

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
 Data File : BM029633.D
 Acq On : 24 Apr 2021 12:29
 Operator : CG/JU
 Sample : M2149-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-39-9.5-10.0

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_M\Methods\8270-BM042121.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM029633.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.222	390	397	406	rBV	1046549	1478408	42.40%	5.507%
2	6.805	659	666	677	rBV	973557	1597609	45.81%	5.951%
3	7.622	798	805	812	rBV	248007	384791	11.03%	1.433%
4	8.775	994	1001	1014	rBV	601962	1027564	29.47%	3.828%
5	10.404	1271	1278	1285	rBV	292646	487243	13.97%	1.815%
6	12.886	1693	1700	1719	rBV	1715630	2445729	70.14%	9.111%
7	13.733	1835	1844	1852	rBV	50957	93094	2.67%	0.347%
8	14.139	1908	1913	1924	rVB2	31332	74308	2.13%	0.277%
9	14.269	1929	1935	1943	rVB	465913	695910	19.96%	2.592%
10	14.845	2029	2033	2040	rVB	30938	49835	1.43%	0.186%
11	15.551	2144	2153	2159	rBV2	24139	76239	2.19%	0.284%
12	15.722	2176	2182	2183	rBV	25804	38188	1.10%	0.142%
13	15.763	2183	2189	2197	rVB	1554638	2083415	59.75%	7.761%
14	16.063	2234	2240	2247	rVB2	56685	91237	2.62%	0.340%
15	16.322	2279	2284	2288	rBV2	27739	46525	1.33%	0.173%
16	16.416	2293	2300	2310	rVB2	118021	265351	7.61%	0.988%
17	16.574	2322	2327	2335	rBV	29168	51217	1.47%	0.191%
18	17.016	2396	2402	2408	rBV	569586	818233	23.46%	3.048%
19	17.680	2511	2515	2522	rBV9	18562	37232	1.07%	0.139%
20	17.927	2549	2557	2564	rBV	892304	1379909	39.57%	5.140%
21	17.986	2564	2567	2575	rVB	139945	208984	5.99%	0.778%
22	18.198	2597	2603	2606	rBV	249836	395915	11.35%	1.475%
23	18.239	2606	2610	2617	rVB	331337	504885	14.48%	1.881%
24	19.210	2770	2775	2792	rVB	463257	730035	20.93%	2.719%
25	19.645	2844	2849	2854	rBV	2819054	3487169	100.00%	12.990%
26	19.721	2854	2862	2875	rVB2	304326	824358	23.64%	3.071%
27	19.939	2894	2899	2913	rVB2	163505	422037	12.10%	1.572%
28	20.662	3015	3022	3031	rBV	459718	971213	27.85%	3.618%
29	20.921	3062	3066	3072	rVB	282822	349577	10.02%	1.302%
30	21.198	3108	3113	3118	rBV	880388	1092399	31.33%	4.069%
31	21.415	3145	3150	3158	rBV	537555	939615	26.94%	3.500%
32	22.133	3267	3272	3280	rVB	500176	978386	28.06%	3.645%
33	22.933	3403	3408	3419	rVB	386862	910041	26.10%	3.390%
34	23.386	3479	3485	3494	rVB2	637658	1177507	33.77%	4.386%
35	23.845	3558	3563	3572	rVB	274741	630500	18.08%	2.349%

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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_M\Methods\8270-BM042121.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

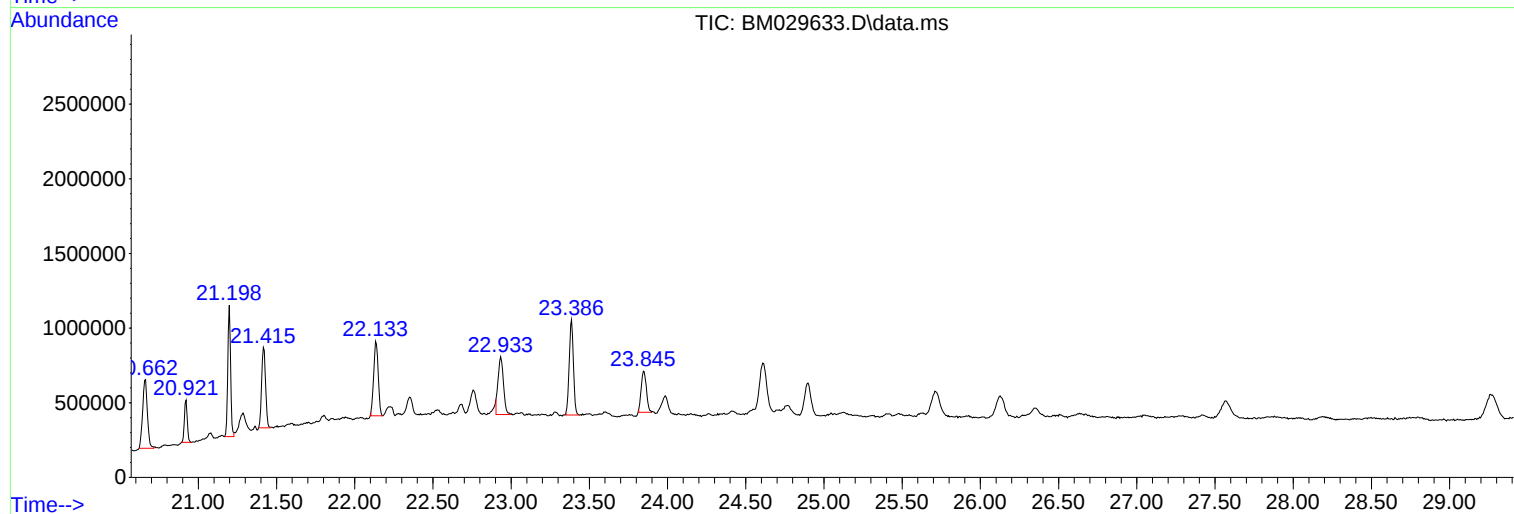
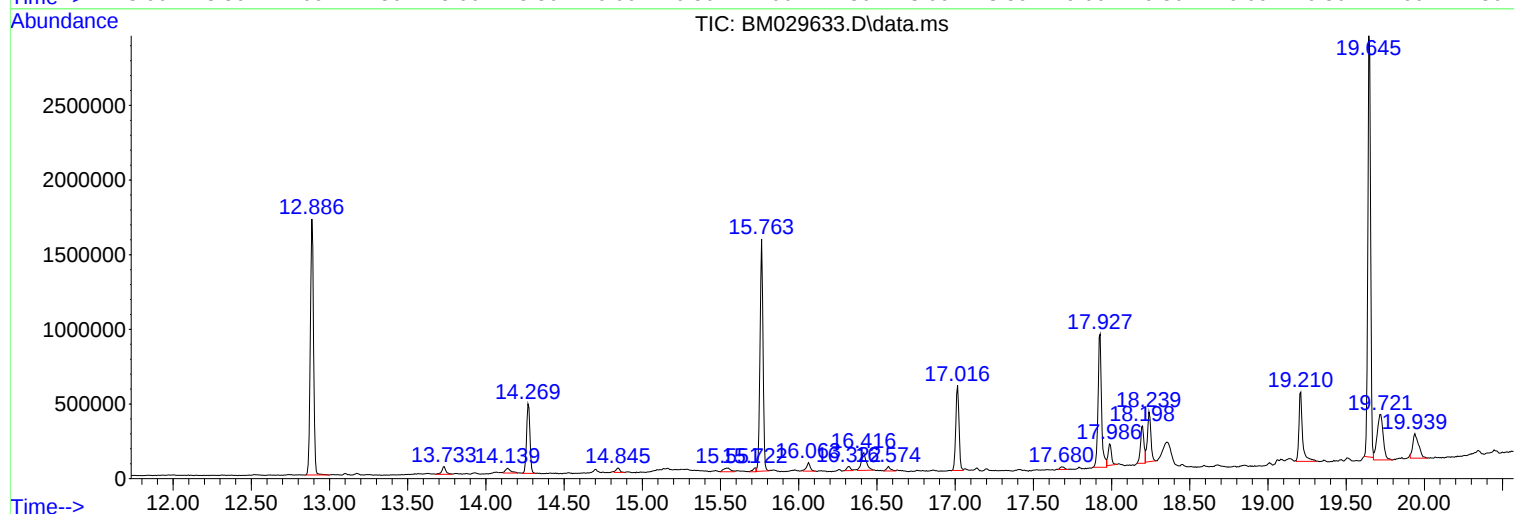
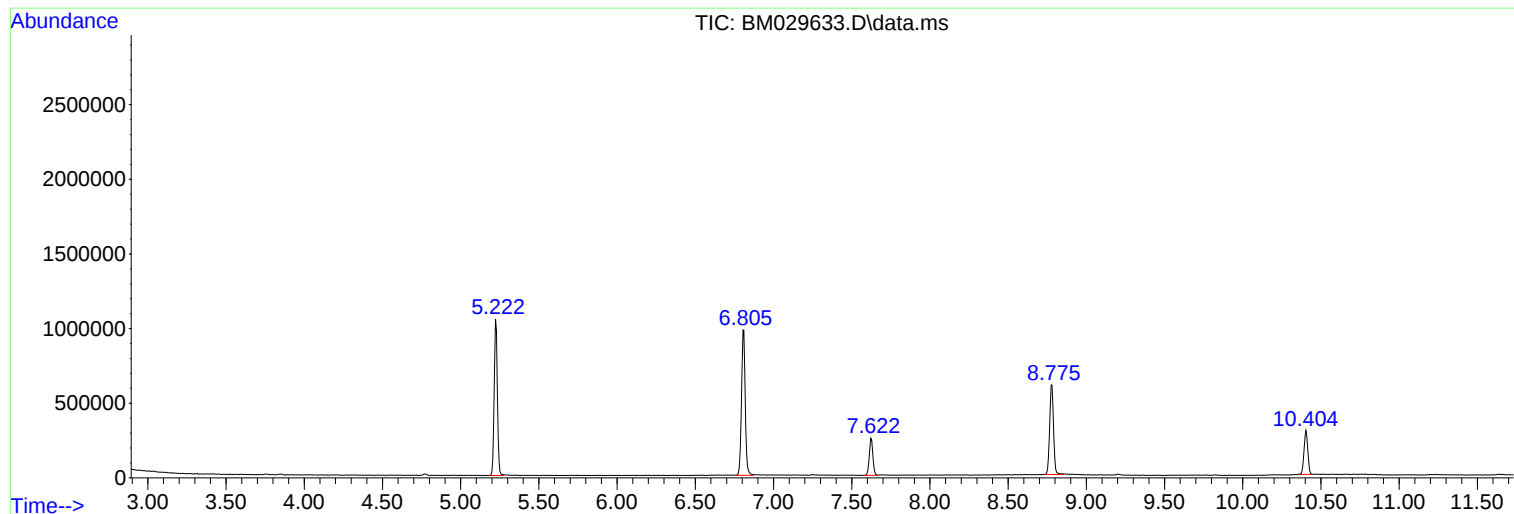
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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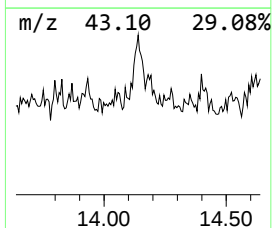
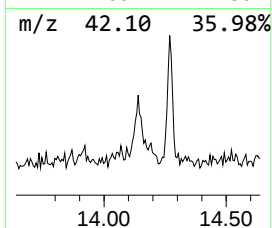
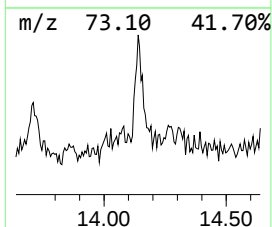
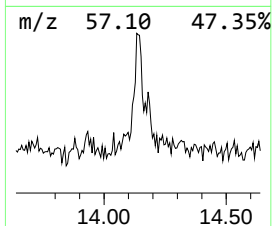
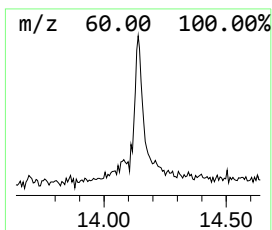
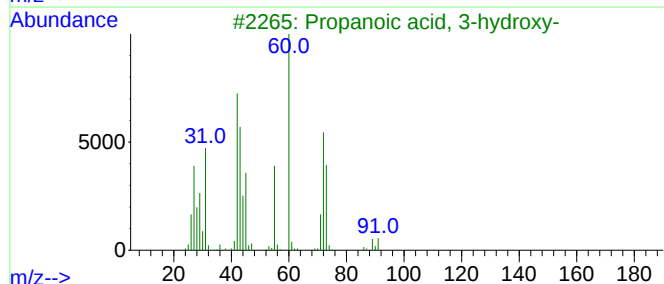
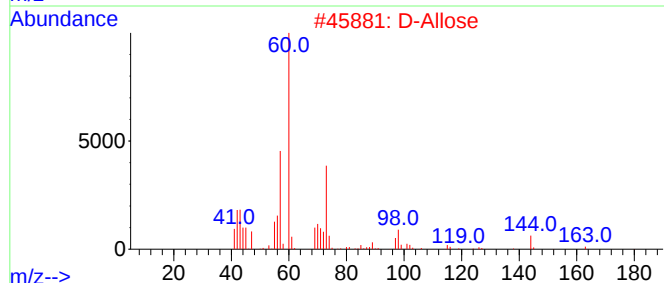
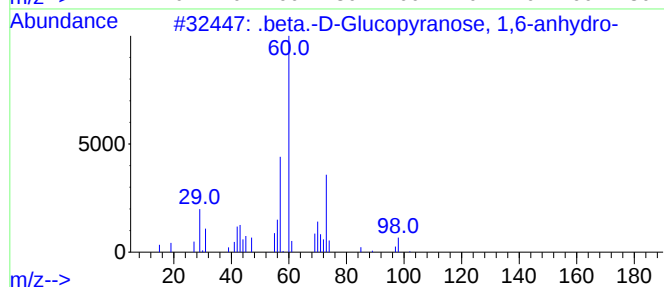
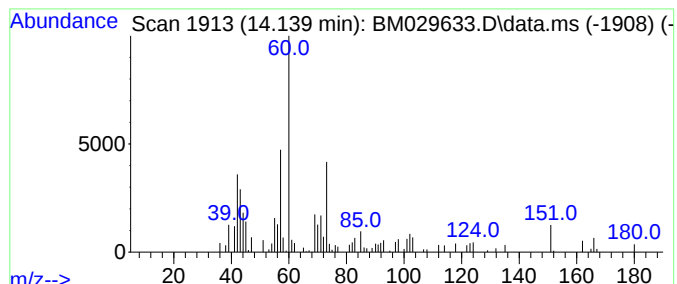
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 .beta.-D-Glucopyranose, 1,6... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.139	2.14 ng	74308	Acenaphthene-d10	14.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-D-Glucopyranose, 1,6-anhy...	162	C6H10O5	000498-07-7	53
2			D-Allose	180	C6H12O6	002595-97-3	49
3			Propanoic acid, 3-hydroxy-	90	C3H6O3	000503-66-2	47
4			Methyl-4-azido-4-desoxy.beta.l-a...	189	C6H11N3O4	1000312-10-3	47
5			Ethyl .alpha.-d-glucopyranoside	208	C8H16O6	1000127-29-4	45



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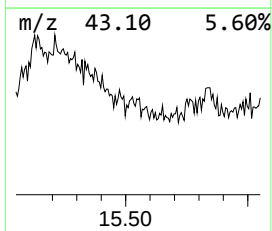
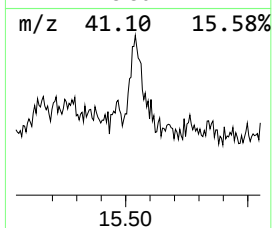
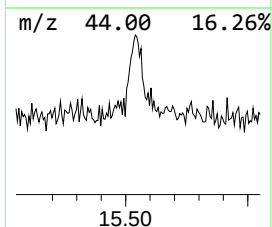
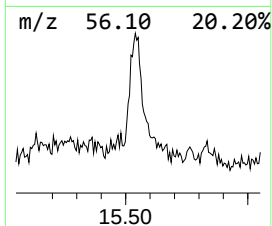
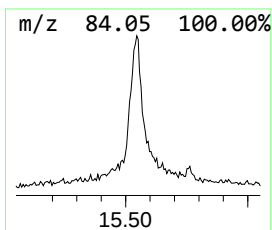
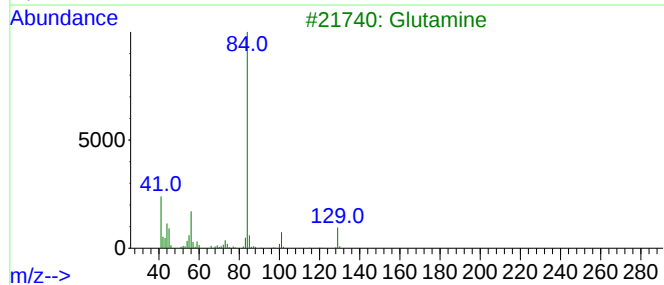
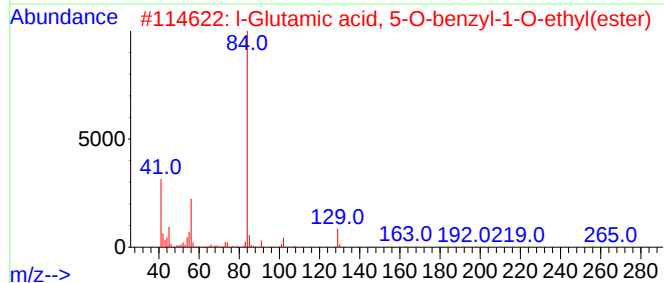
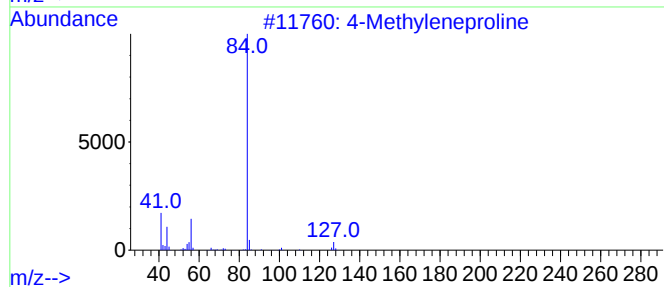
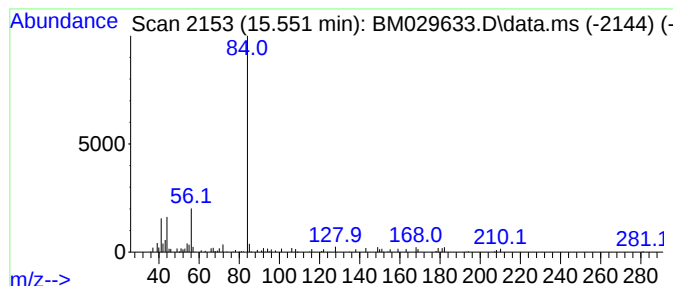
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Peak Number 2 4-Methyleneproline Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.551	2.19 ng	76239	Acenaphthene-d10	14.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methyleneproline	127	C6H9NO2	1000131-14-7	64
2			l-Glutamic acid, 5-O-benzyl-1-O-...	265	C14H19NO4	1000130-30-6	64
3			Glutamine	146	C5H10N2O3	006899-04-3	56
4			Pyrrolidine-5-one, 2-[3-hydroxyp...	143	C7H13NO2	080243-73-8	53
5			Propane, 2-isocyanato-2-methyl-	99	C5H9NO	001609-86-5	50



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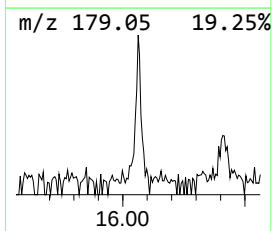
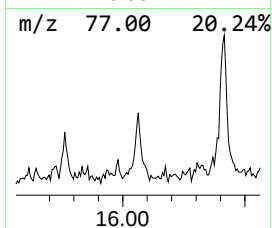
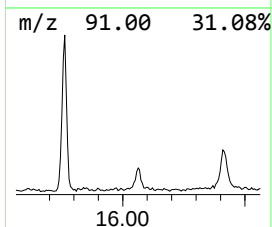
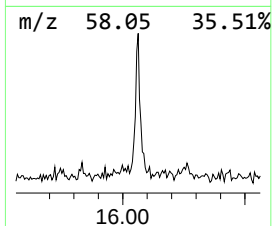
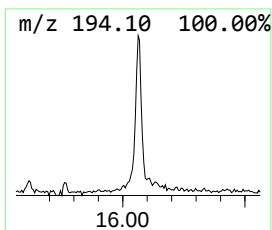
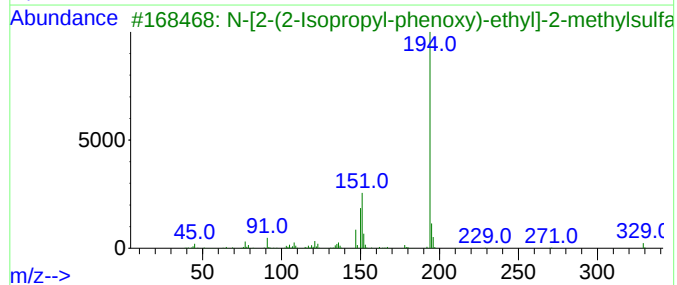
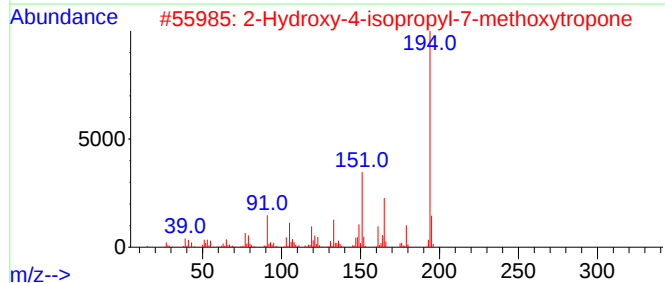
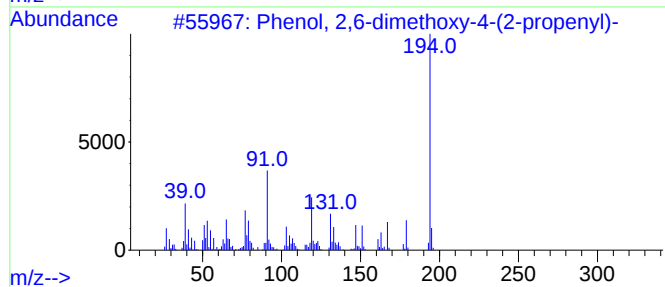
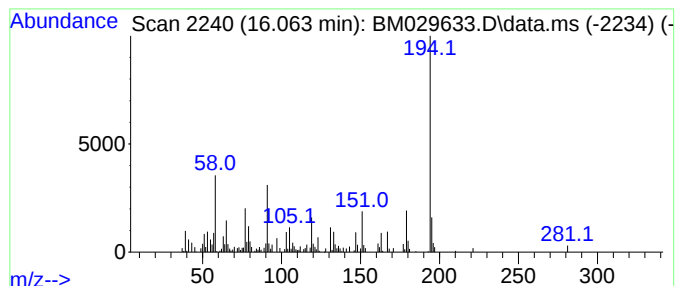
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Phenol, 2,6-dimethoxy-4-(2-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	2.23 ng	91237	Phenanthrene-d10	17.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-dimethoxy-4-(2-prope...	194	C11H14O3	006627-88-9	87
2			2-Hydroxy-4-isopropyl-7-methoxyt...	194	C11H14O3	089647-85-8	64
3			N-[2-(2-Isopropyl-phenoxy)-ethyl]...	329	C19H23NO2S	1000275-37-1	52
4			2-Propenoic acid, 3-(4-hydroxy-3...	194	C10H10O4	000537-98-4	50
5			2-Propenoic acid, 3-(4-hydroxy-3...	194	C10H10O4	001135-24-6	50



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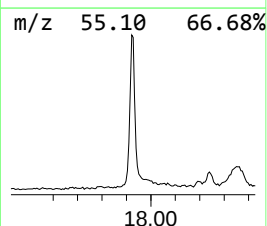
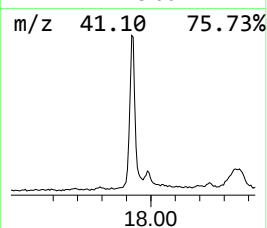
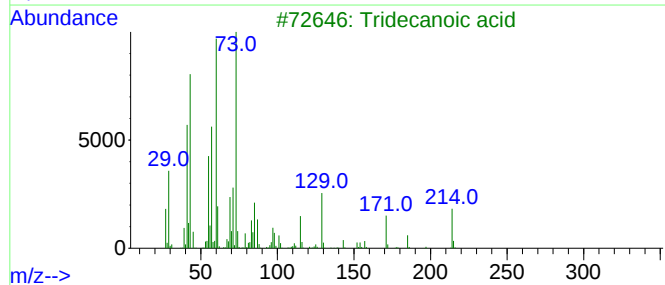
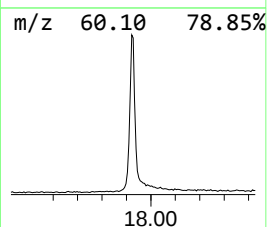
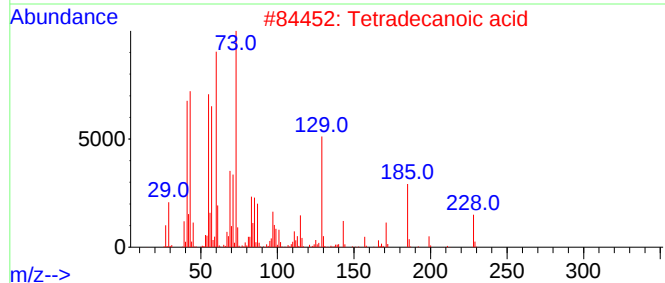
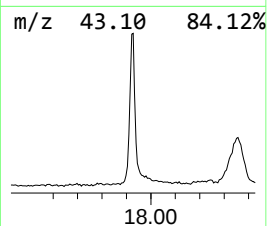
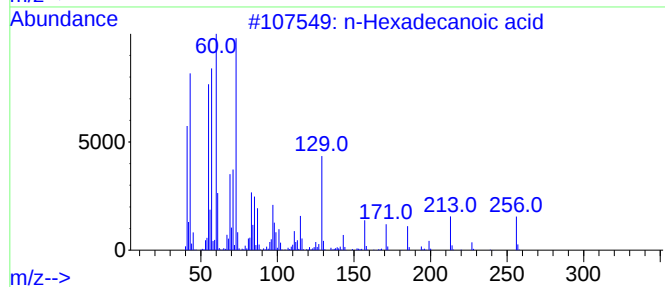
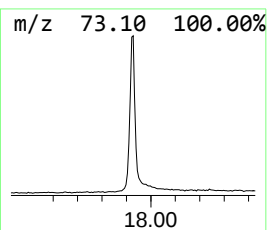
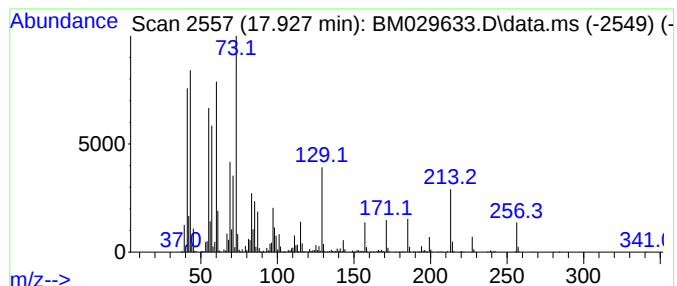
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.927	33.73 ng	1379910	Phenanthrene-d10	17.016

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	91
3			Tridecanoic acid	214	C13H26O2	000638-53-9	90
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			n-Decanoic acid	172	C10H20O2	000334-48-5	53



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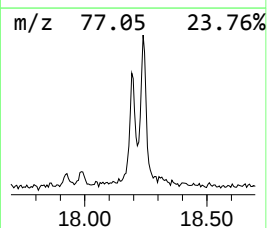
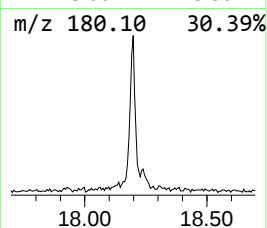
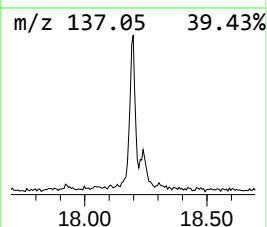
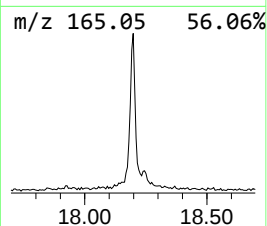
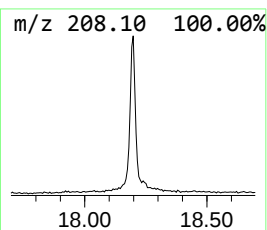
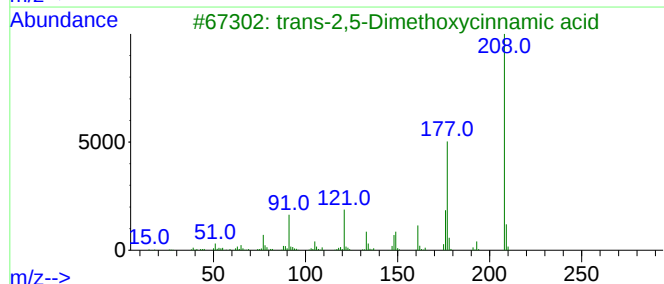
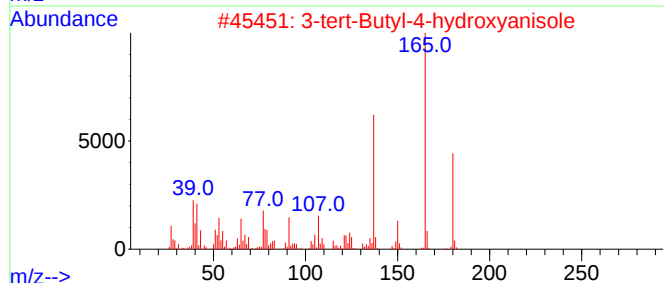
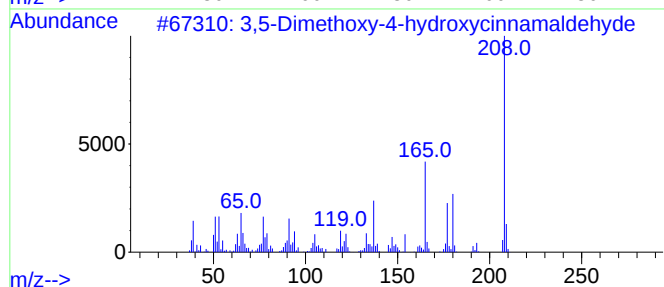
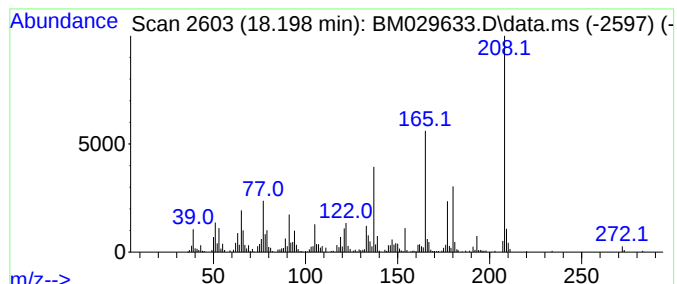
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.198	9.68 ng	395915	Phenanthrene-d10	17.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Dimethoxy-4-hydroxycinnamal... 208	C11H12O4	087345-53-7	92	
2			3-tert-Butyl-4-hydroxyanisole 180	C11H16O2	000121-00-6	42	
3			trans-2,5-Dimethoxycinnamic acid 208	C11H12O4	038489-74-6	38	
4			Benzene, 1,2,3-trimethoxy-5-(2-p... 208	C12H16O3	000487-11-6	38	
5			Asarone 208	C12H16O3	002883-98-9	38	



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
Data File : BM029633.D
Acq On : 24 Apr 2021 12:29
Operator : CG/JU
Sample : M2149-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

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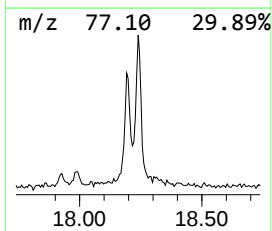
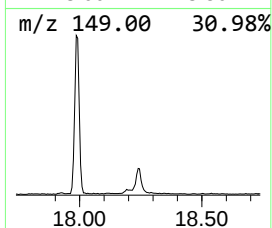
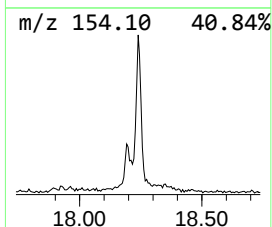
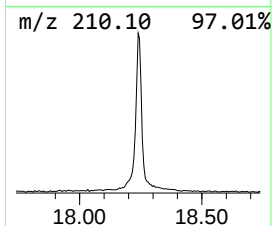
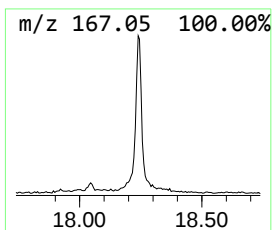
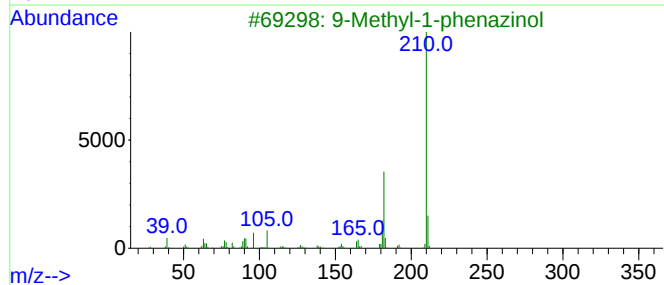
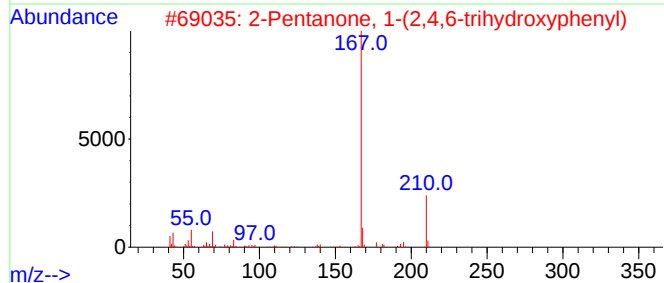
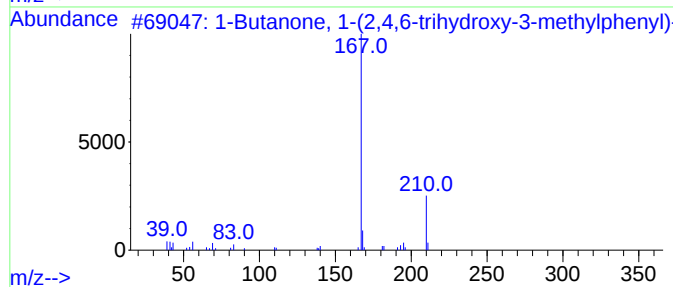
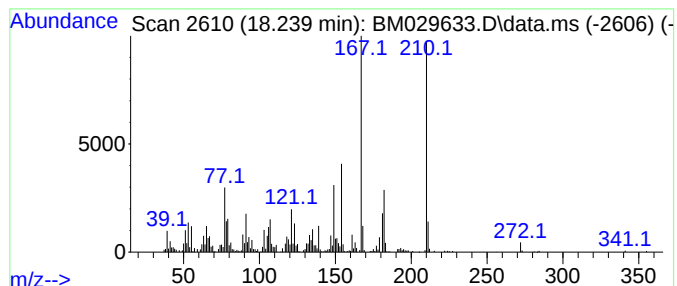
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1-Butanone, 1-(2,4,6-trihyd... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	12.34 ng	504885	Phenanthrene-d10	17.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butanone, 1-(2,4,6-trihydroxy-...	210	C11H14O4	001509-06-4	52
2			2-Pentanone, 1-(2,4,6-trihydroxy...	210	C11H14O4	1000116-22-3	50
3			9-Methyl-1-phenazinol	210	C13H10N2O	013860-43-0	38
4			Imidazo[4,5-e][1,4]diazepin-8(1H...	210	C8H10N4OS	1000142-41-3	35
5			Benzene, 1,1'-ethyldienebis-	182	C14H14	000612-00-0	30



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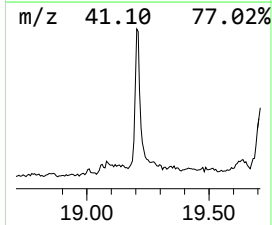
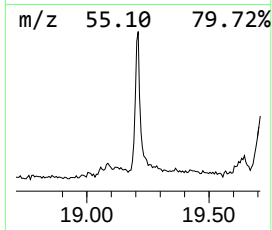
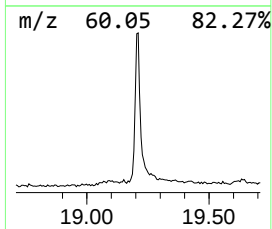
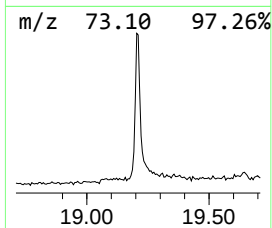
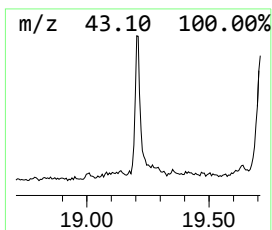
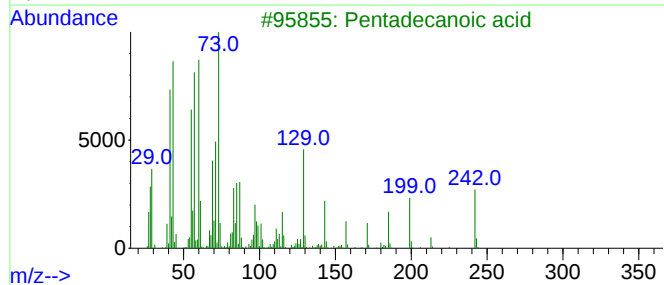
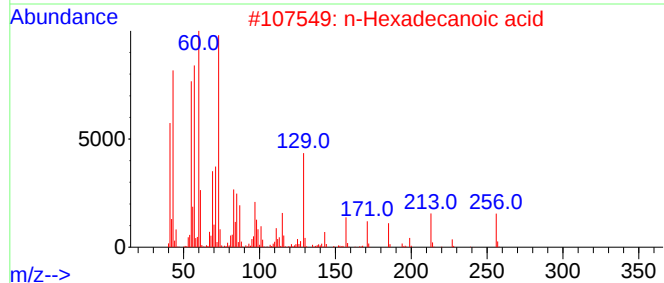
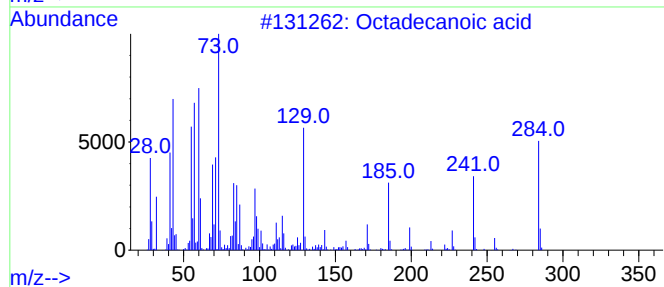
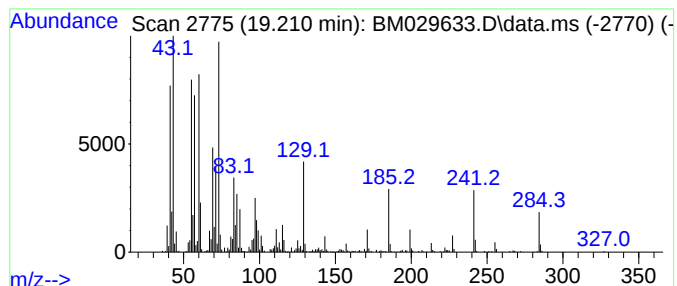
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.210	13.37 ng	730035	Chrysene-d12	21.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72



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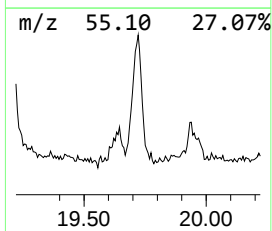
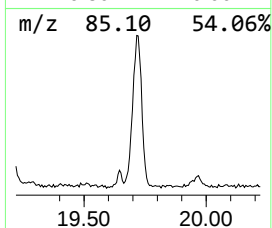
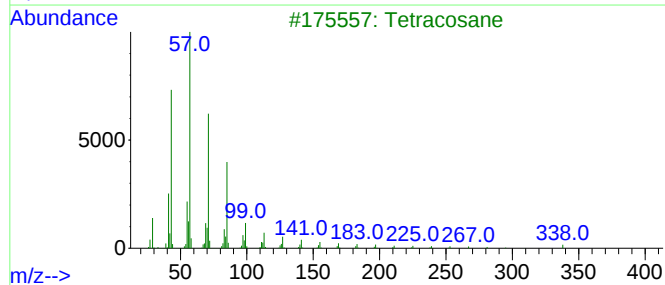
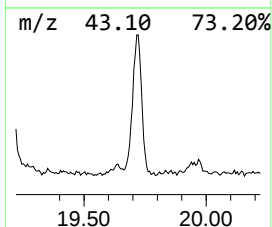
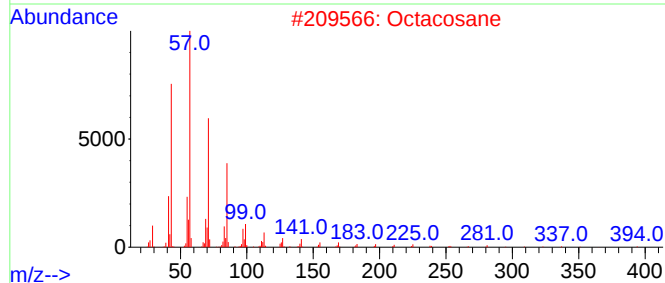
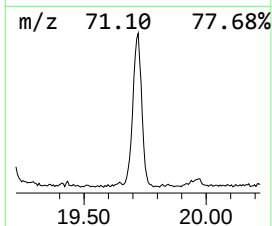
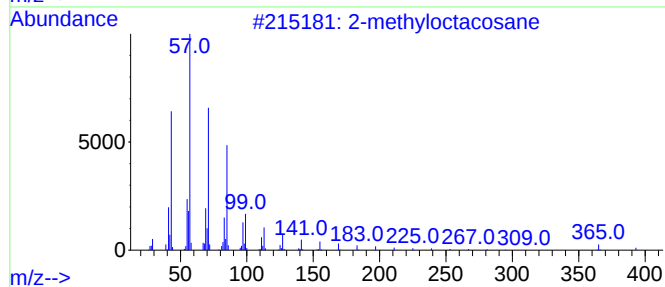
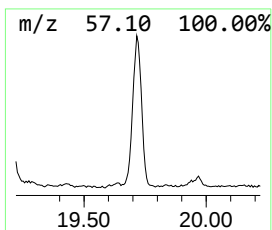
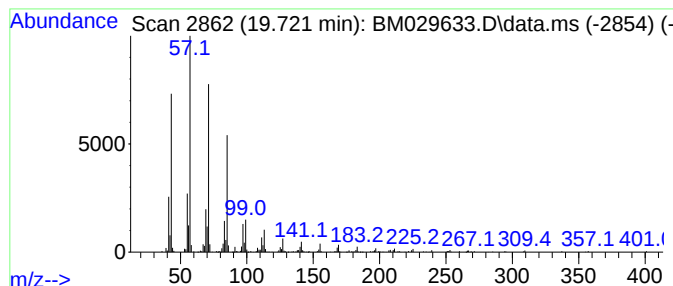
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 2-methyloctacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.721	15.09 ng	824358	Chrysene-d12	21.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Docosane	310	C22H46	000629-97-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
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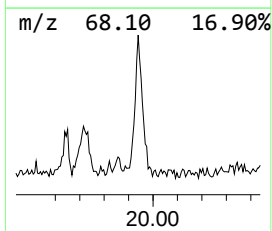
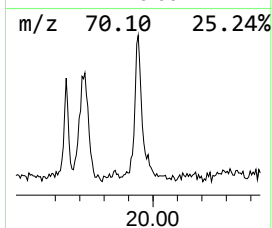
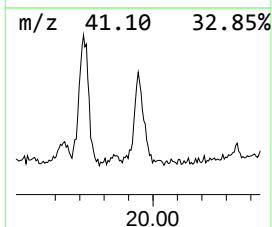
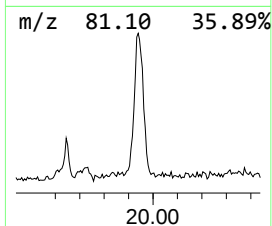
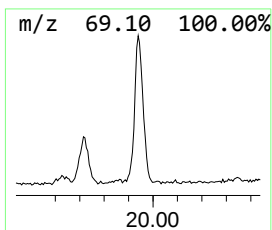
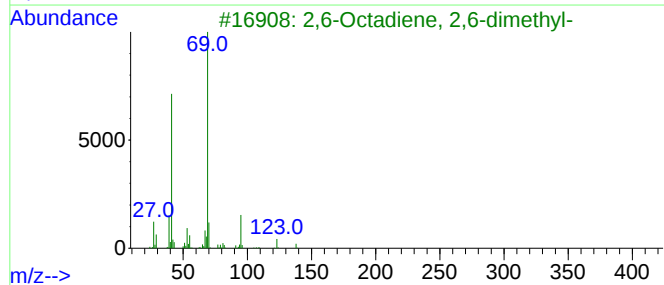
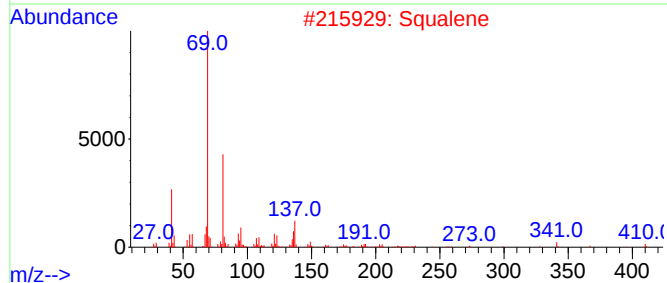
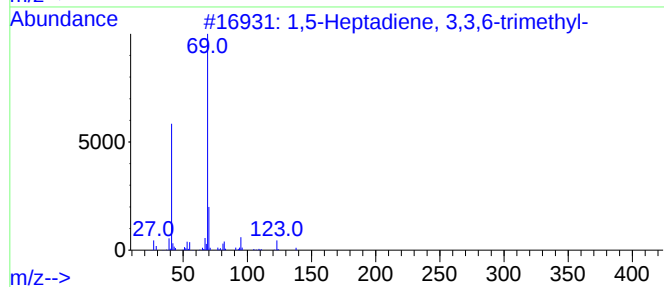
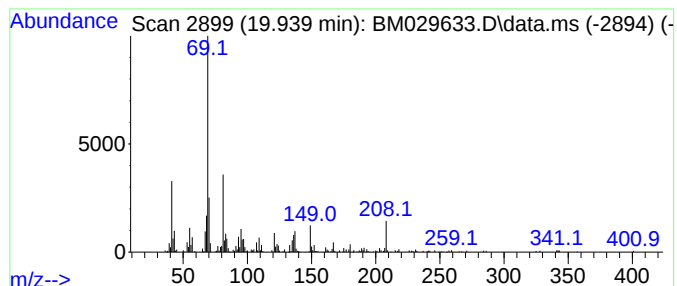
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1,5-Heptadiene, 3,3,6-trime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.939	7.73 ng	422037	Chrysene-d12	21.198	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	53
2	Squalene	410	C30H50	000111-02-4	50
3	2,6-Octadiene, 2,6-dimethyl-	138	C10H18	002792-39-4	47
4	3,7,11-Trimethyl-dodeca-2,6,10-t...	236	C15H24O2	007548-13-2	47
5	4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	47



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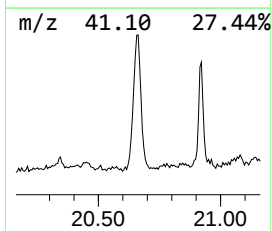
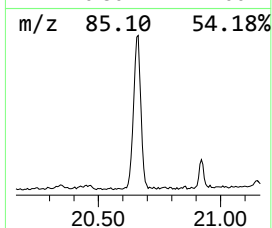
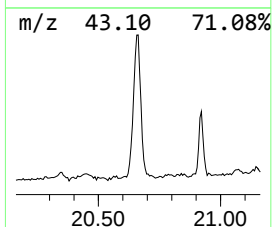
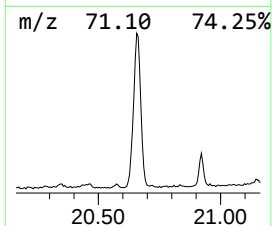
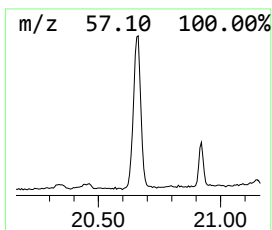
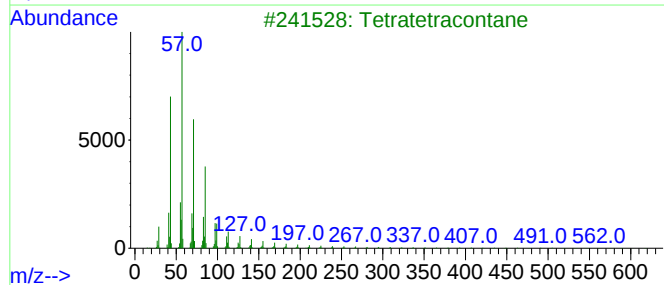
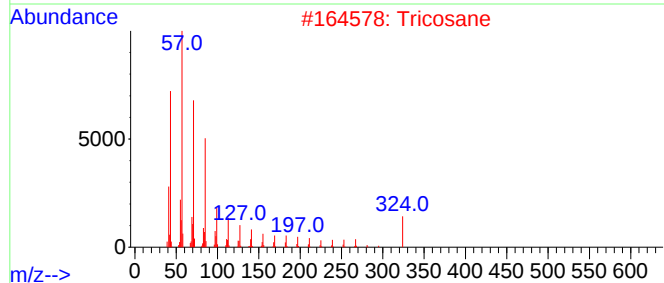
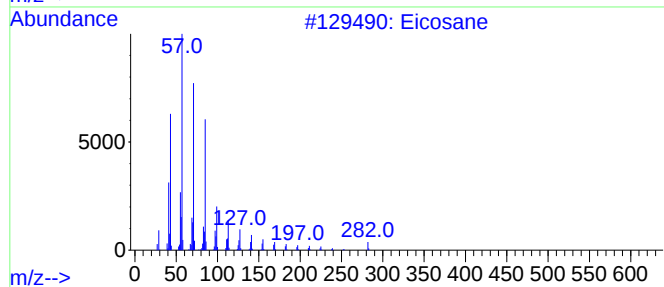
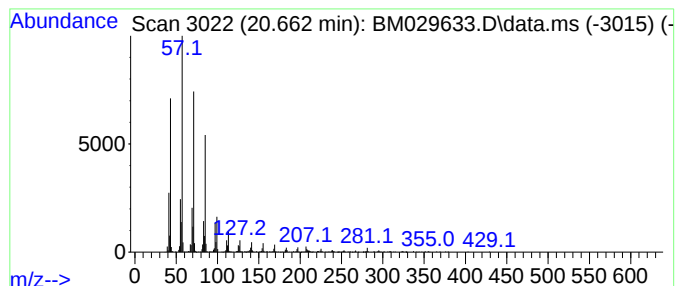
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Eicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.662	17.78 ng	971213	Chrysene-d12	21.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	97
2	Tricosane			324	C23H48	000638-67-5	94
3	Tetratetracontane			619	C44H90	007098-22-8	94
4	2-Bromo dodecane			248	C12H25Br	013187-99-0	93
5	Hexacosane			366	C26H54	000630-01-3	91



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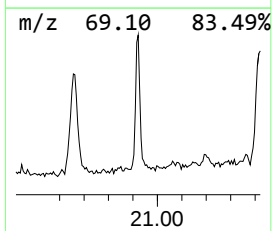
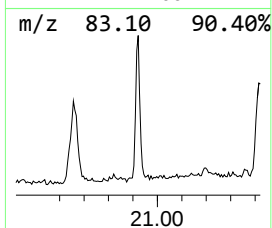
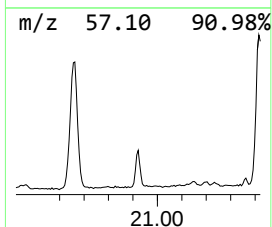
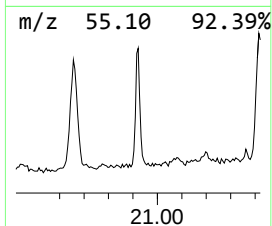
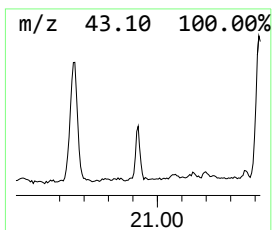
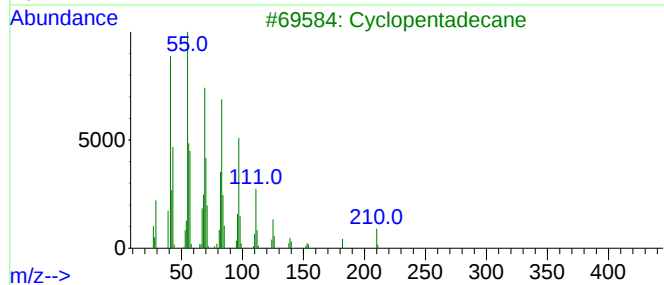
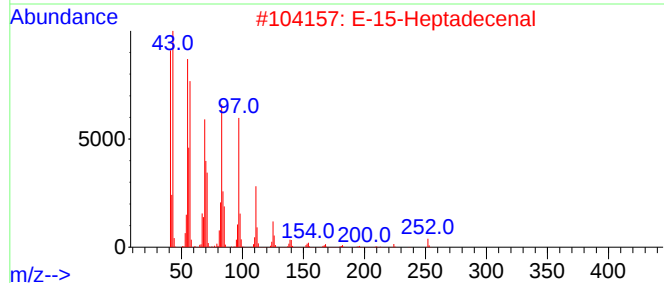
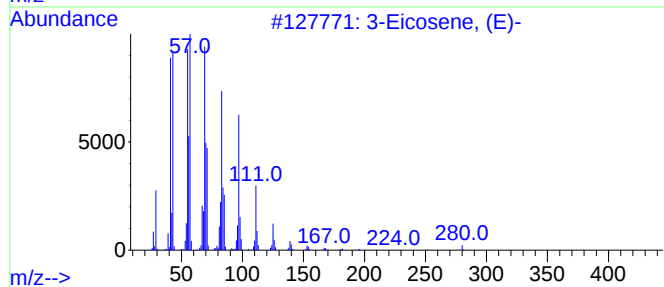
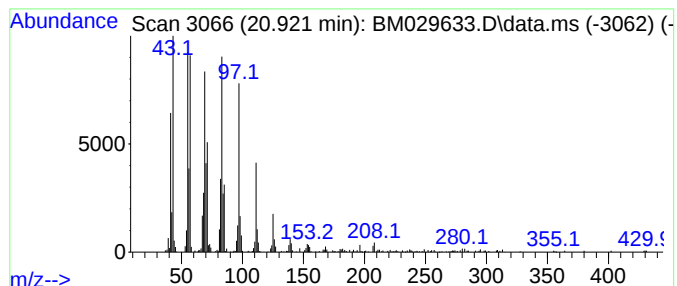
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 3-Eicosene, (E)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.921	6.40 ng	349577	Chrysene-d12	21.198

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Eicosene, (E)-	280	C20H40	074685-33-9	97
2		E-15-Heptadecenal	252	C17H32O	1000130-97-9	94
3		Cyclopentadecane	210	C15H30	000295-48-7	93
4		n-Heptadecanol-1	256	C17H36O	001454-85-9	93
5		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



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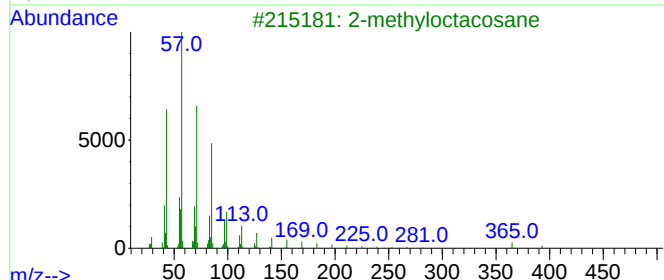
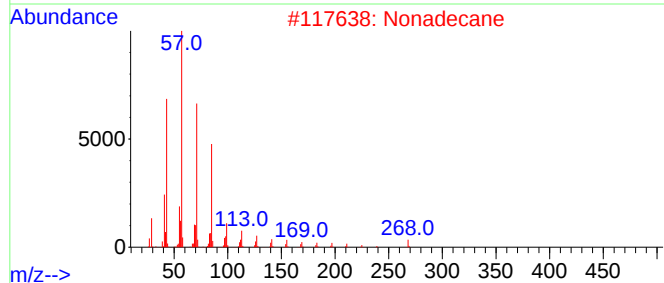
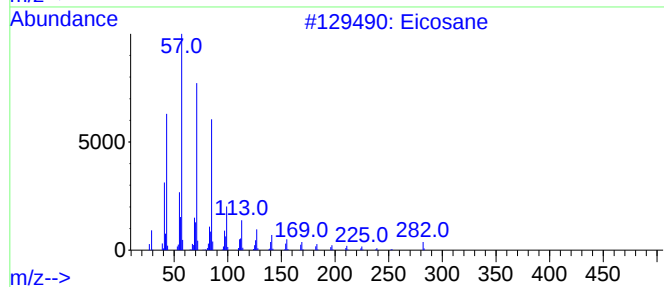
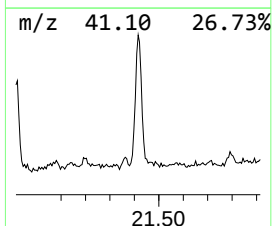
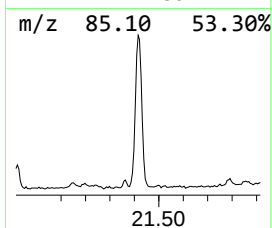
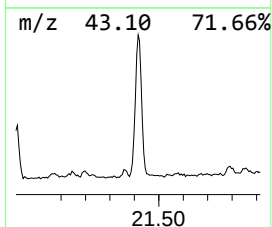
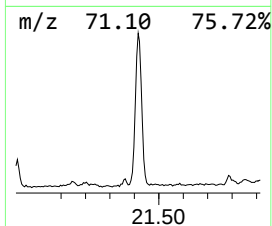
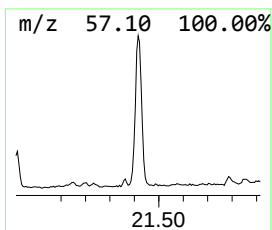
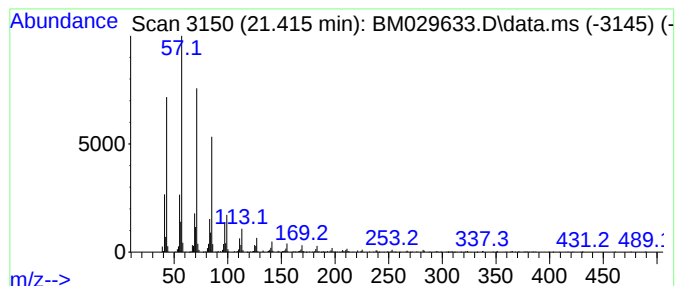
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Nonadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.415	17.20 ng	939615	Chrysene-d12	21.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Nonadecane	268	C19H40	000629-92-5	93
3			2-methyloctacosane	408	C29H60	1000376-72-8	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
Data File : BM029633.D
Acq On : 24 Apr 2021 12:29
Operator : CG/JU
Sample : M2149-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SB-39-9.5-10.0

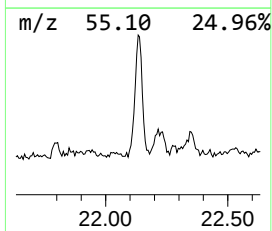
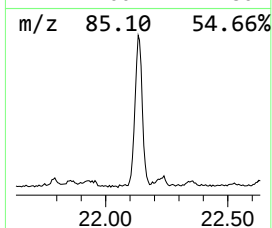
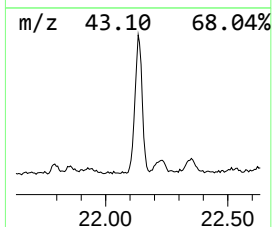
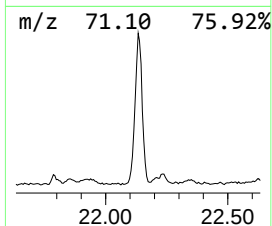
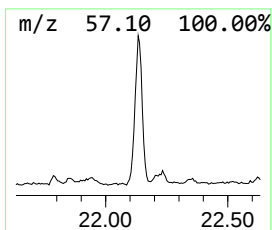
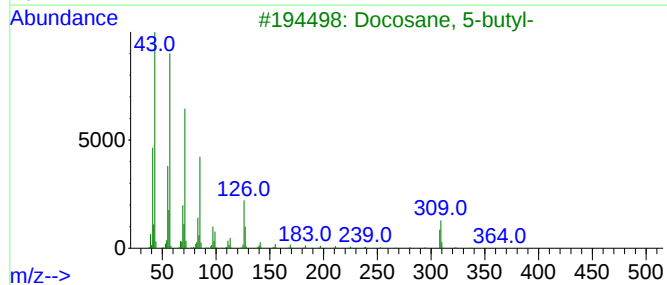
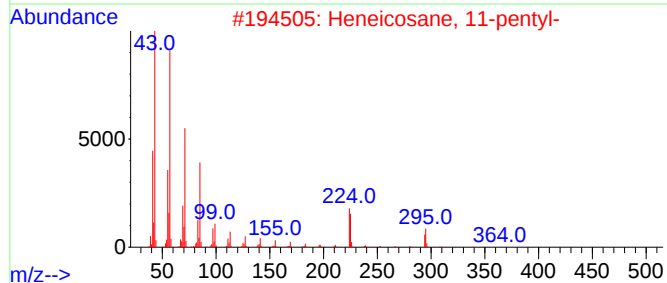
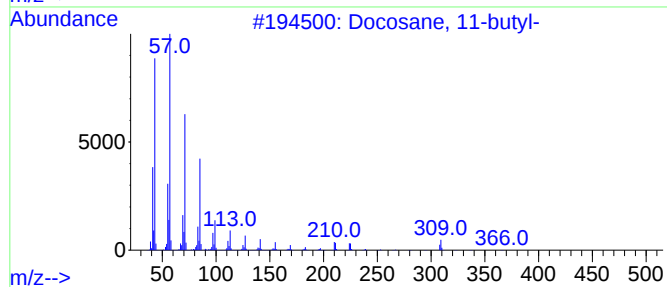
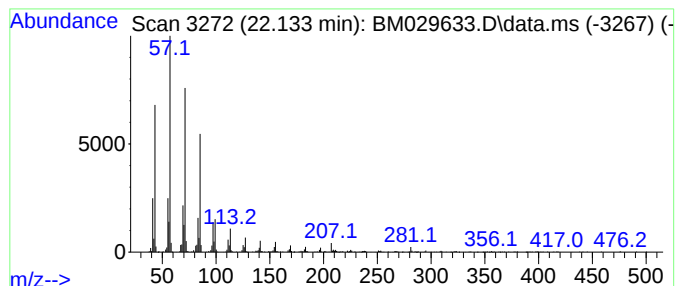
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Docosane, 11-butyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.133	17.91 ng	978386	Chrysene-d12	21.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
3			Docosane, 5-butyl-	366	C26H54	055282-16-1	94
4			Hexacosane	366	C26H54	000630-01-3	94
5			Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
Data File : BM029633.D
Acq On : 24 Apr 2021 12:29
Operator : CG/JU
Sample : M2149-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SB-39-9.5-10.0

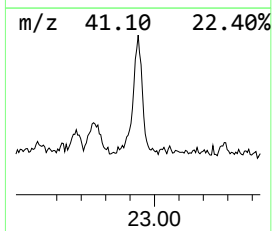
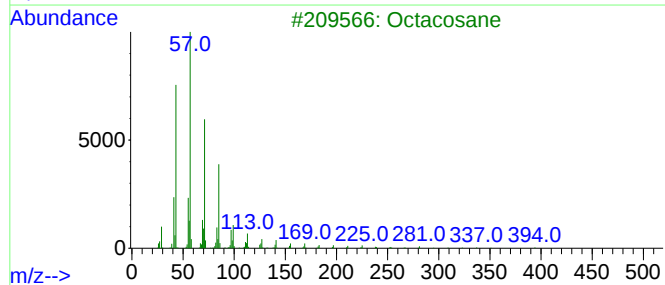
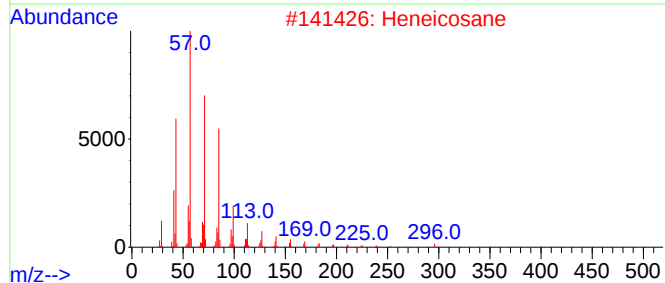
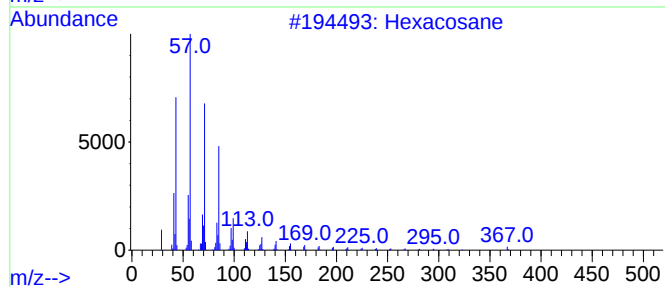
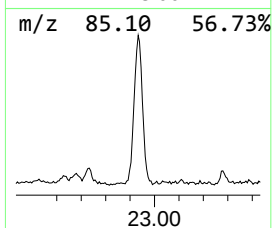
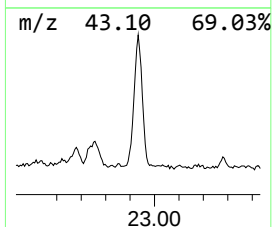
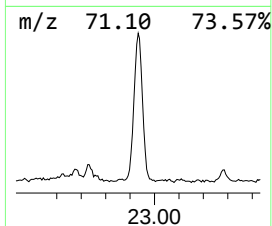
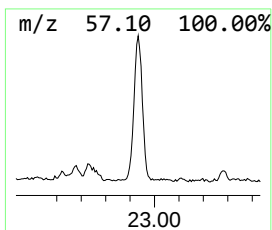
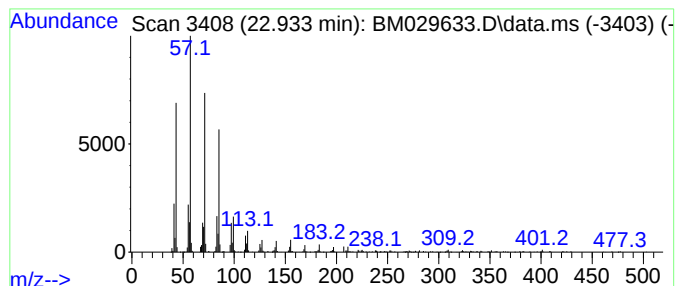
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.933	15.46 ng	910041	Perylene-d12	23.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Octacosane	394	C28H58	000630-02-4	91
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	90
5			Tridecane, 6-propyl-	226	C16H34	055045-10-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
Data File : BM029633.D
Acq On : 24 Apr 2021 12:29
Operator : CG/JU
Sample : M2149-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-39-9.5-10.0

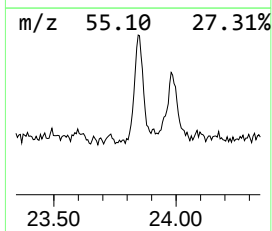
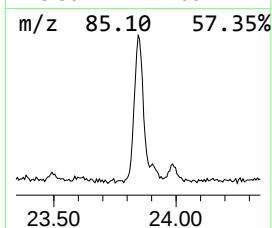
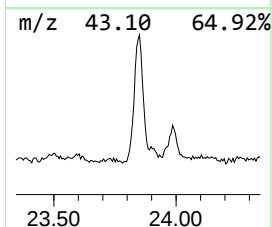
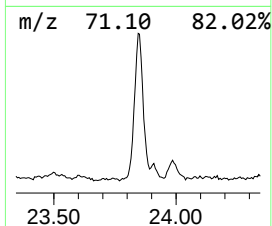
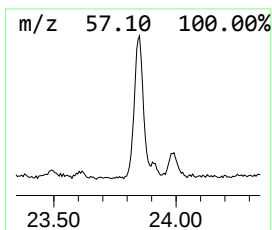
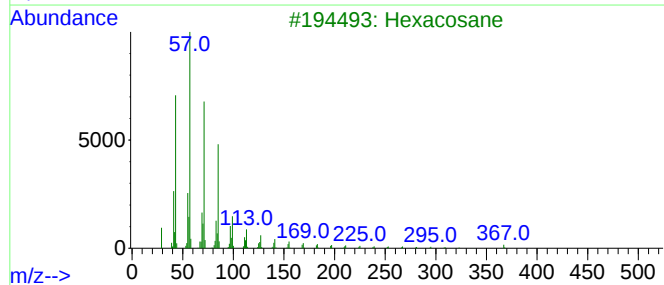
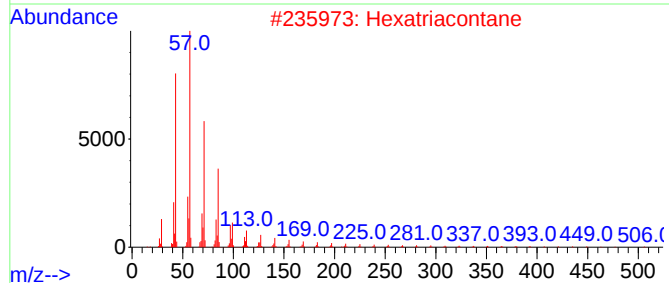
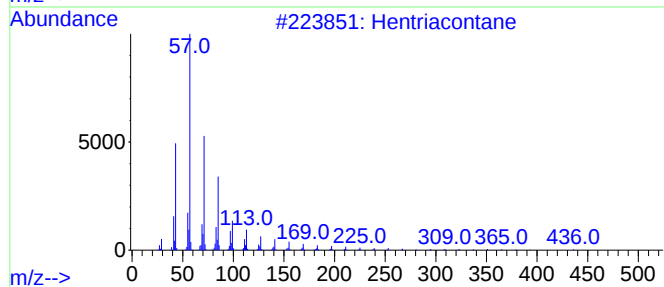
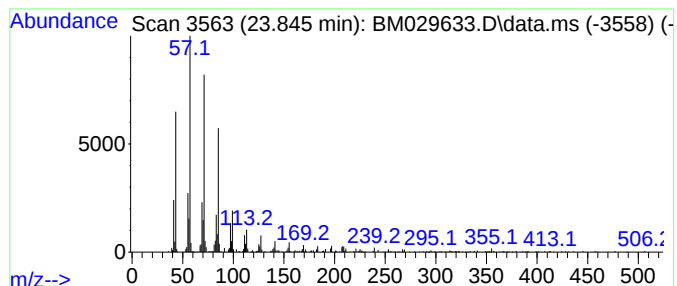
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Hentriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.845	10.71 ng	630500	Perylene-d12	23.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	95
2			Hexatriacontane	507	C36H74	000630-06-8	95
3			Hexacosane	366	C26H54	000630-01-3	94
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	93
5			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042421\
 Data File : BM029633.D
 Acq On : 24 Apr 2021 12:29
 Operator : CG/JU
 Sample : M2149-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-39-9.5-10.0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
.beta.-D-Glucop...	14.139	2.1	ng	74308	3	14.269	695910	20.0
4-Methyleneproline	15.551	2.2	ng	76239	3	14.269	695910	20.0
Phenol, 2,6-dim...	16.063	2.2	ng	91237	4	17.016	818233	20.0
4-((1E)-3-Hydro...	16.416	6.5	ng	265351	4	17.016	818233	20.0
n-Hexadecanoic ...	17.927	33.7	ng	1379910	4	17.016	818233	20.0
3,5-Dimethoxy-4...	18.198	9.7	ng	395915	4	17.016	818233	20.0
1-Butanone, 1-(...	18.239	12.3	ng	504885	4	17.016	818233	20.0
Octadecanoic acid	19.210	13.4	ng	730035	5	21.198	1092400	20.0
2-methyloctacosane	19.721	15.1	ng	824358	5	21.198	1092400	20.0
1,5-Heptadiene,...	19.939	7.7	ng	422037	5	21.198	1092400	20.0
Eicosane	20.662	17.8	ng	971213	5	21.198	1092400	20.0
3-Eicosene, (E)-	20.921	6.4	ng	349577	5	21.198	1092400	20.0
Nonadecane	21.415	17.2	ng	939615	5	21.198	1092400	20.0
Docosane, 11-bu...	22.133	17.9	ng	978386	5	21.198	1092400	20.0
Hexacosane	22.933	15.5	ng	910041	6	23.386	1177510	20.0
Hentriacontane	23.845	10.7	ng	630500	6	23.386	1177510	20.0