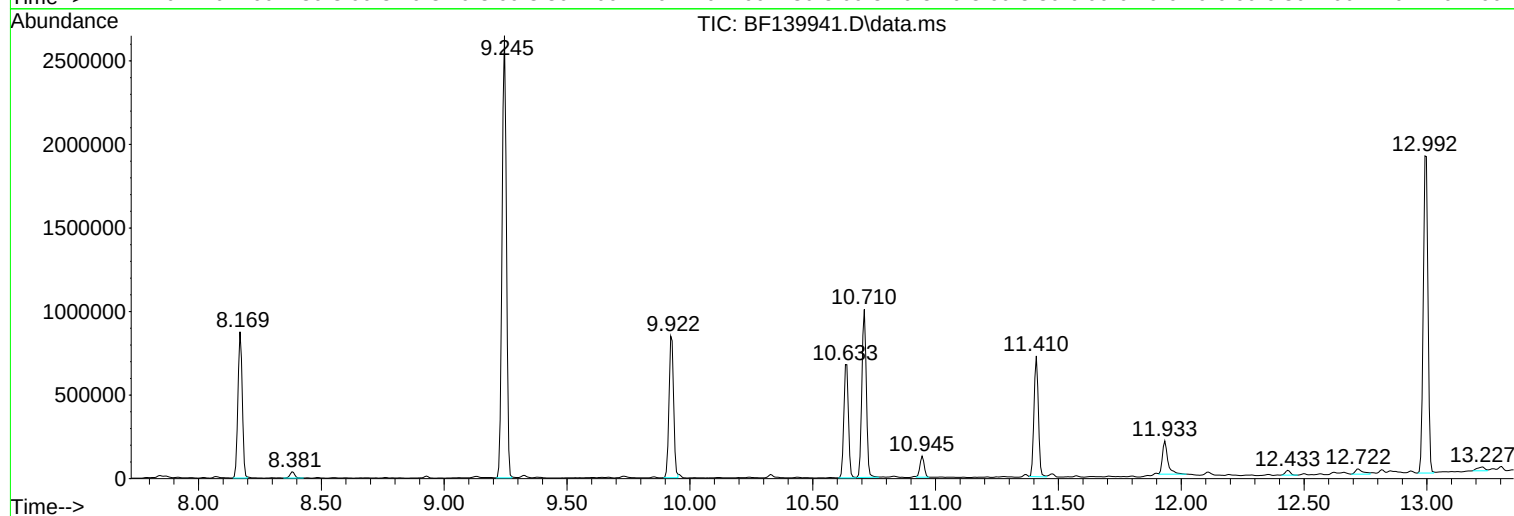
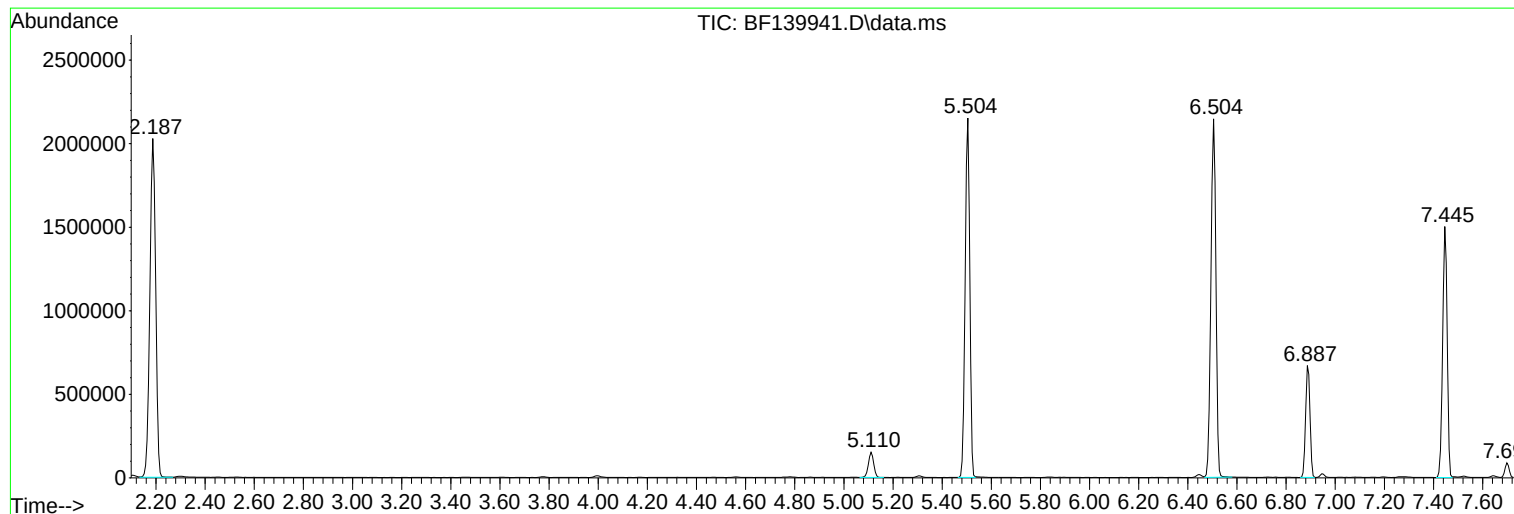
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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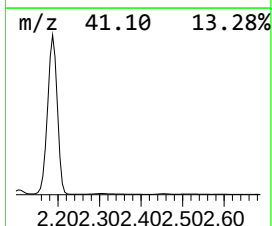
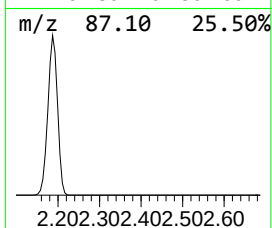
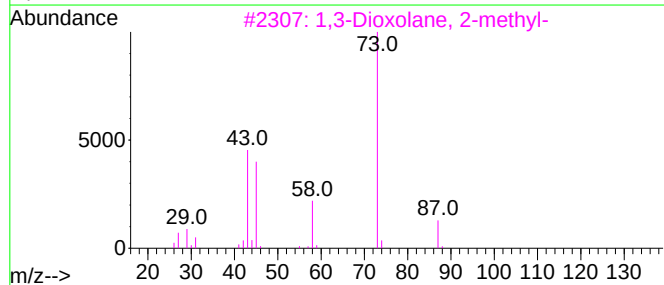
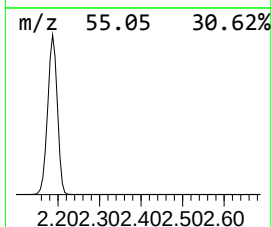
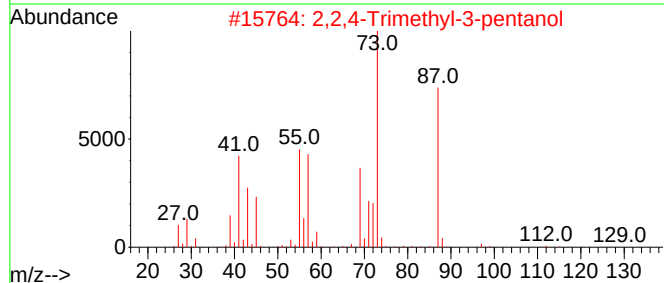
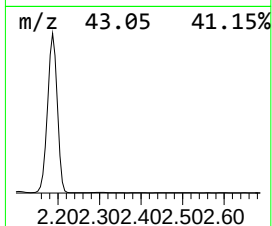
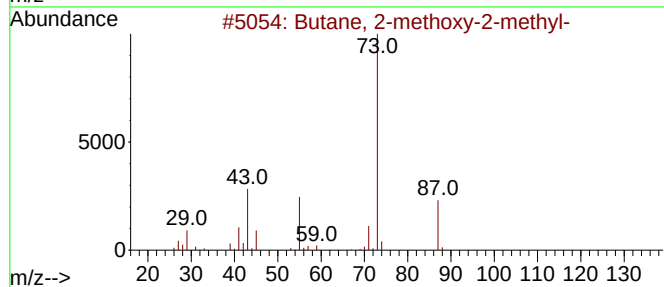
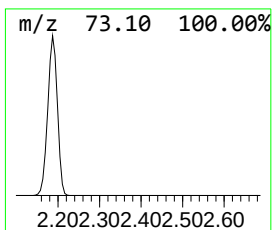
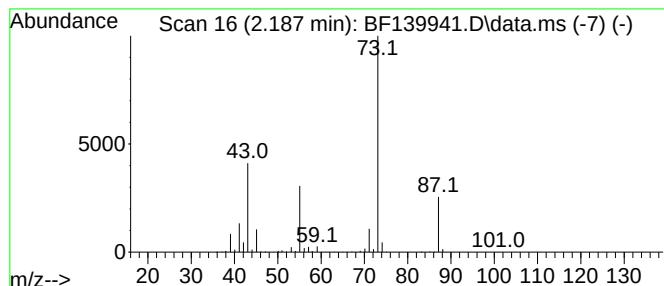
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.187	78.81 ng	3279150	1,4-Dichlorobenzene-d4	6.887

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	40
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17



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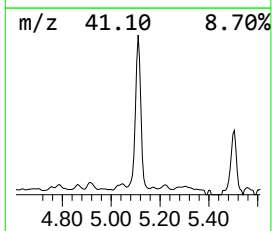
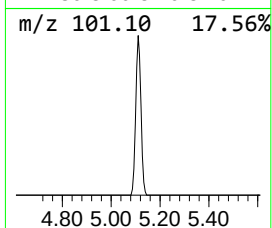
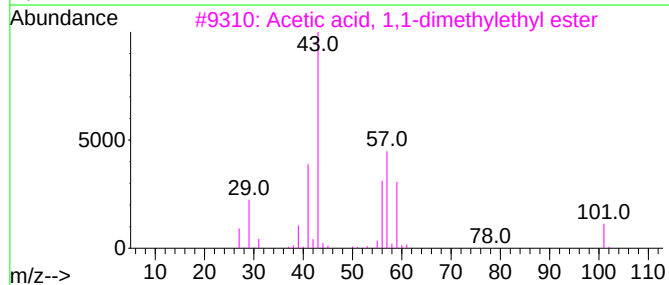
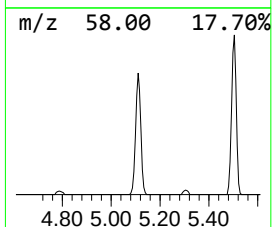
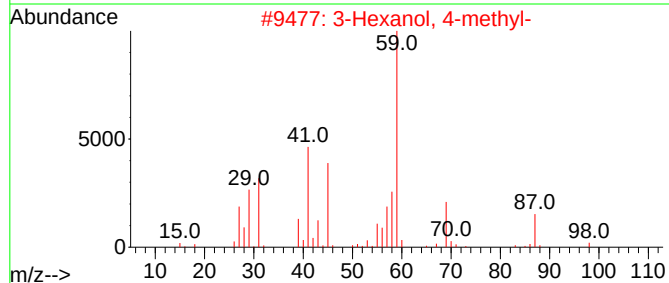
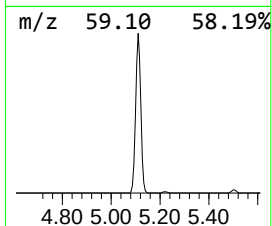
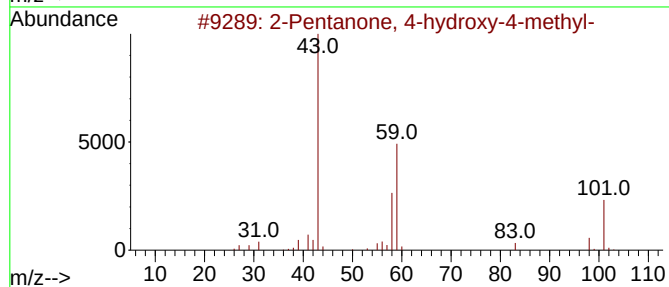
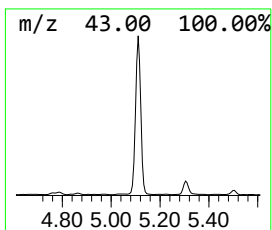
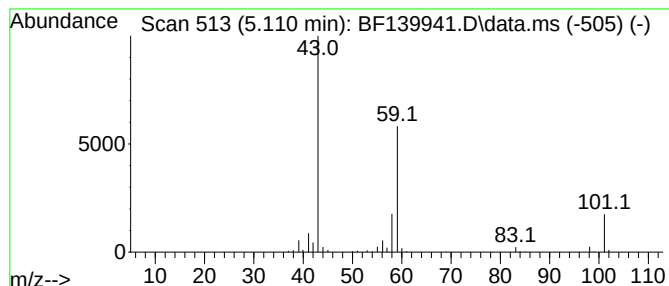
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.110	5.42 ng	225704	1,4-Dichlorobenzene-d4	6.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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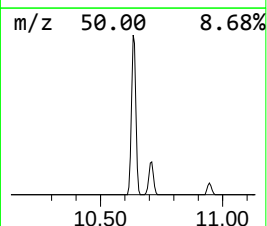
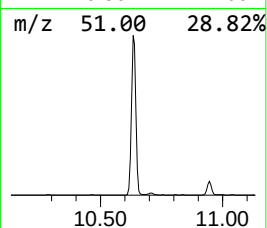
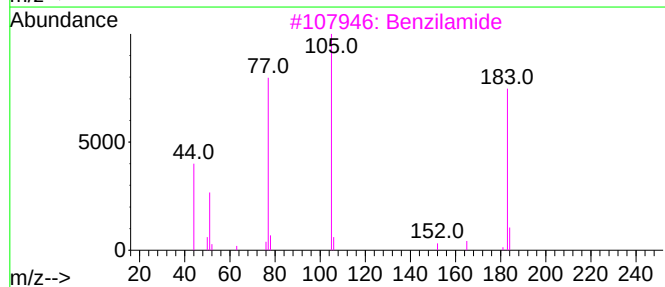
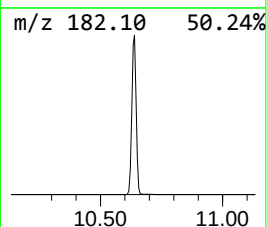
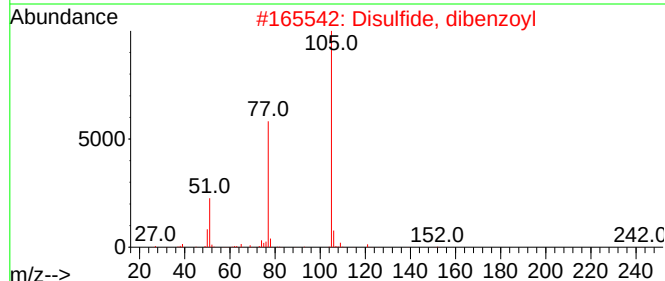
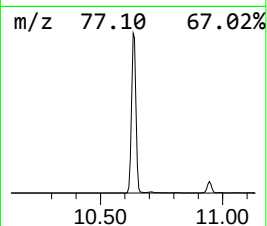
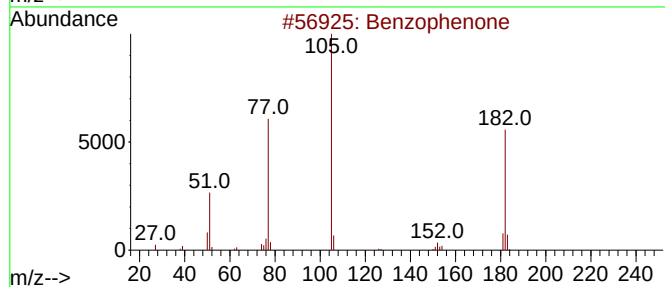
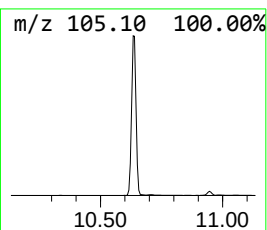
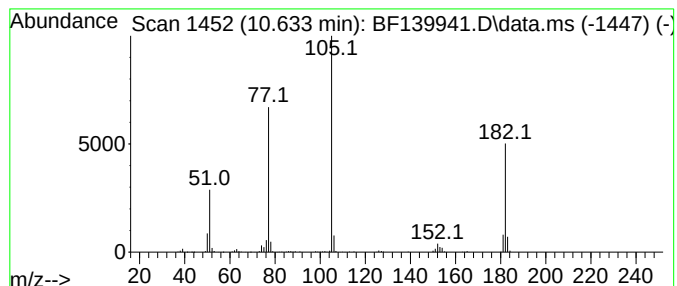
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Peak Number 3 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.633	16.07 ng	882492	Acenaphthene-d10	9.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Disulfide, dibenzoyl	274	C14H10O2S2	000644-32-6	47
3			Benzilamide	227	C14H13NO2	004746-87-6	47
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	47
5			Benzene, (trimethoxymethyl)-	182	C10H14O3	000707-07-3	47



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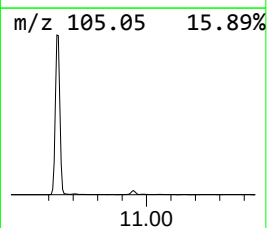
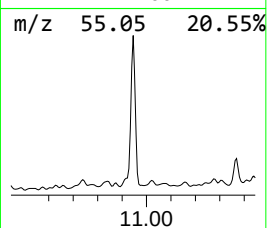
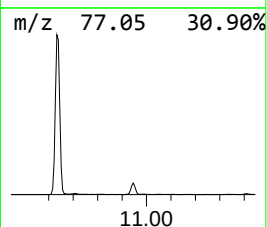
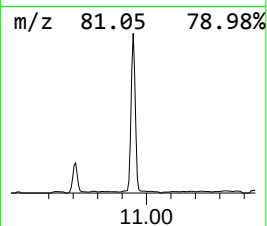
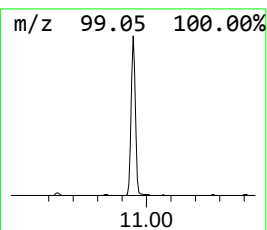
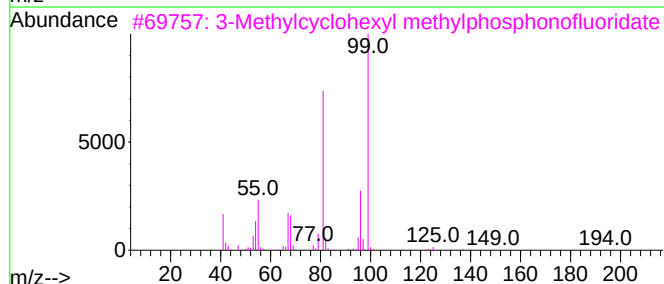
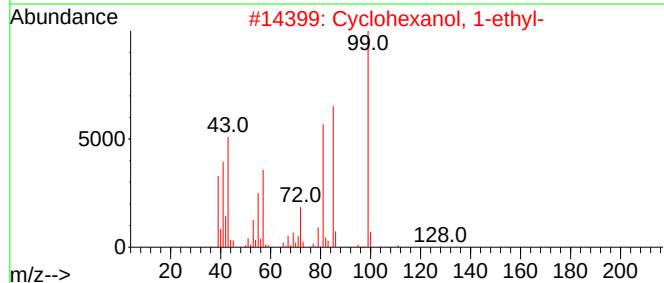
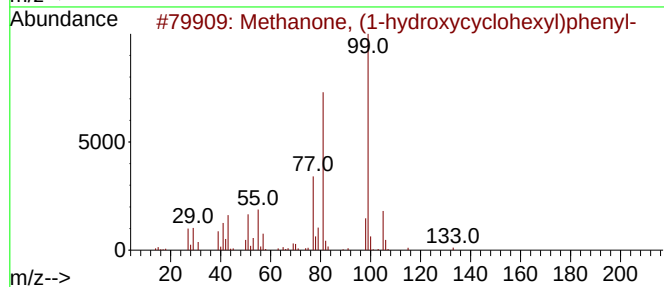
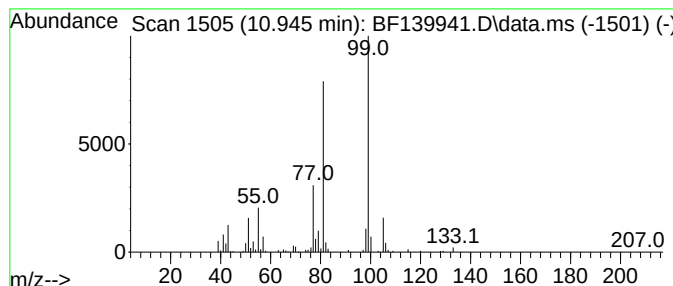
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Peak Number 4 Methanone, (1-hydroxycyclohexyl) Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.945	3.40 ng	152533	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	91
2			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	59
3			3-Methylcyclohexyl methylphospho...	194	C8H16FO2P	113548-86-0	59
4			1-Hydroxycyclohexanecarboxylic acid	144	C7H12O3	001123-28-0	45
5			2-Methylcyclosarin	194	C8H16FO2P	085473-32-1	38



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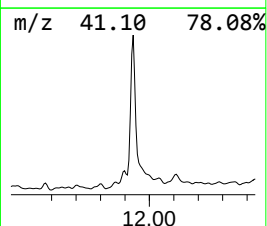
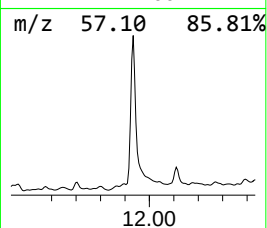
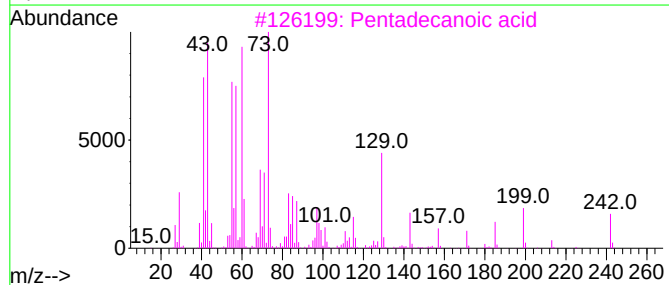
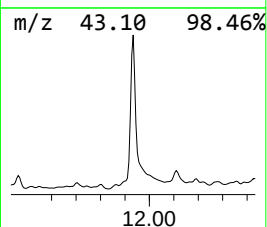
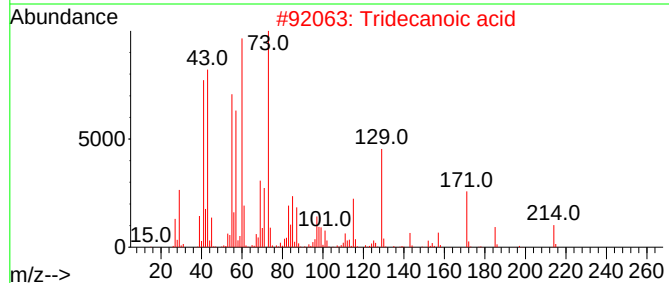
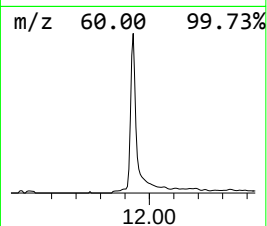
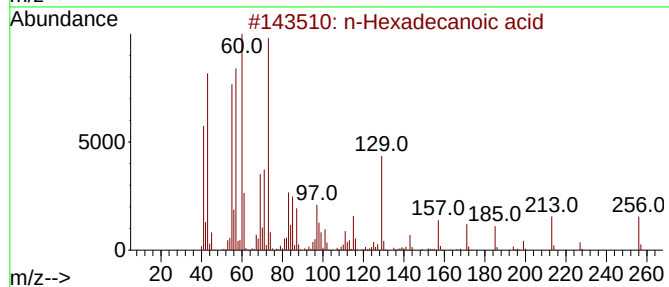
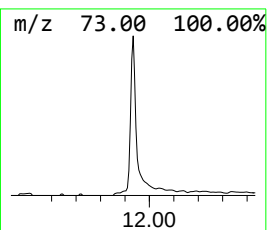
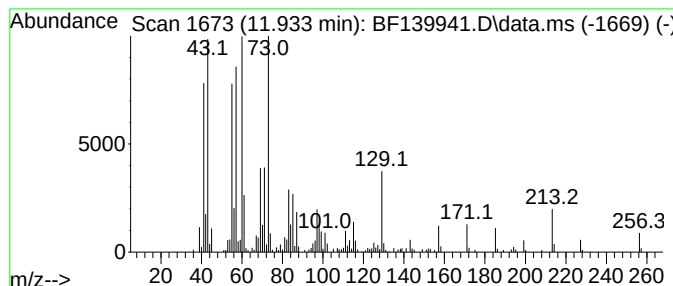
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	6.72 ng	301586	Phenanthrene-d10	11.410

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	93
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	78



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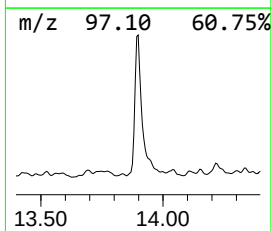
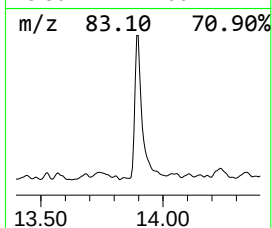
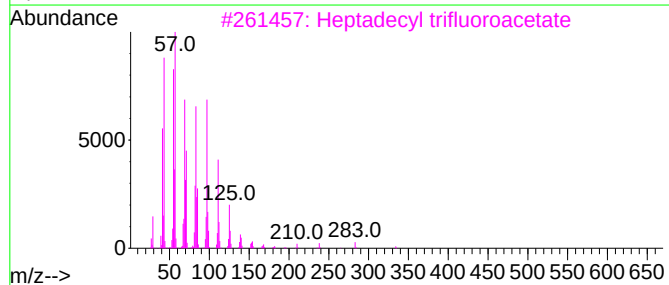
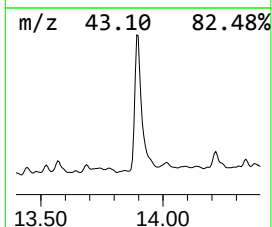
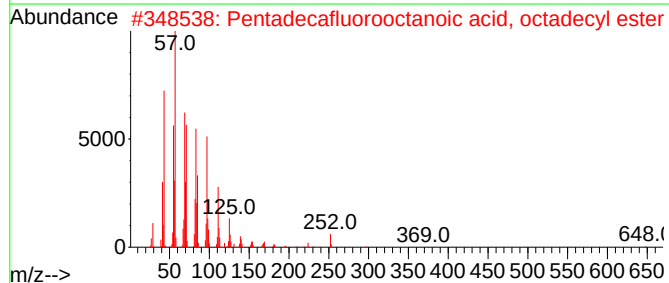
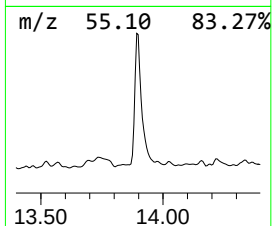
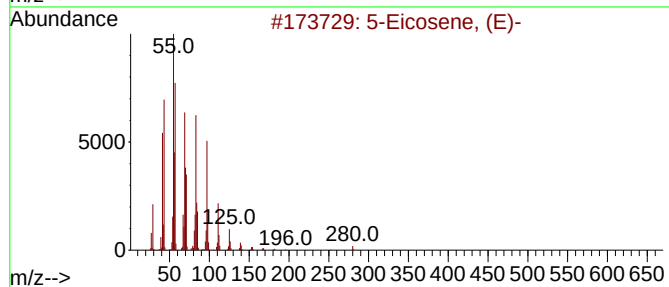
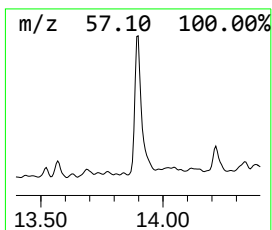
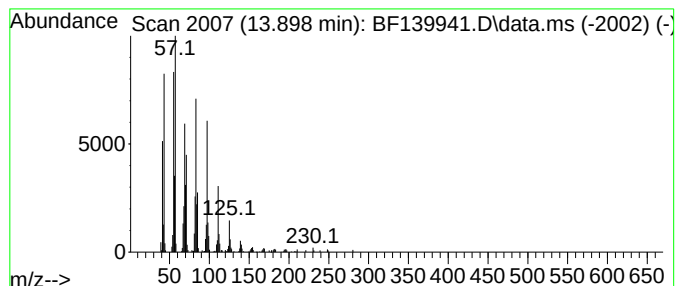
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 5-Eicosene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.898	8.42 ng	362160	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	95
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
4			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

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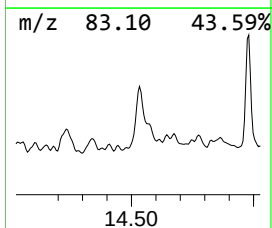
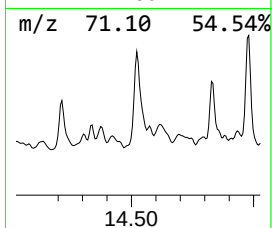
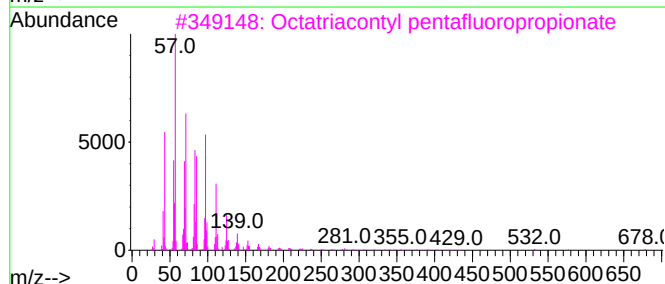
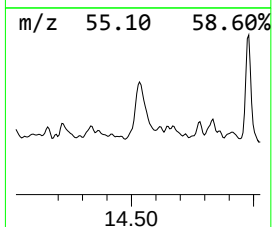
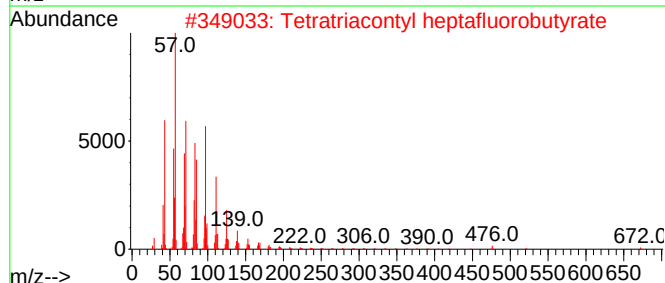
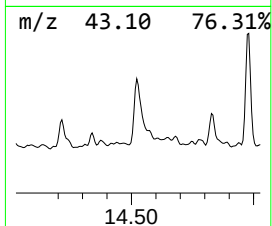
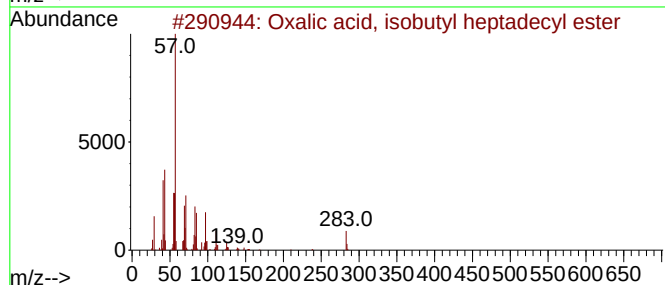
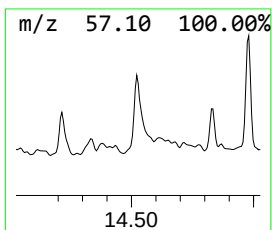
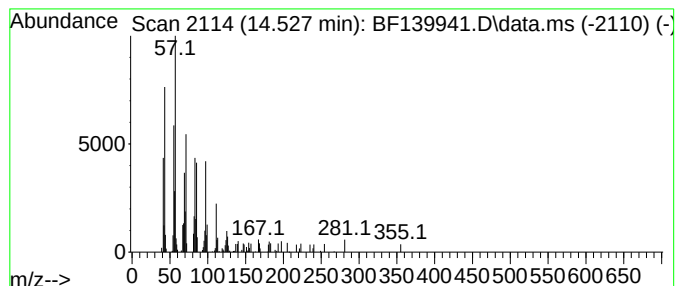
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Oxalic acid, isobutyl hepta... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.527	2.17 ng	93159	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	83
2			Tetratriacontyl heptafluorobutyrate	691	C38H69F7O2	1000351-84-1	83
3			Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	83
4			Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1010351-81-4	83
5			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

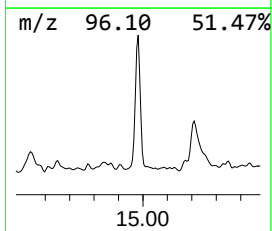
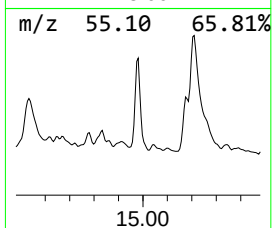
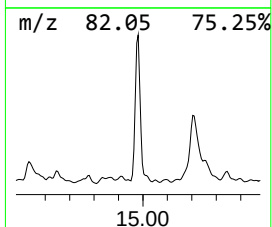
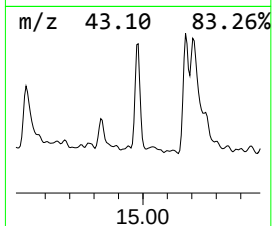
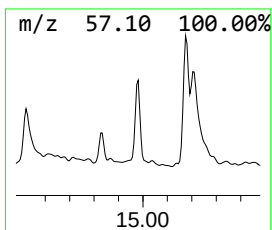
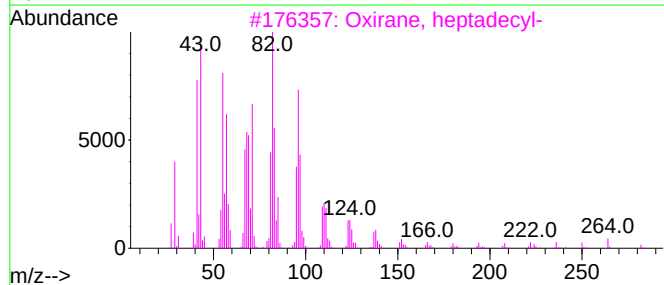
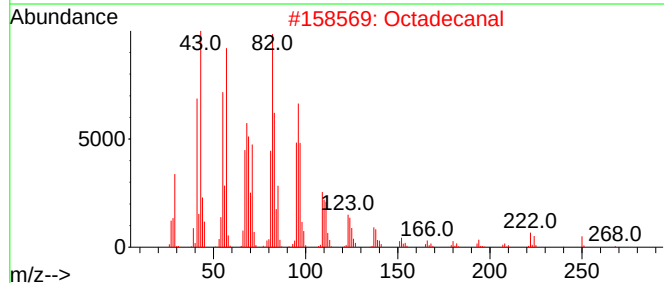
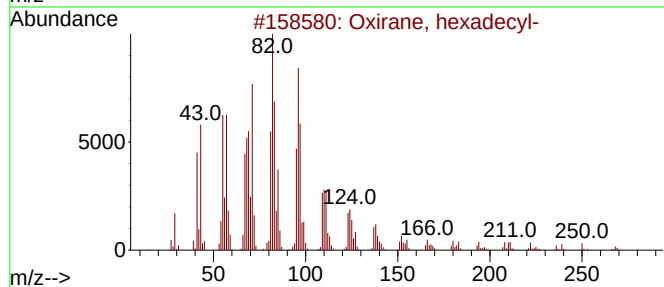
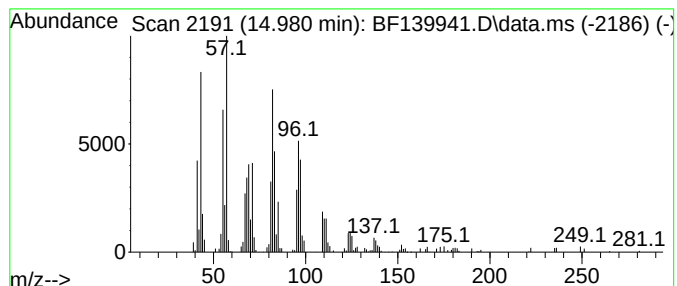
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Oxirane, hexadecyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.980	3.10 ng	139426	Perylene-d12	15.527	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2	Octadecanal	268	C18H36O	000638-66-4	91
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	90
4	1,19-Eicosadiene	278	C20H38	014811-95-1	74
5	Pentadecanal-	226	C15H30O	002765-11-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
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Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

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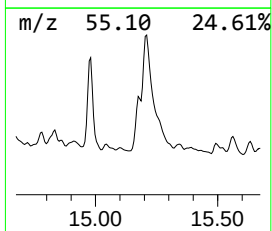
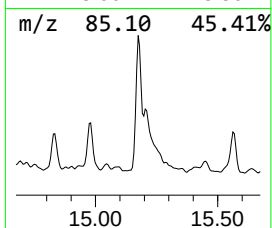
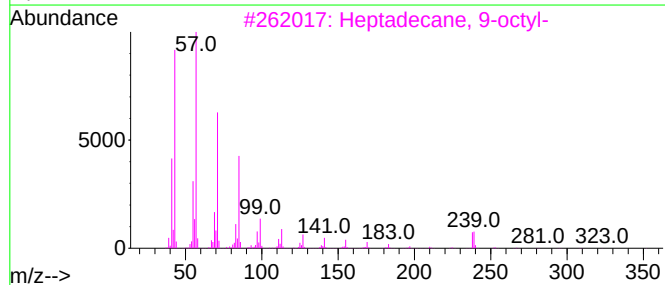
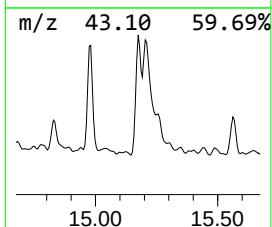
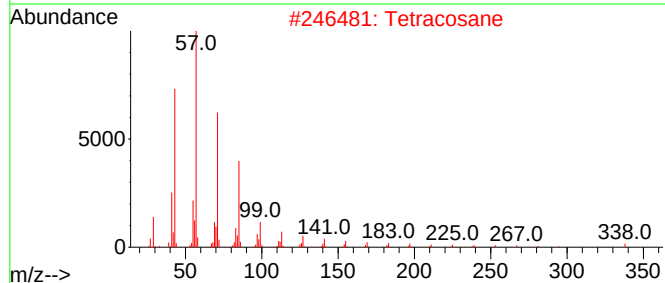
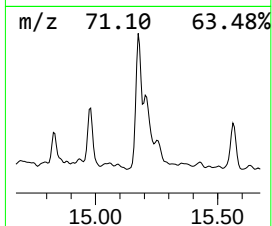
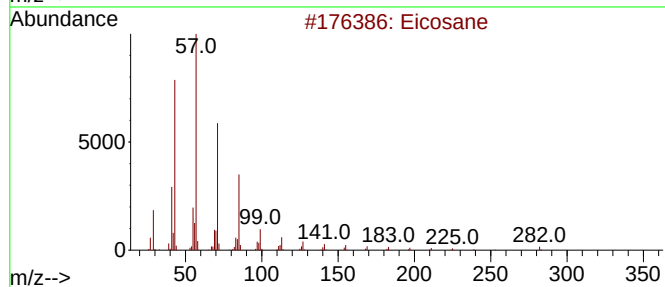
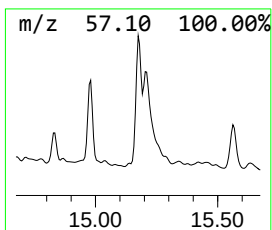
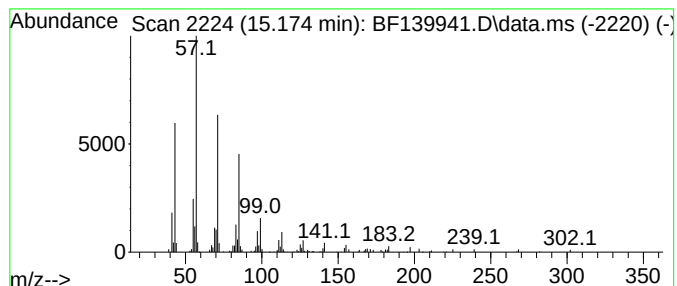
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.174	2.27 ng	102232	Perylene-d12	15.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	94
2			Tetracosane	338	C24H50	000646-31-1	91
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

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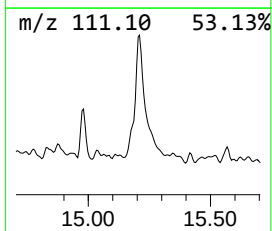
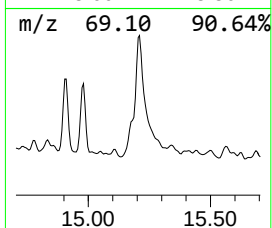
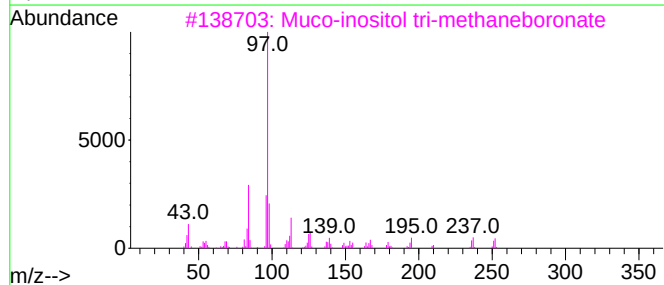
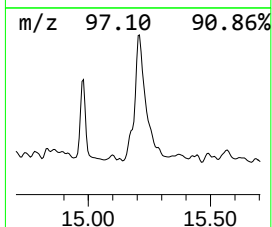
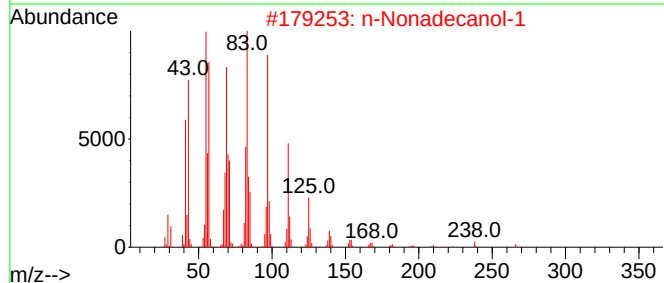
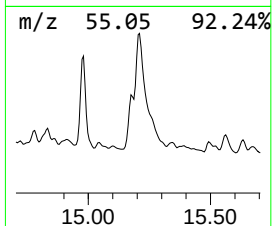
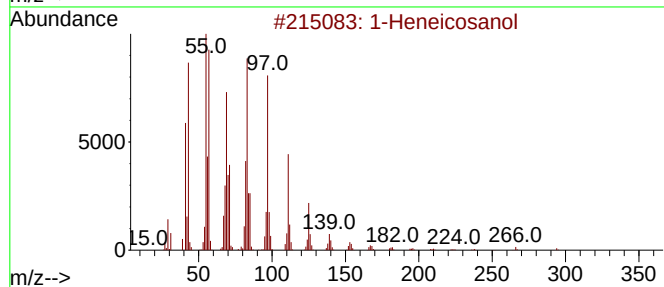
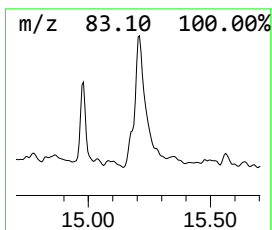
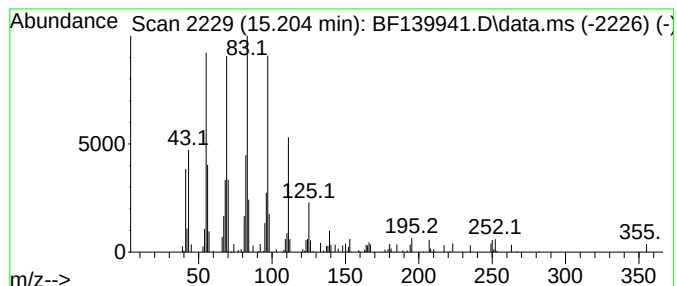
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.204	6.35 ng	285476	Perylene-d12	15.527

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	91
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	52
3		Muco-inositol tri-methaneboronate	252	C9H15B3O6	1000064-40-0	47
4		Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	46
5		Disparlure	282	C19H38O	029804-22-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

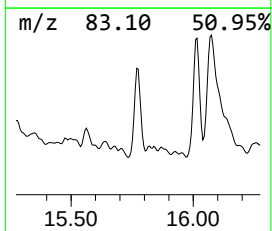
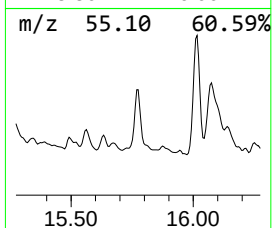
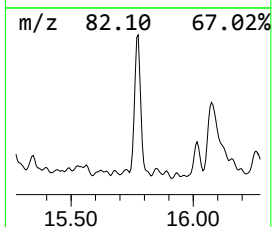
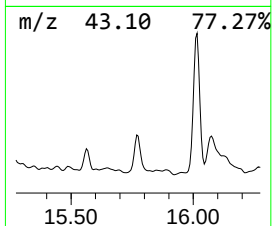
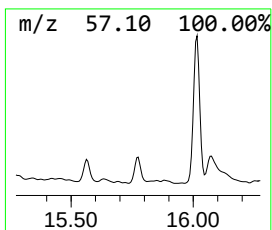
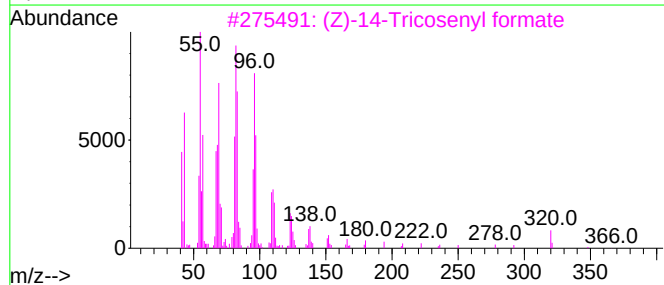
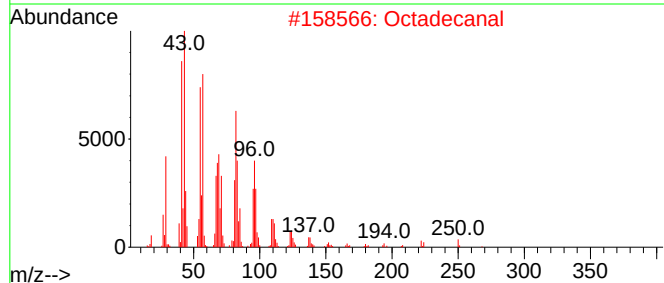
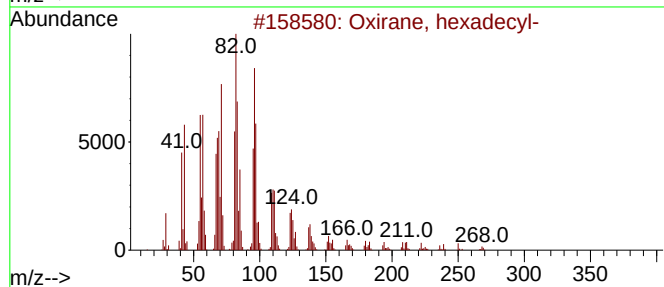
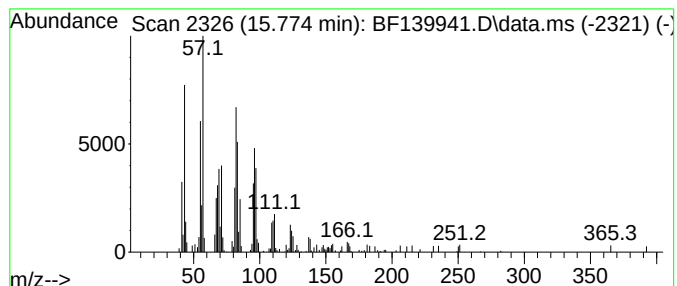
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.774	2.28 ng	102481	Perylene-d12	15.527	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
2	Octadecanal	268	C18H36O	000638-66-4	86
3	(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	70
4	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	68
5	1,19-Eicosadiene	278	C20H38	014811-95-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
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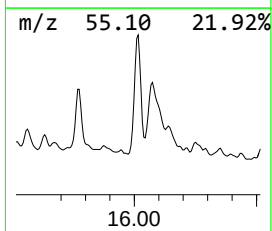
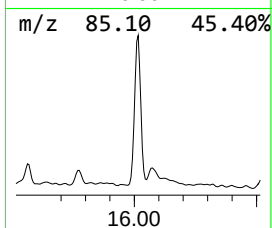
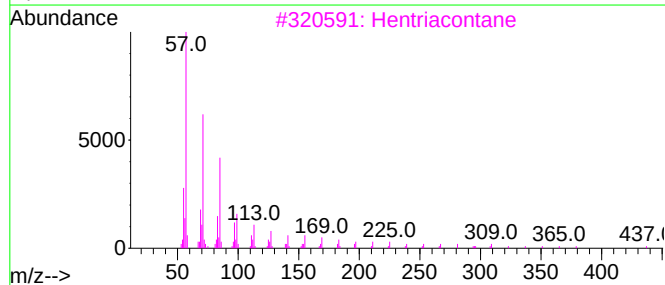
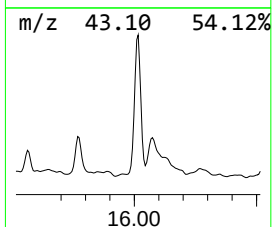
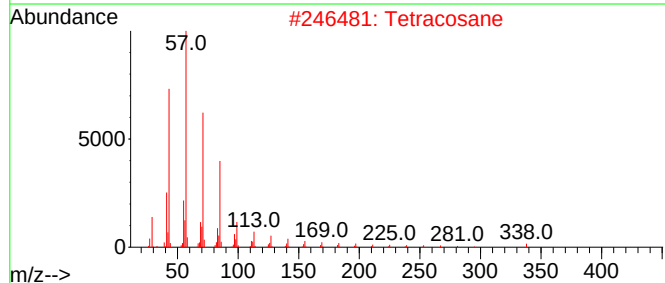
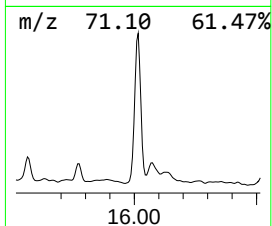
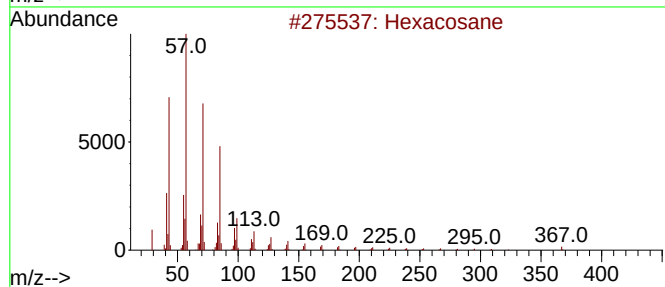
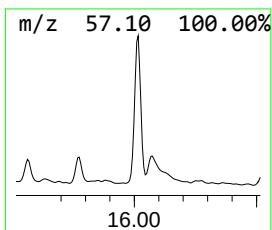
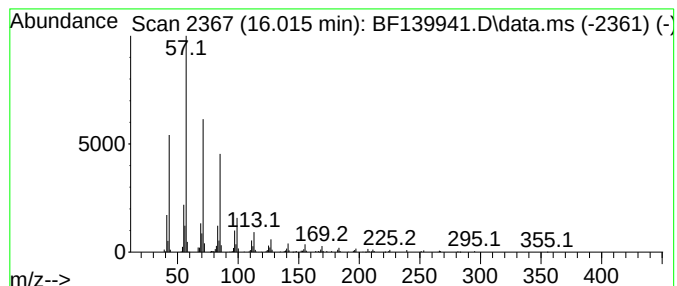
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.015	6.35 ng	285667	Perylene-d12	15.527	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	94
2	Tetracosane	338	C24H50	000646-31-1	91
3	Hentriacontane	437	C31H64	000630-04-6	91
4	Octacosane	394	C28H58	000630-02-4	91
5	Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

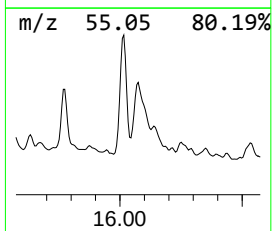
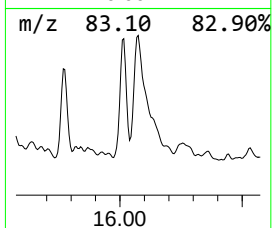
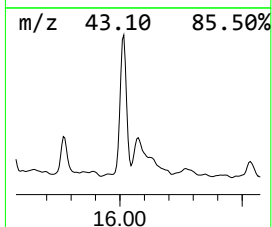
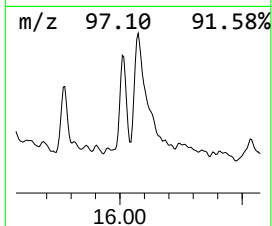
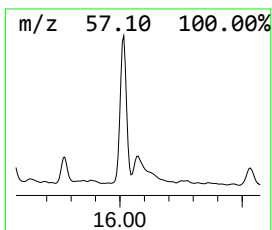
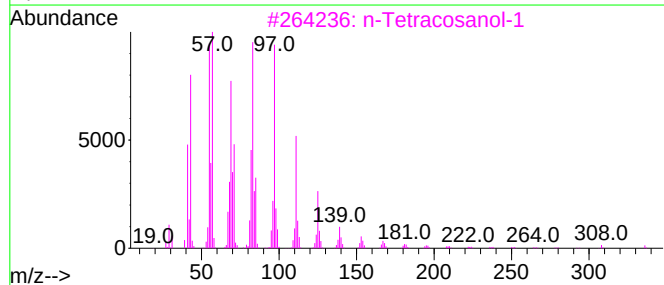
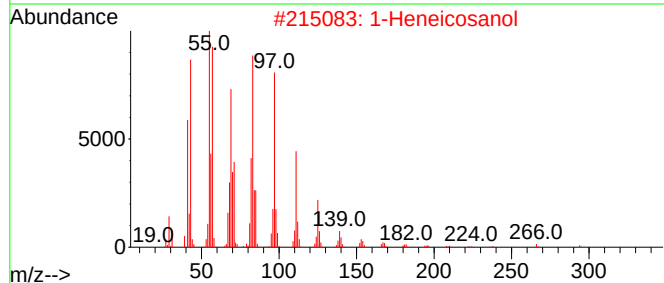
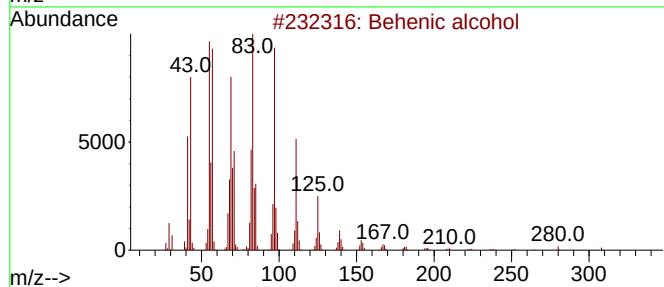
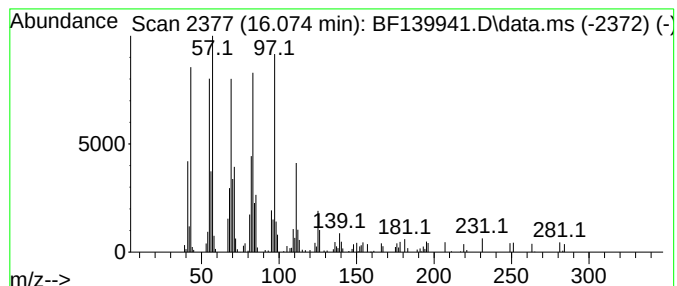
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Behenic alcohol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.074	3.07 ng	137826	Perylene-d12	15.527	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4	Cyclotetradecane	196	C14H28	000295-17-0	93
5	1-Hexacosene	364	C26H52	018835-33-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-1

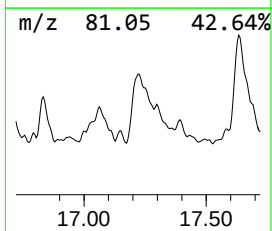
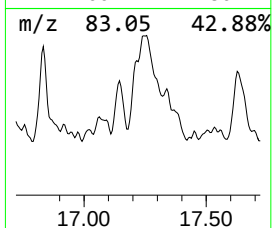
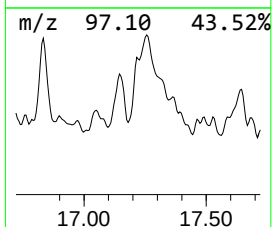
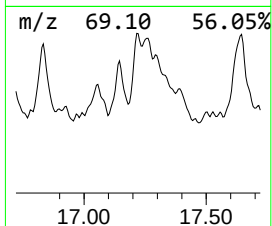
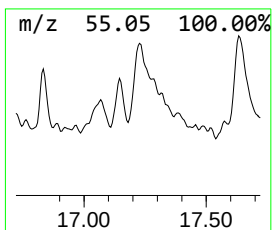
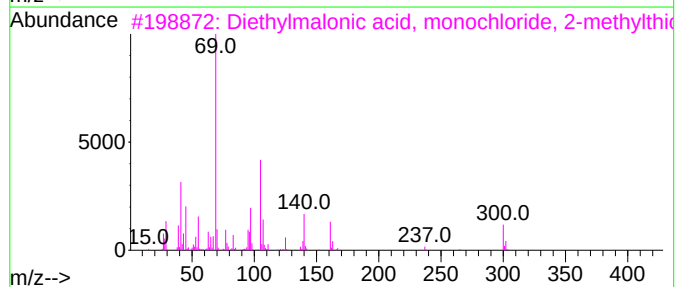
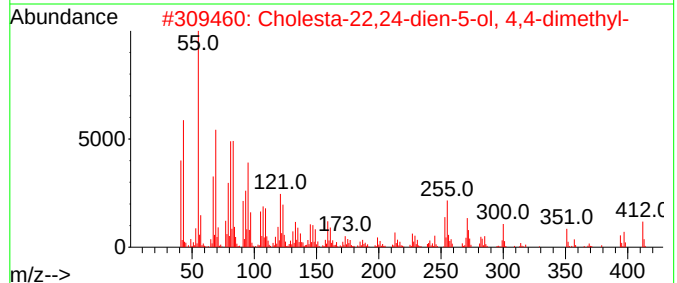
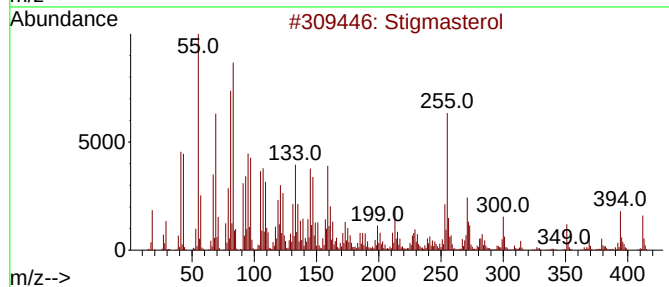
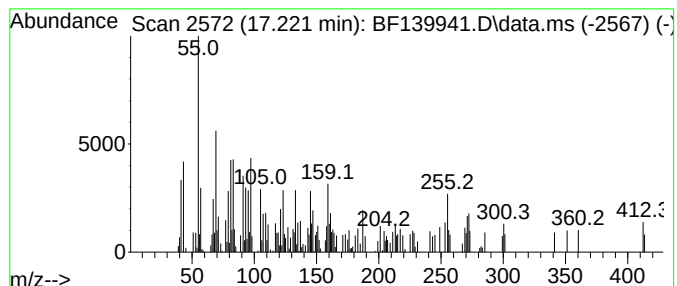
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Stigmasterol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.221	3.82 ng	171859	Perylene-d12	15.527

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Stigmasterol	412	C29H48O	000083-48-7	87
2		Cholesta-22,24-dien-5-ol, 4,4-di...	412	C29H48O	1000128-66-1	53
3		Diethylmalonic acid, monochlorid...	300	C14H17ClO3S	1000369-54-4	11
4		22-Stigmasten-3-one	412	C29H48O	004736-95-2	11
5		Chondrillasterol	412	C29H48O	000481-17-4	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

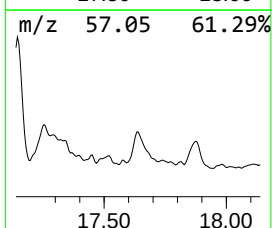
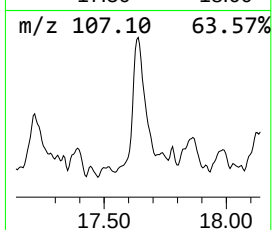
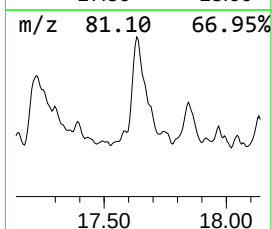
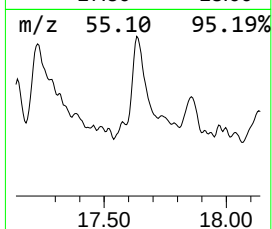
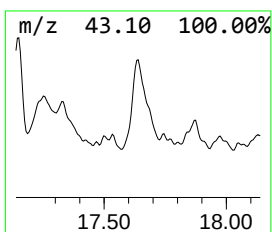
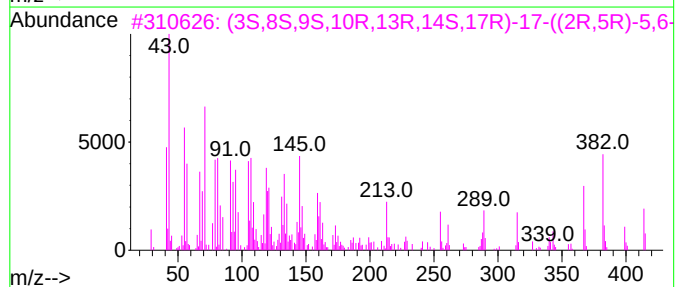
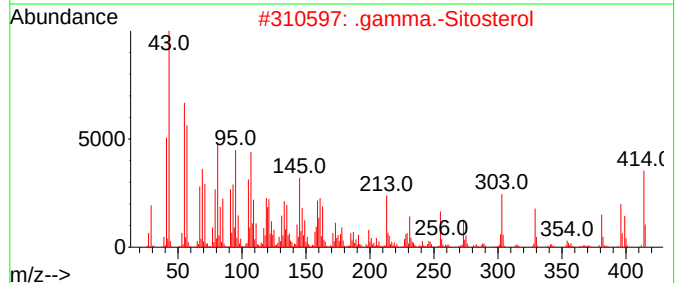
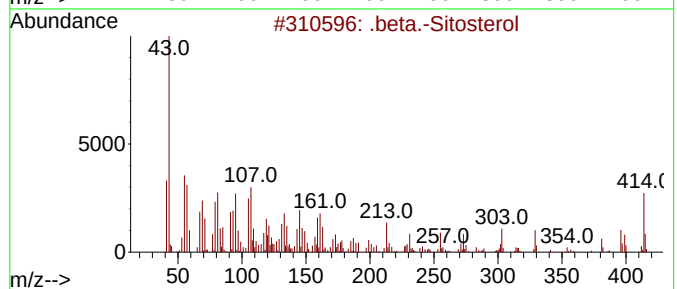
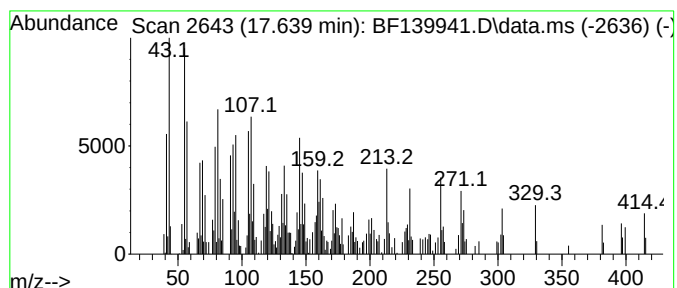
Instrument :
BNA_F
ClientSampleId :
SP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 .beta.-Sitosterol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.639	5.83 ng	262088	Perylene-d12	15.527	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Sitosterol	414	C29H50O	000083-46-5	99
2	.gamma.-Sitosterol	414	C29H50O	000083-47-6	87
3	(3S,8S,9S,10R,13R,14S,17R)-17-((...)	414	C29H50O	029365-29-5	83
4	Boscactol F	300	C20H28O2	1486443-17-7	74
5	Androst-5,16-diene-3.beta.-ol	272	C19H28O	001224-94-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102224\
Data File : BF139941.D
Acq On : 22 Oct 2024 21:20
Operator : RC/JU
Sample : P4385-02
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.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
Butane, 2-metho...	2.187	78.8	ng	3279150	1	6.887	832153	20.0
2-Pentanone, 4-...	5.110	5.4	ng	225704	1	6.887	832153	20.0
Benzophenone	10.633	16.1	ng	882492	3	9.922	1098310	20.0
Methanone, (1-h...	10.945	3.4	ng	152533	4	11.410	897512	20.0
n-Hexadecanoic ...	11.933	6.7	ng	301586	4	11.410	897512	20.0
5-Eicosene, (E)-	13.898	8.4	ng	362160	5	14.045	860025	20.0
Oxalic acid, is...	14.527	2.2	ng	93159	5	14.045	860025	20.0
Oxirane, hexade...	14.980	3.1	ng	139426	6	15.527	899329	20.0
Eicosane	15.174	2.3	ng	102232	6	15.527	899329	20.0
1-Heneicosanol	15.204	6.3	ng	285476	6	15.527	899329	20.0
Octadecanal	15.774	2.3	ng	102481	6	15.527	899329	20.0
Hexacosane	16.015	6.3	ng	285667	6	15.527	899329	20.0
Behenic alcohol	16.074	3.1	ng	137826	6	15.527	899329	20.0
Stigmasterol	17.221	3.8	ng	171859	6	15.527	899329	20.0
.beta.-Sitosterol	17.639	5.8	ng	262088	6	15.527	899329	20.0