

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
 Data File : BF126068.D
 Acq On : 8 Nov 2021 16:42
 Operator : JU/CG
 Sample : M4540-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-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-BF110321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126068.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.163	5	13	24	rVB	1853989	2359496	46.78%	6.493%
2	2.975	145	151	156	rBV2	37192	58892	1.17%	0.162%
3	5.075	499	508	512	rBV	89259	108954	2.16%	0.300%
4	5.469	570	575	579	rBV	3705292	3771608	74.78%	10.379%
5	6.481	741	747	750	rBV	3983401	3827229	75.88%	10.532%
6	6.851	807	810	814	rBV	1183089	939154	18.62%	2.584%
7	7.416	900	906	909	rBV	2670858	2441085	48.40%	6.717%
8	8.033	1008	1011	1022	rBV	316891	378448	7.50%	1.041%
9	8.133	1024	1028	1033	rVV	1376493	1106034	21.93%	3.044%
10	9.210	1206	1211	1214	rBV	5150416	4099392	81.28%	11.281%
11	9.327	1228	1231	1235	rBV2	63457	56045	1.11%	0.154%
12	9.516	1260	1263	1266	rBV2	137728	124350	2.47%	0.342%
13	9.551	1266	1269	1272	rVV	132849	106189	2.11%	0.292%
14	9.592	1272	1276	1278	rVB	69717	62628	1.24%	0.172%
15	9.804	1309	1312	1314	rBV	140661	120747	2.39%	0.332%
16	9.863	1318	1322	1324	rBV	1189278	1036813	20.56%	2.853%
17	9.886	1324	1326	1333	rVB	1313248	1148998	22.78%	3.162%
18	10.010	1344	1347	1350	rBV	253750	201916	4.00%	0.556%
19	10.051	1350	1354	1358	rVB2	122396	109611	2.17%	0.302%
20	10.674	1455	1460	1463	rBV	2293174	2052674	40.70%	5.649%
21	10.710	1464	1466	1470	rVB3	56248	55043	1.09%	0.151%
22	10.868	1490	1493	1504	rVB3	60003	85902	1.70%	0.236%
23	11.368	1574	1578	1585	rBV	1075514	937158	18.58%	2.579%
24	12.957	1844	1848	1851	rBV	3974987	3141307	62.28%	8.644%
25	13.162	1880	1883	1886	rBV	112124	102405	2.03%	0.282%
26	13.498	1934	1940	1944	rBV2	109606	135704	2.69%	0.373%
27	13.609	1955	1959	1963	rBV	84905	77203	1.53%	0.212%
28	13.727	1974	1979	1981	rBV	5292224	5043570	100.00%	13.879%
29	13.862	1998	2002	2015	rBV2	208851	348484	6.91%	0.959%
30	14.009	2022	2027	2030	rBV	1051057	917326	18.19%	2.524%
31	14.504	2104	2111	2115	rBV5	76707	161940	3.21%	0.446%
32	14.933	2182	2184	2190	rVB3	59969	67951	1.35%	0.187%
33	15.133	2214	2218	2221	rBV	65848	75950	1.51%	0.209%
34	15.168	2221	2224	2229	rBV6	36347	62929	1.25%	0.173%
35	15.468	2271	2275	2280	rBV	567608	694704	13.77%	1.912%
36	15.509	2280	2282	2290	rVB2	40209	56952	1.13%	0.157%

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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-BF110321.M
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37	15.715	2314	2317	2324	rVB5	51578	71243	1.41%	0.196%
38	15.951	2353	2357	2362	rVB	85628	135472	2.69%	0.373%
39	16.456	2439	2443	2448	rBV3	36142	58381	1.16%	0.161%

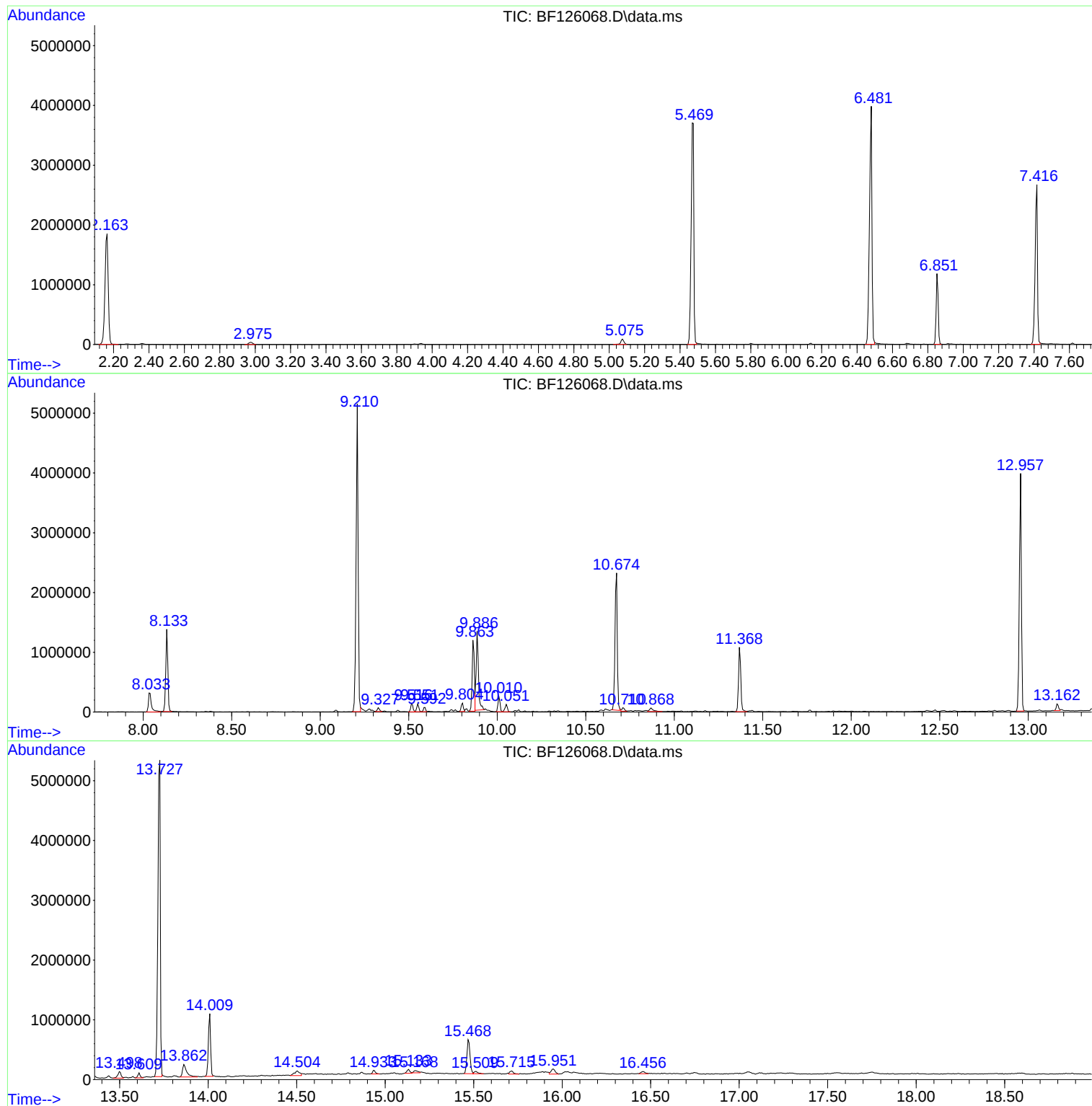
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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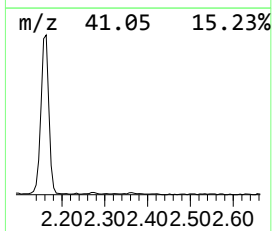
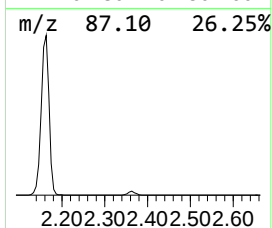
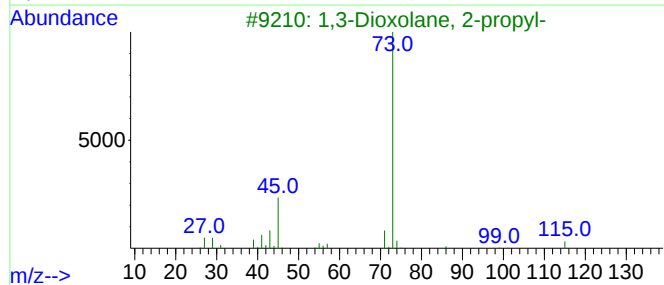
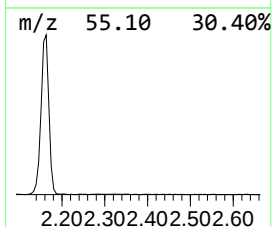
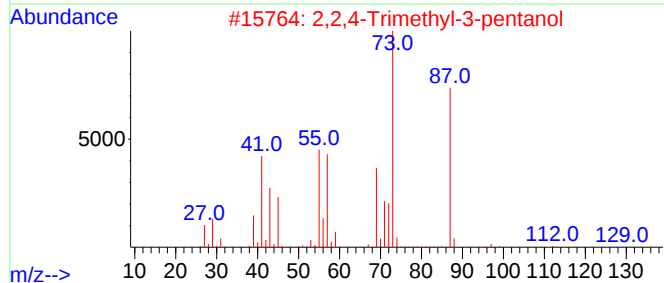
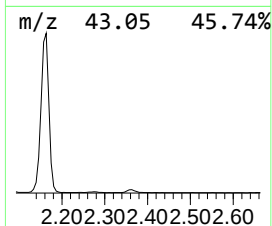
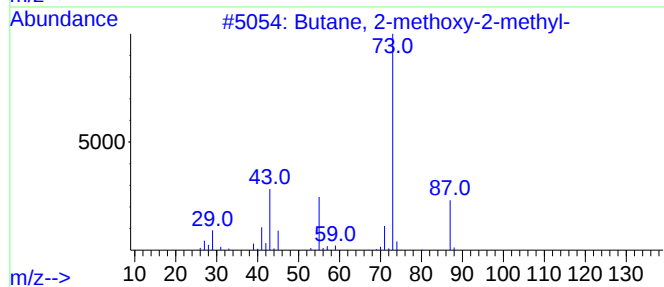
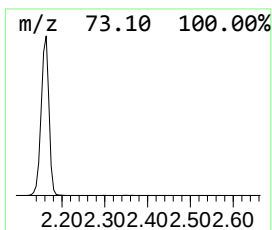
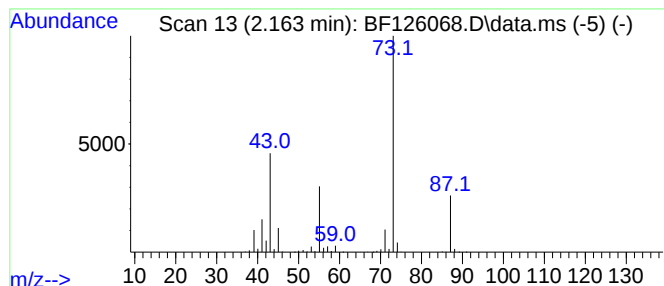
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.163	50.25 ng	2359500	1,4-Dichlorobenzene-d4	6.851

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-propyl-	116	C6H12O2	003390-13-4	23
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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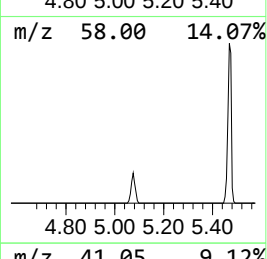
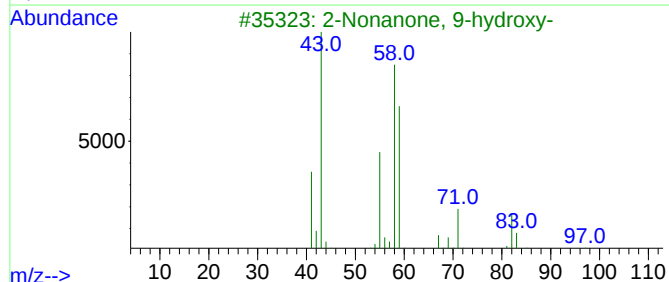
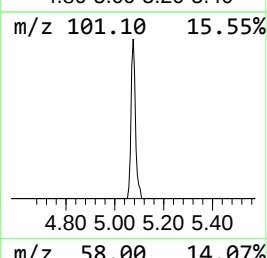
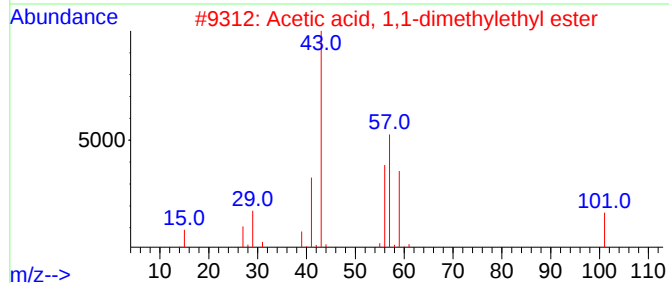
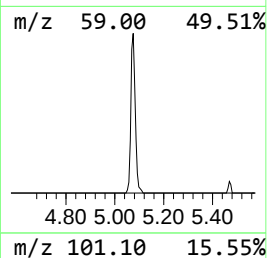
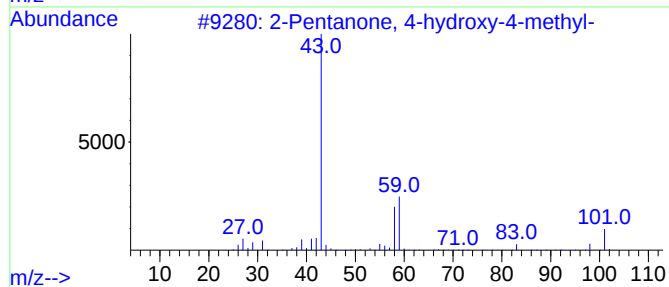
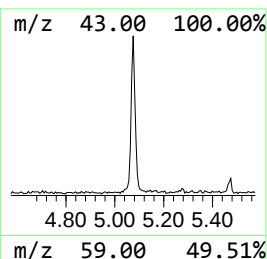
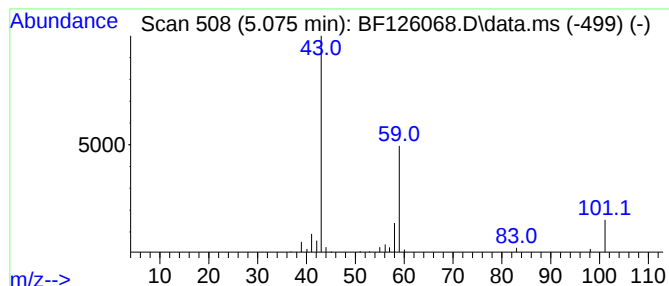
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.075	2.32 ng	108954	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	36
4			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9
5			Diacetamide	101	C4H7NO2	000625-77-4	9



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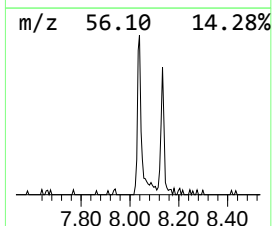
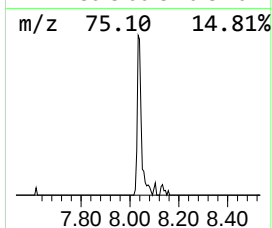
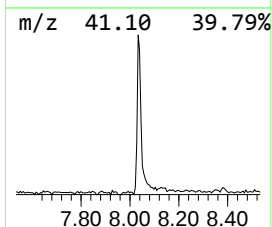
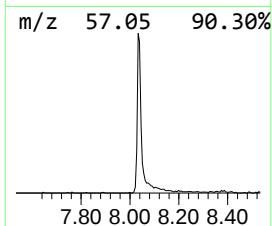
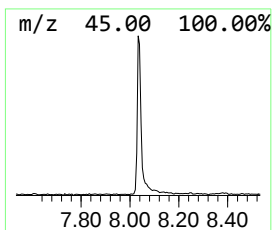
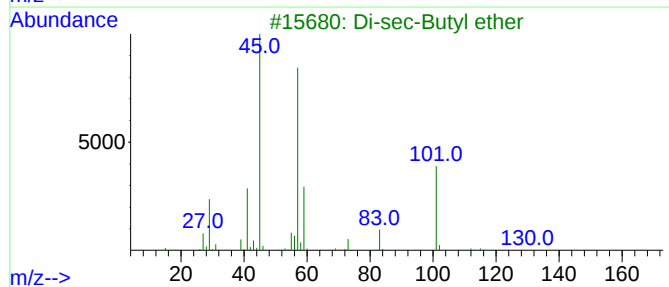
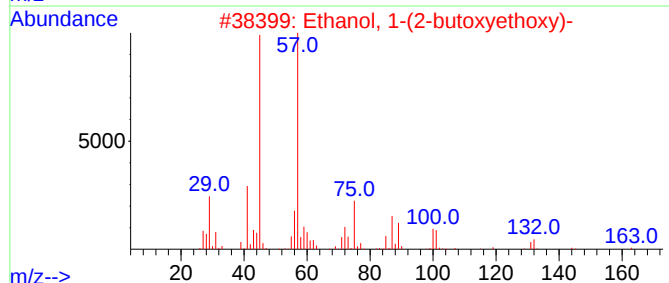
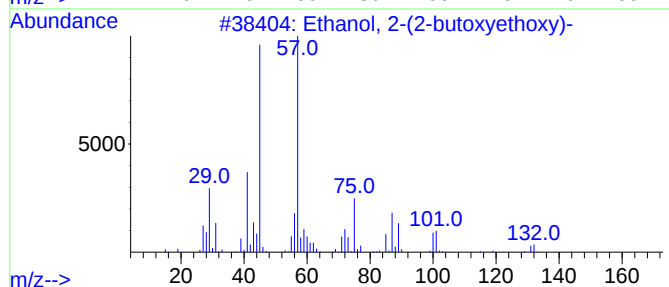
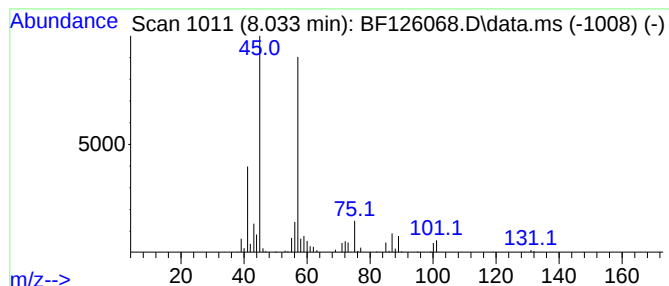
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.033	6.84 ng	378448	Naphthalene-d8	8.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Di-sec-Butyl ether	130	C8H18O	006863-58-7	59
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50
5			Methoxyacetic acid, heptyl ester	188	C10H20O3	168920-36-3	50



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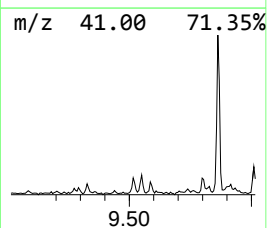
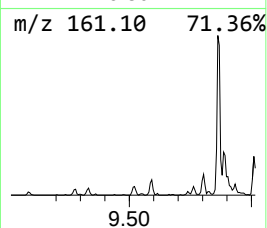
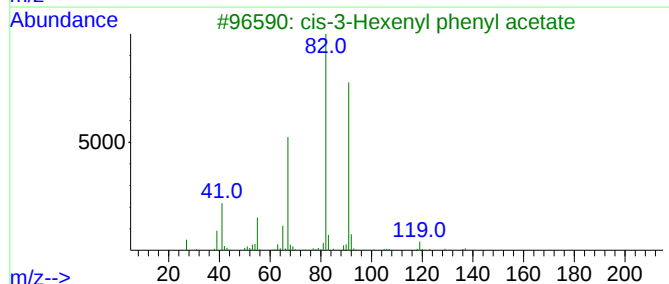
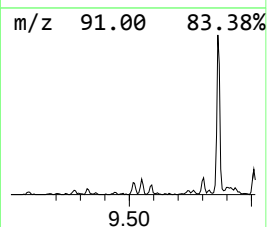
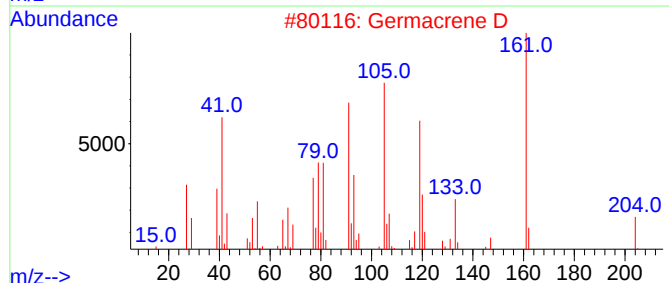
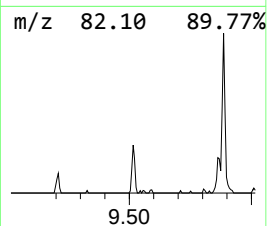
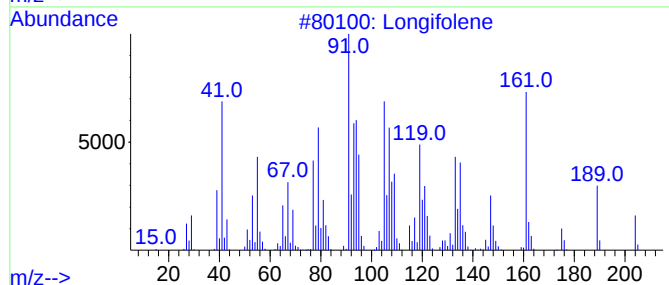
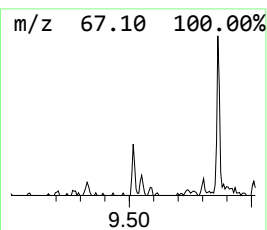
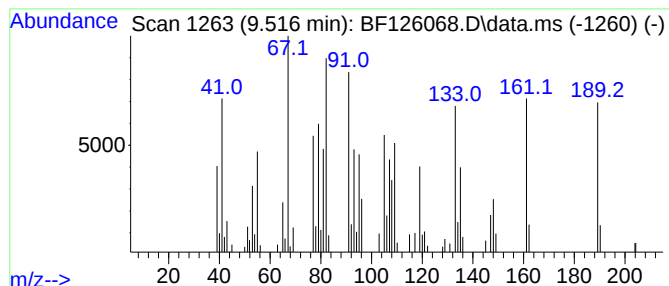
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Longifolene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.516	2.16 ng	124350	Acenaphthene-d10	9.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Longifolene	204	C15H24	000475-20-7	86
2			Germacrene D	204	C15H24	023986-74-5	25
3			cis-3-Hexenyl phenyl acetate	218	C14H18O2	042436-07-7	25
4			Cis-bicyclo[4.2.0]octane	110	C8H14	028282-35-1	22
5			3-Hexyne	82	C6H10	000928-49-4	20



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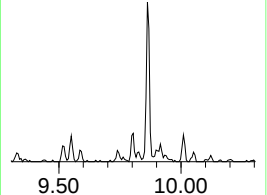
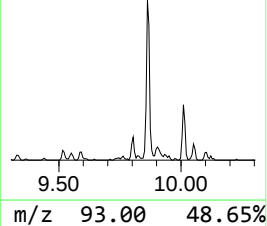
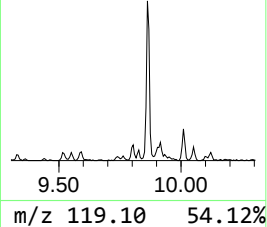
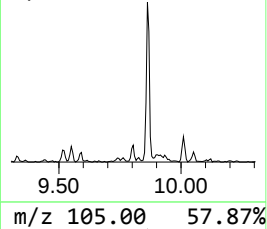
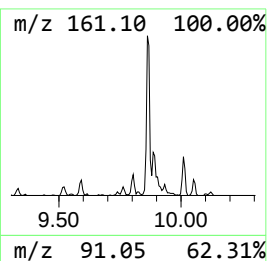
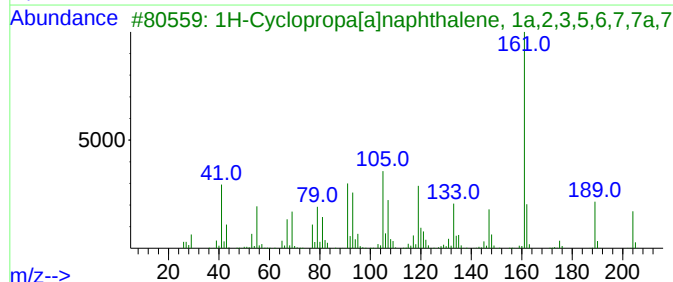
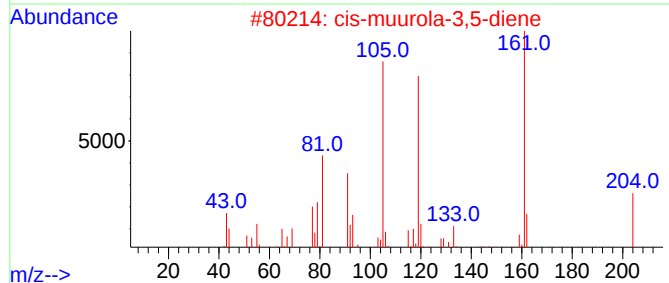
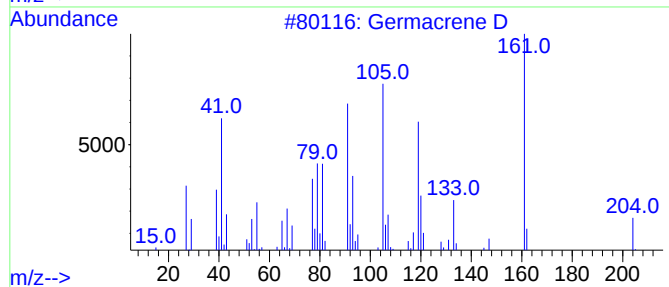
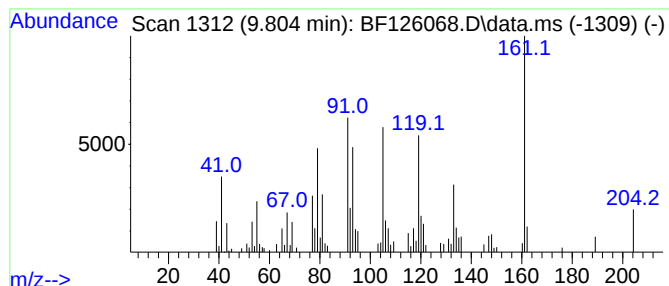
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Peak Number 5 Germacrene D Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.804	2.10 ng	120747	Acenaphthene-d10	9.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Germacrene D	204	C15H24	023986-74-5	95
2			cis-muurolo-3,5-diene	204	C15H24	1000365-95-4	87
3			1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	017334-55-3	87
4			1H-Cyclopenta[1,3]cyclopropa[1,2...	204	C15H24	013744-15-5	70
5			Copaene	204	C15H24	003856-25-5	64



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Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
Operator : JU/CG
Sample : M4540-01
Misc :
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Instrument :
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ClientSampleId :
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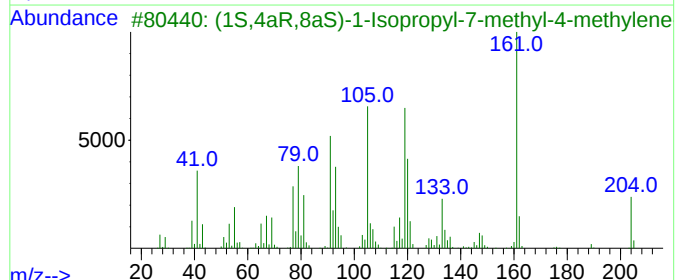
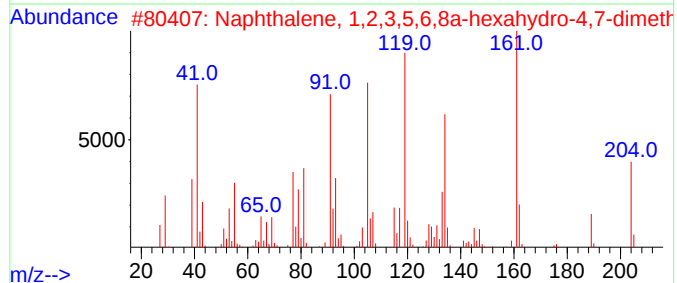
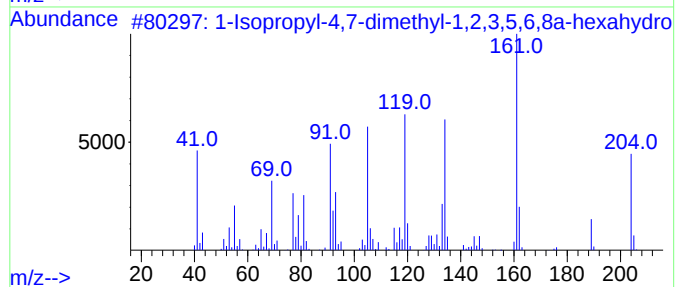
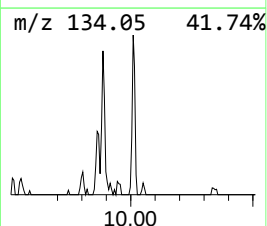
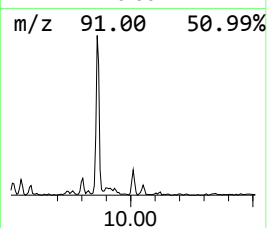
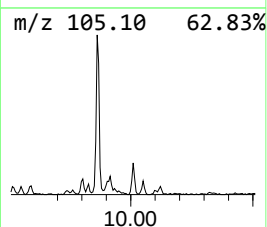
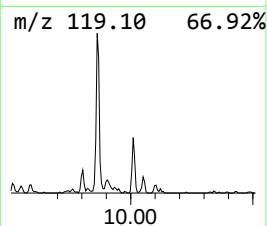
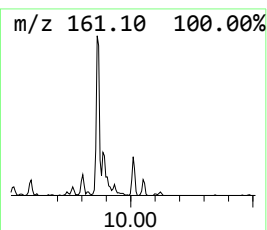
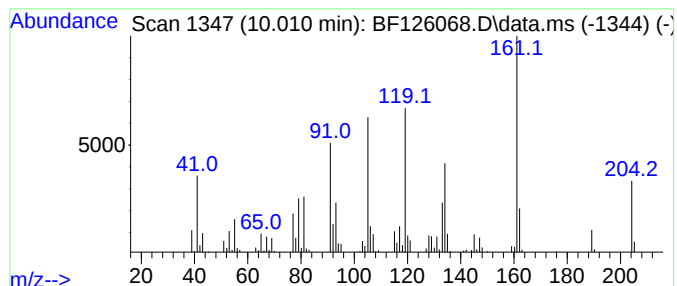
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1-Isopropyl-4,7-dimethyl-1,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.010	3.51 ng	201916	Acenaphthene-d10	9.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Isopropyl-4,7-dimethyl-1,2,3,5...	204	C15H24	016729-01-4	98
2			Naphthalene, 1,2,3,5,6,8a-hexahy...	204	C15H24	000483-76-1	97
3			(1S,4aR,8aS)-1-Isopropyl-7-methy...	204	C15H24	006980-46-7	86
4			1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	83
5			cis-muurola-3,5-diene	204	C15H24	1000365-95-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
Operator : JU/CG
Sample : M4540-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

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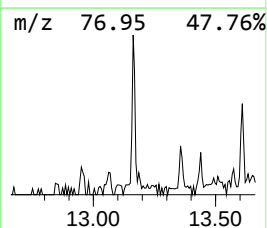
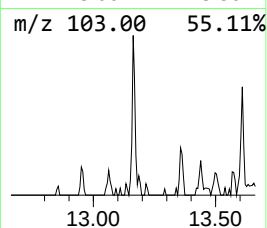
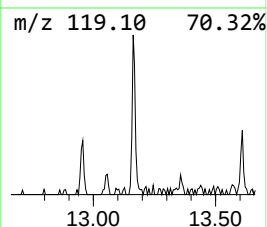
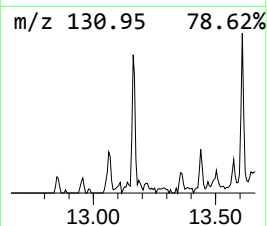
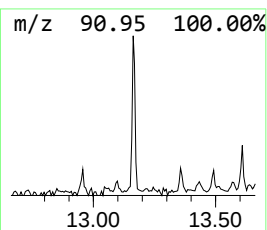
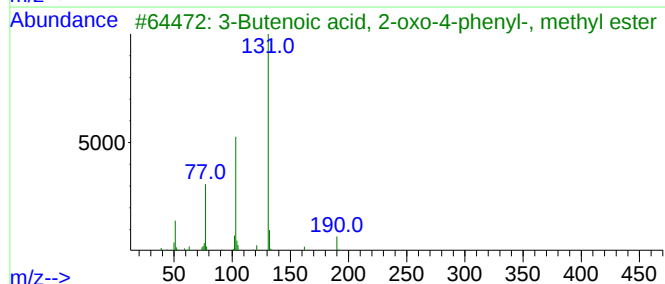
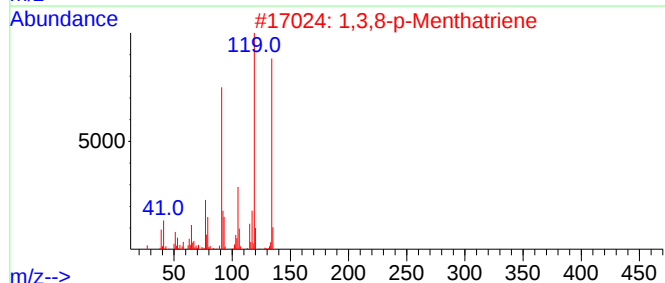
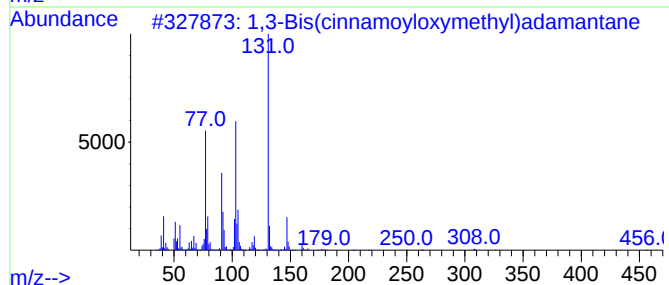
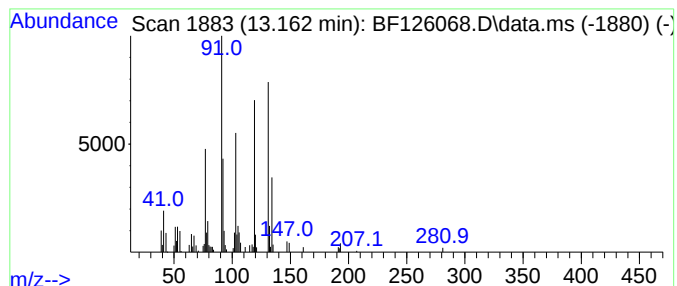
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,3-Bis(cinnamoyloxymethyl)... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.163	2.23 ng	102405	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Bis(cinnamoyloxymethyl)adama...	456	C30H32O4	303797-58-2	53
2			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	43
3			3-Butenoic acid, 2-oxo-4-phenyl-...	190	C11H10O3	1000142-22-4	38
4			7-Oxo-1,3,5-cycloheptatriene-1-c...	131	C8H5NO	000934-19-0	38
5			Myrtenyl isovalerate	236	C15H24O2	033900-84-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
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Sample : M4540-01
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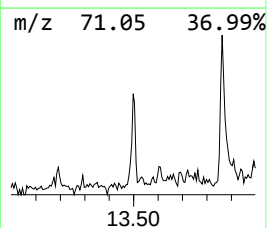
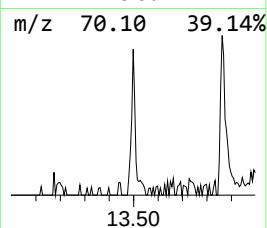
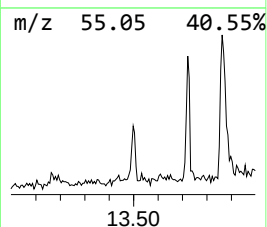
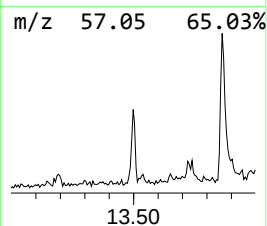
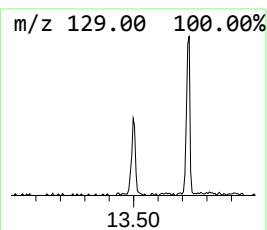
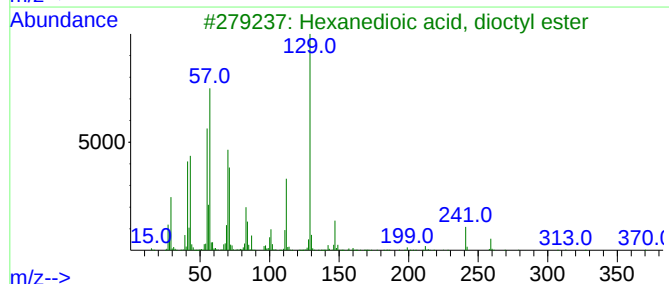
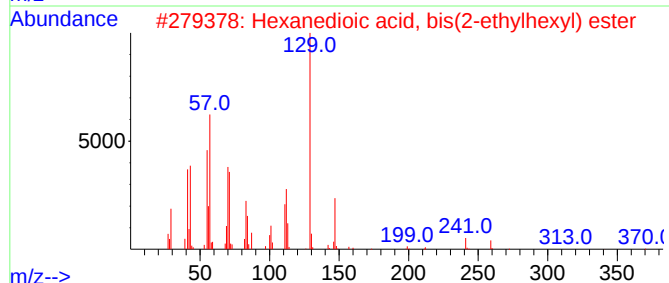
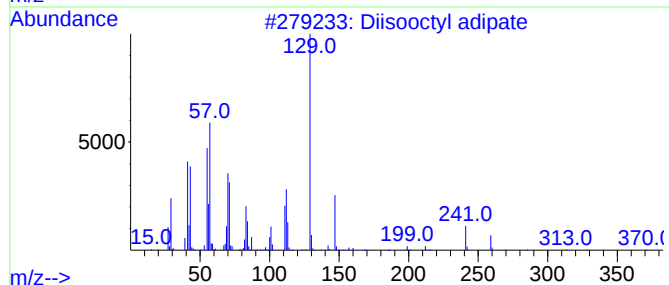
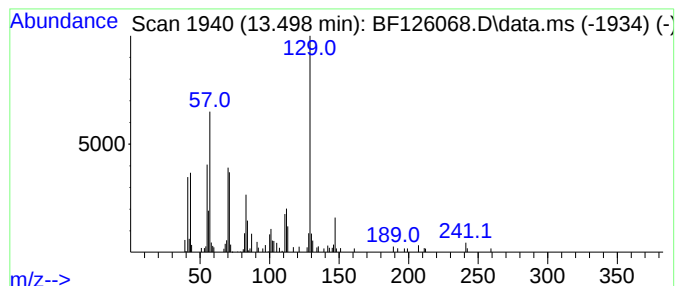
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Diisooctyl adipate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.498	2.96 ng	135704	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisooctyl adipate	370	C22H42O4	001330-86-5	90
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	83
3			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	72
4			Adipic acid, 2-ethylhexyl isobut...	314	C18H34O4	1010324-48-0	64
5			Adipic acid, octyl 2-octyl ester	370	C22H42O4	1000324-45-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
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Sample : M4540-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

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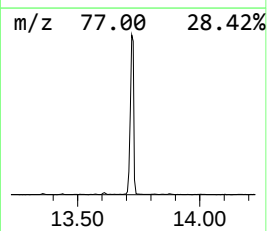
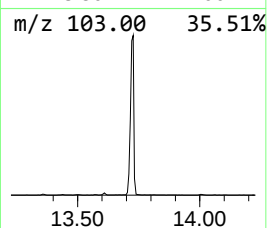
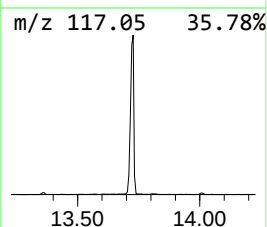
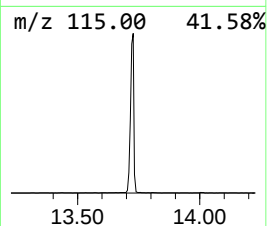
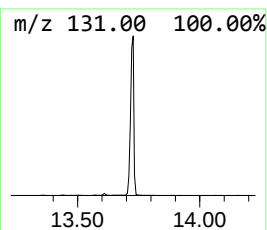
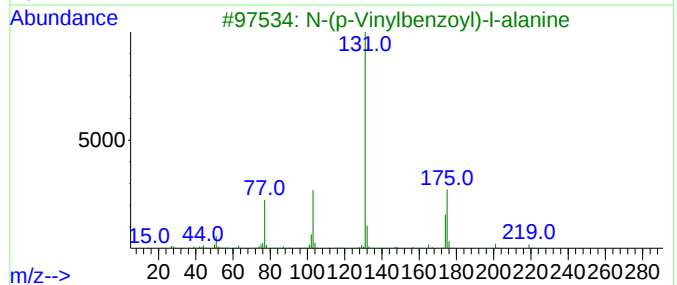
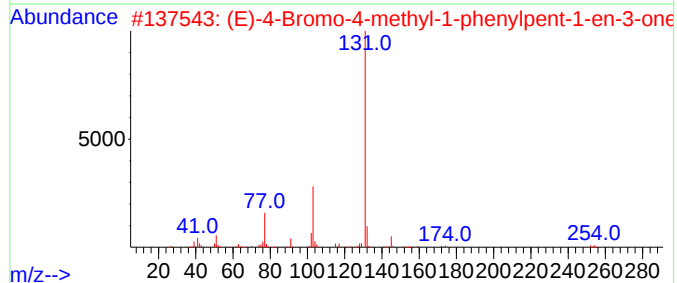
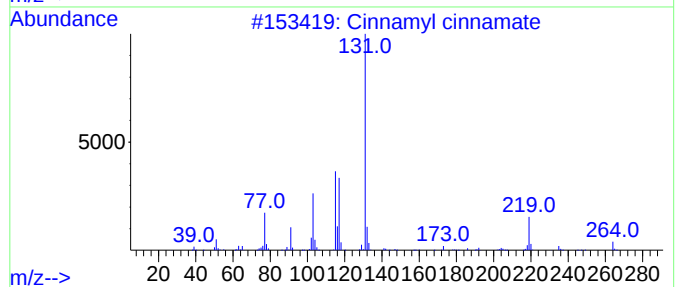
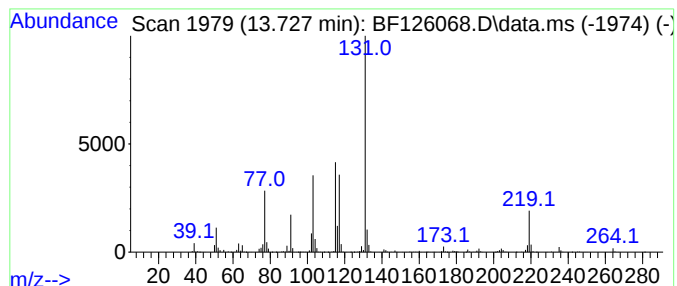
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cinnamyl cinnamate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.727	109.96 ng	5043570	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2			(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	49
3			N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	43
4			(2E)-N-(2-Fluorophenyl)-3-phenyl...	241	C15H12FNO	025893-50-9	43
5			Propanedioic acid, nitrile, hydr...	229	C12H11N3O2	294878-35-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
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ClientSampleId :
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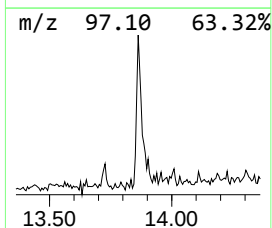
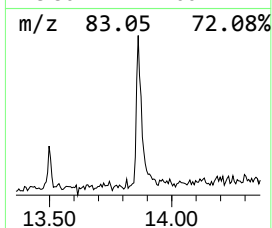
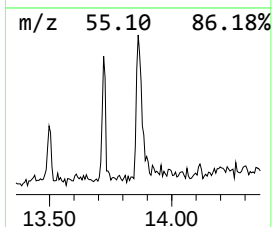
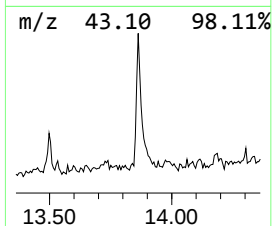
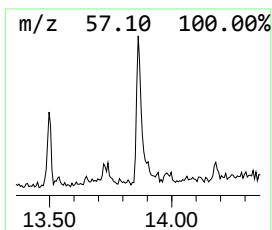
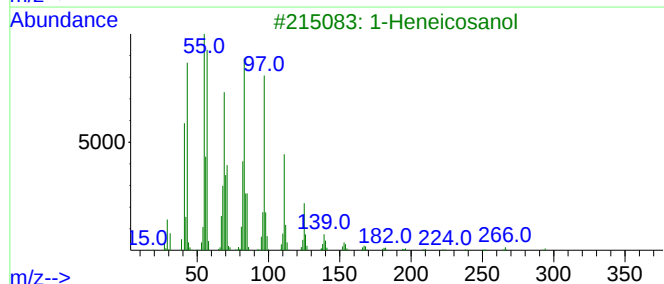
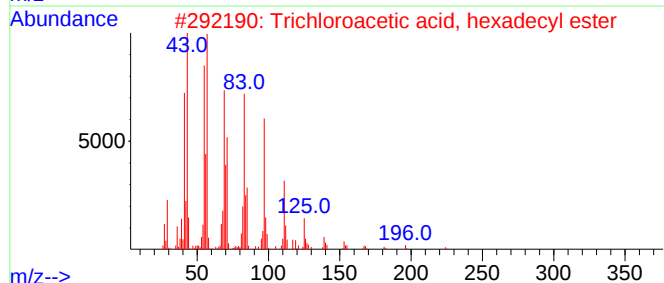
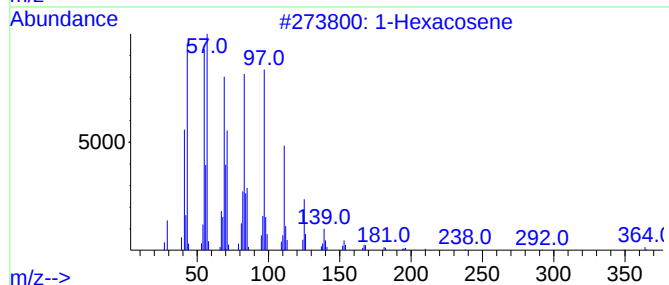
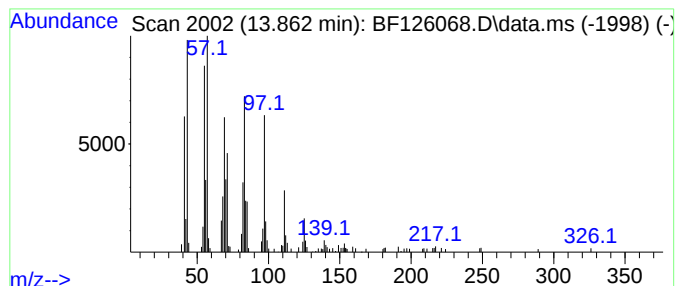
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1-Hexacosene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.862	7.60 ng	348484	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexacosene	364	C26H52	018835-33-1	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
3			1-Heneicosanol	312	C21H44O	015594-90-8	93
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
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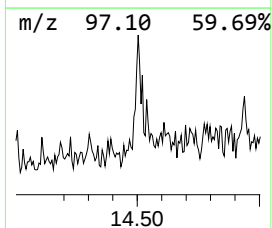
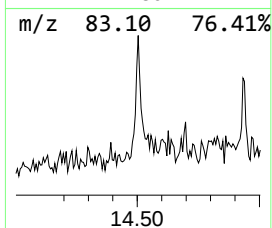
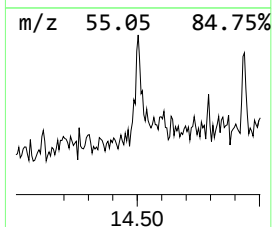
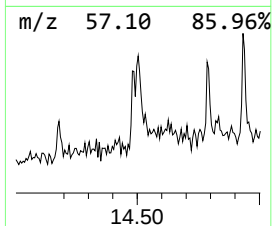
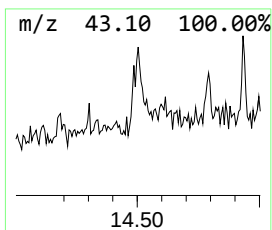
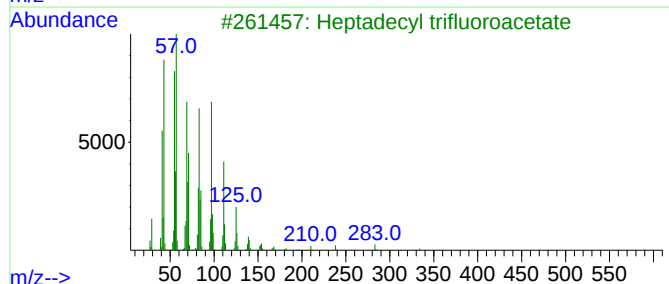
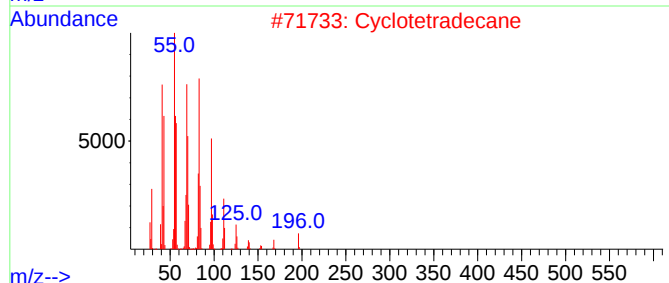
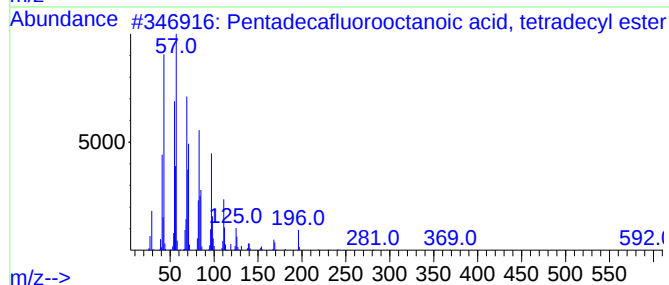
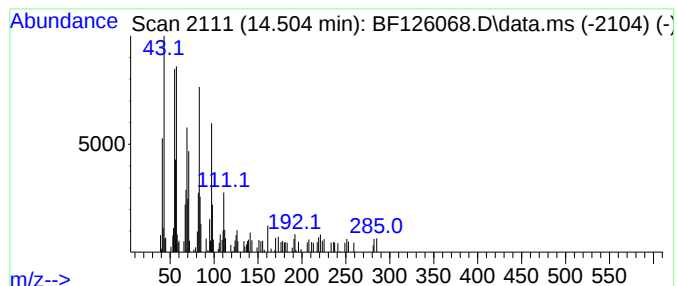
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentadecafluorooctanoic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.504	3.53 ng	161940	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, te...	610	C22H29F15O2	1000406-04-4	90
2			Cyclotetradecane	196	C14H28	000295-17-0	86
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	72
4			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	64
5			2- Bromopropionic acid, pentadec...	362	C18H35BrO2	1000292-44-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
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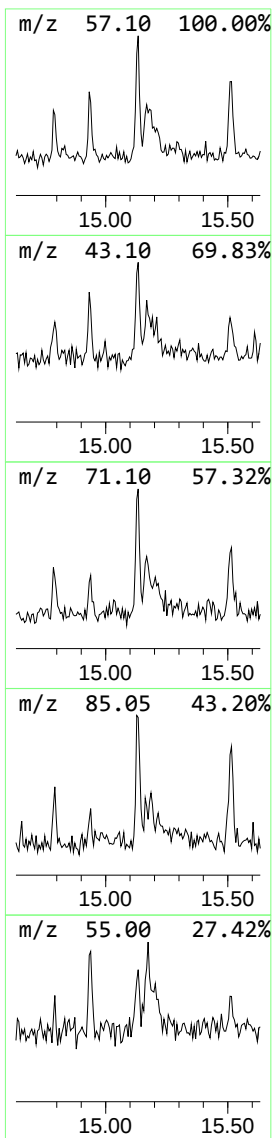
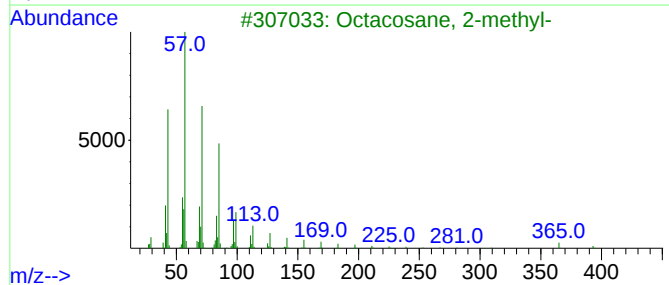
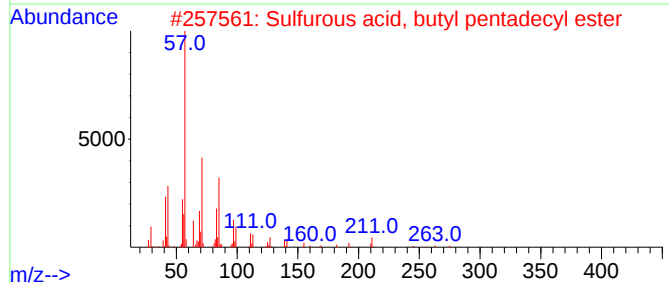
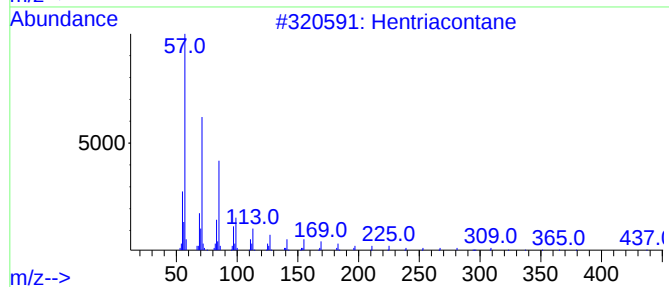
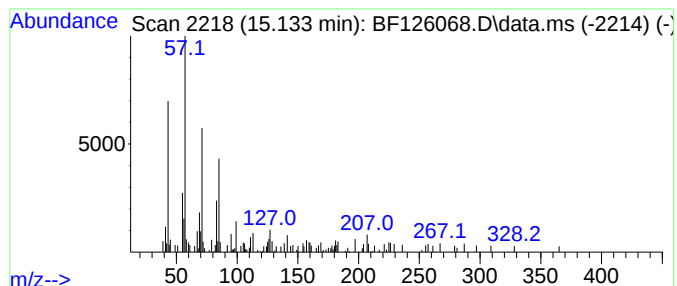
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hentriacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.133	2.19 ng	75950	Perylene-d12	15.468	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	81
2	Sulfurous acid, butyl pentadecyl...	348	C19H40O3S	1000309-18-2	80
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	80
4	Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	80
5	Heptacosane	380	C27H56	000593-49-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
Operator : JU/CG
Sample : M4540-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-1

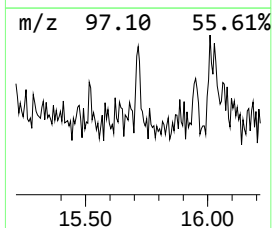
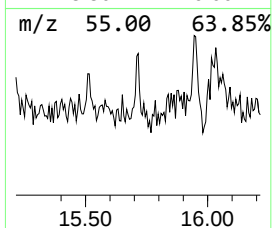
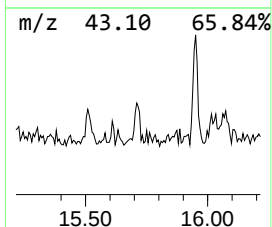
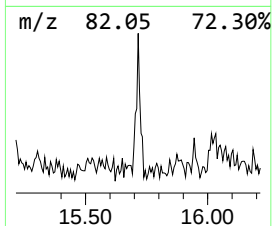
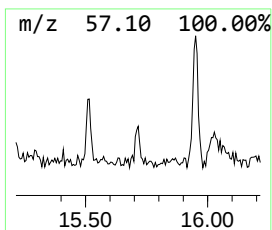
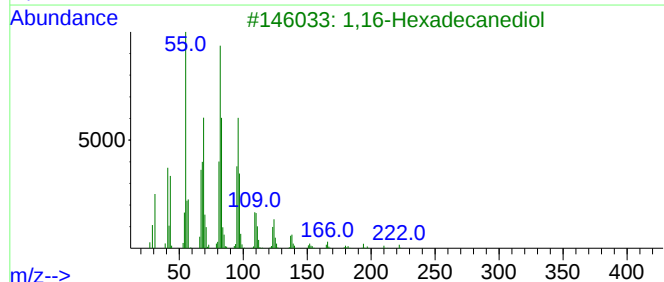
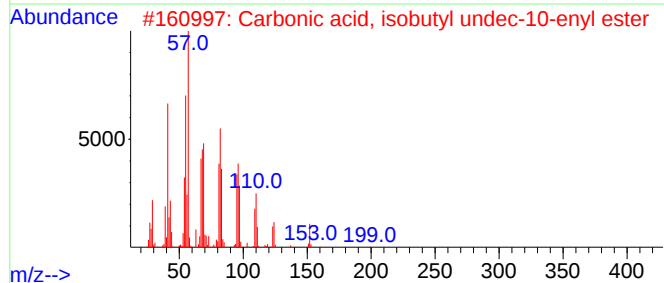
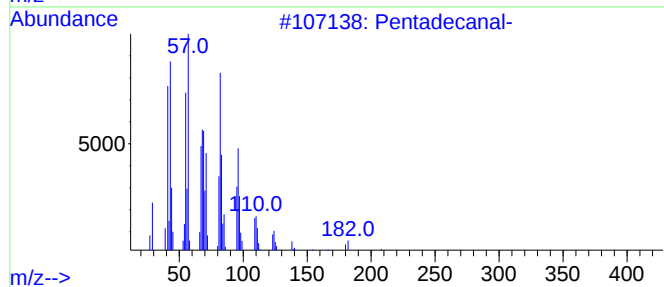
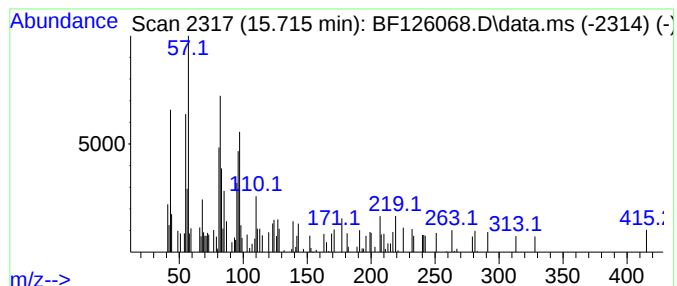
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.715 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.715	2.05 ng	71243	Perylene-d12	15.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanal-	226	C15H30O	002765-11-9	38
2			Carbonic acid, isobutyl undec-10...	270	C16H30O3	1000314-60-8	38
3			1,16-Hexadecanediol	258	C16H34O2	007735-42-4	38
4			N-(1-Methylpiperidin-4-yl)-4-nit...	263	C13H17N3O3	210643-97-3	38
5			Tetradecanal	212	C14H28O	000124-25-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
Operator : JU/CG
Sample : M4540-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-1

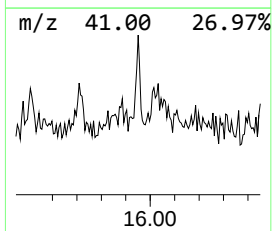
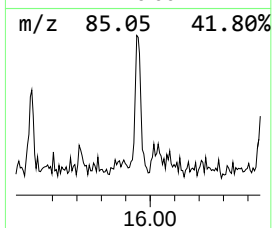
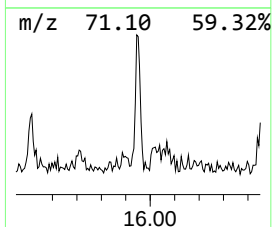
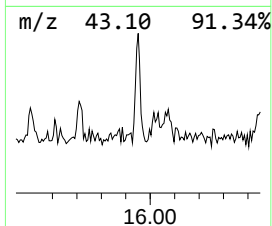
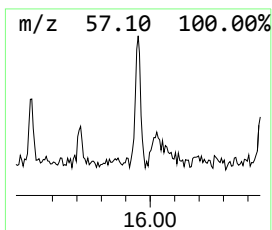
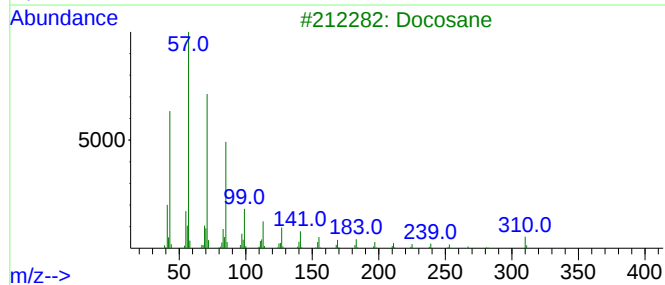
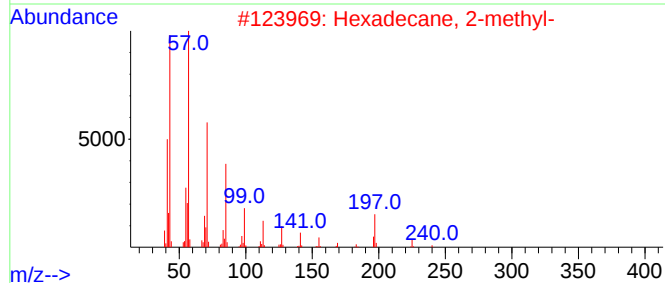
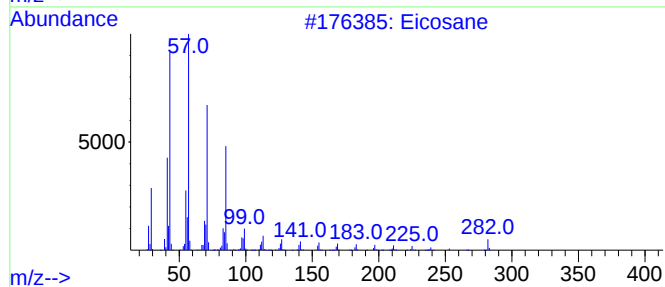
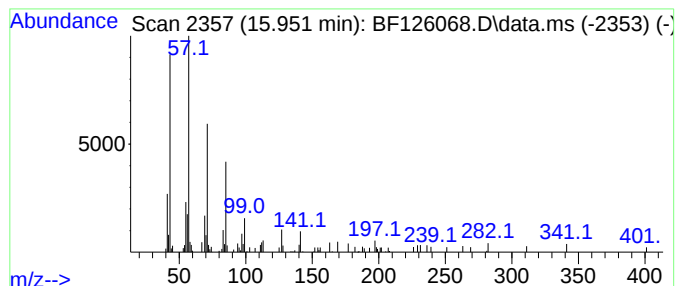
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.951	3.90 ng	135472	Perylene-d12	15.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	90
2			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	86
3			Docosane	310	C22H46	000629-97-0	86
4			Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	86
5			Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110821\
Data File : BF126068.D
Acq On : 8 Nov 2021 16:42
Operator : JU/CG
Sample : M4540-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.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...	2.163	50.3	ng	2359500	1	6.851	939154	20.0
2-Pentanone, 4-...	5.075	2.3	ng	108954	1	6.851	939154	20.0
Ethanol, 2-(2-b...	8.033	6.8	ng	378448	2	8.133	1106030	20.0
Longifolene	9.516	2.2	ng	124350	3	9.886	1149000	20.0
Germacrene D	9.804	2.1	ng	120747	3	9.886	1149000	20.0
1-Isopropyl-4,7...	10.010	3.5	ng	201916	3	9.886	1149000	20.0
1,3-Bis(cinnamo...	13.163	2.2	ng	102405	5	14.009	917326	20.0
Diisooctyl adipate	13.498	3.0	ng	135704	5	14.009	917326	20.0
Cinnamyl cinnamate	13.727	110.0	ng	5043570	5	14.009	917326	20.0
1-Hexacosene	13.862	7.6	ng	348484	5	14.009	917326	20.0
Pentadecafluoro...	14.504	3.5	ng	161940	5	14.009	917326	20.0
Hentriacontane	15.133	2.2	ng	75950	6	15.468	694704	20.0
unknown15.715	15.715	2.0	ng	71243	6	15.468	694704	20.0
Eicosane	15.951	3.9	ng	135472	6	15.468	694704	20.0