

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
 Data File : BM031817.D
 Acq On : 18 Aug 2021 22:08
 Operator : CG/JU
 Sample : M3438-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CHRT-26855

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-BM081621.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM031817.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.175	348	355	364	rBV	813566	1156533	31.03%	6.650%
2	6.734	613	620	632	rBV	847470	1297007	34.80%	7.458%
3	7.540	751	757	766	rVB	159303	252460	6.77%	1.452%
4	8.687	944	952	965	rBV	497124	829134	22.24%	4.768%
5	9.722	1120	1128	1134	rBV2	33315	52412	1.41%	0.301%
6	10.098	1186	1192	1204	rBV	117611	227028	6.09%	1.305%
7	10.298	1219	1226	1233	rVB	218427	355769	9.54%	2.046%
8	12.786	1643	1649	1658	rVB	1249643	1871211	50.20%	10.760%
9	13.080	1693	1699	1704	rBV	30946	48787	1.31%	0.281%
10	13.580	1779	1784	1791	rBV3	33585	61444	1.65%	0.353%
11	13.904	1833	1839	1844	rBV7	24458	56432	1.51%	0.324%
12	13.992	1850	1854	1862	rVB2	46367	74362	2.00%	0.428%
13	14.069	1862	1867	1872	rBV5	21148	39234	1.05%	0.226%
14	14.127	1872	1877	1880	rBV2	29811	47312	1.27%	0.272%
15	14.175	1880	1885	1893	rVB	355922	522663	14.02%	3.005%
16	14.257	1893	1899	1906	rBV6	28830	66408	1.78%	0.382%
17	14.716	1973	1977	1981	rBV4	30389	46439	1.25%	0.267%
18	14.804	1987	1992	1993	rBV4	27998	43727	1.17%	0.251%
19	14.827	1993	1996	2002	rVV4	50855	77014	2.07%	0.443%
20	14.892	2003	2007	2014	rVB7	43741	86091	2.31%	0.495%
21	14.963	2014	2019	2025	rBV	93541	149629	4.01%	0.860%
22	15.674	2135	2140	2147	rBV	1230967	1770474	47.50%	10.180%
23	16.933	2349	2354	2358	rBV	518255	777291	20.85%	4.470%
24	17.851	2507	2510	2519	rVB3	264310	452623	12.14%	2.603%
25	18.998	2702	2705	2713	rVB	207921	290841	7.80%	1.672%
26	19.139	2726	2729	2733	rVB2	121777	132379	3.55%	0.761%
27	19.362	2764	2767	2770	rVB	150581	177542	4.76%	1.021%
28	19.580	2799	2804	2809	rVB	3164134	3727365	100.00%	21.433%
29	20.868	3019	3023	3027	rBV2	166209	200266	5.37%	1.152%
30	21.127	3061	3067	3071	rBV2	681246	959584	25.74%	5.518%
31	21.162	3071	3073	3081	rVB	119810	153604	4.12%	0.883%
32	21.745	3169	3172	3177	rVB2	63803	87419	2.35%	0.503%
33	22.380	3277	3280	3291	rVB4	60180	97122	2.61%	0.558%
34	22.639	3320	3324	3329	rBV	131923	241105	6.47%	1.386%
35	23.180	3412	3416	3423	rVB	94843	154281	4.14%	0.887%
36	23.274	3426	3432	3438	rVV2	394848	708793	19.02%	4.076%

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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

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37	23.780	3516	3518	3526	rVB3	62194	99100	2.66%	0.570%
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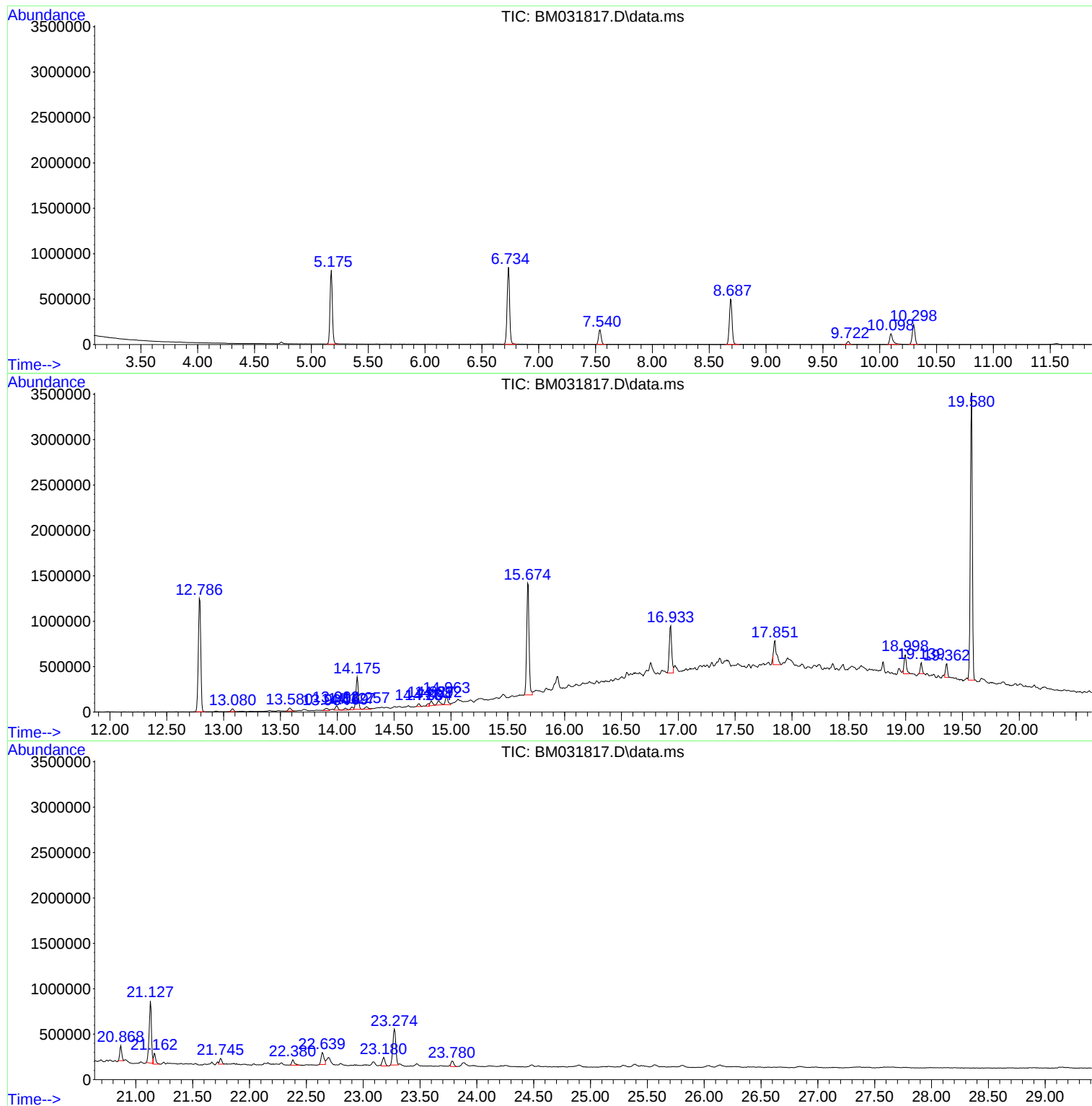
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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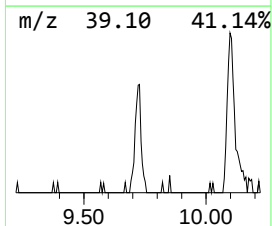
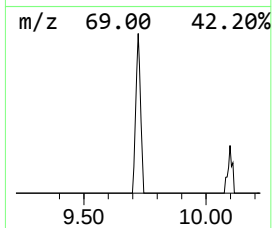
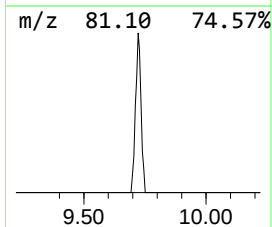
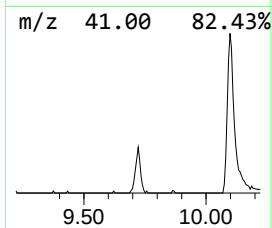
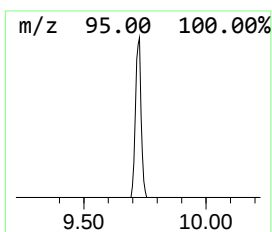
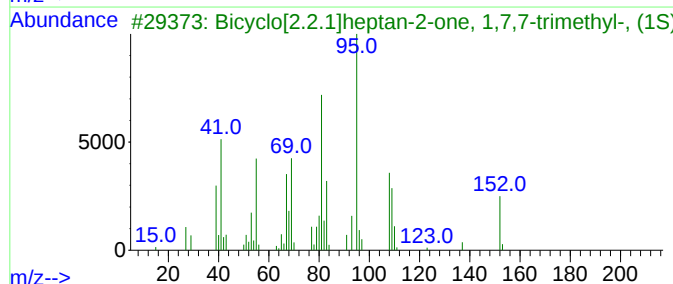
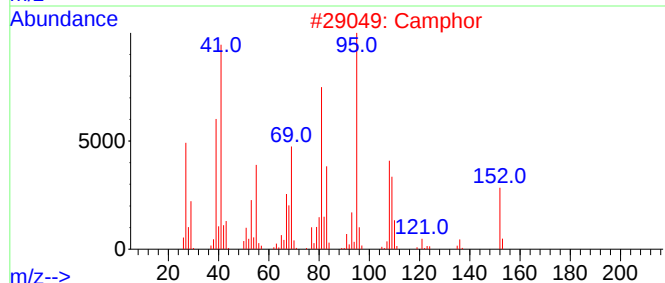
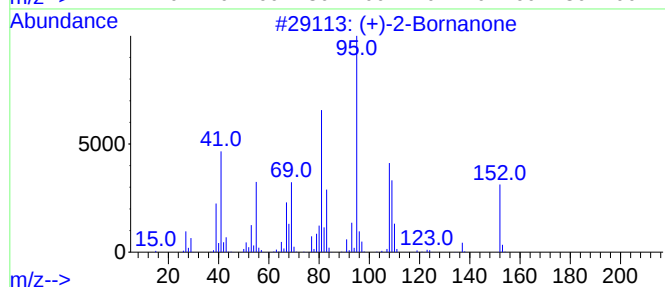
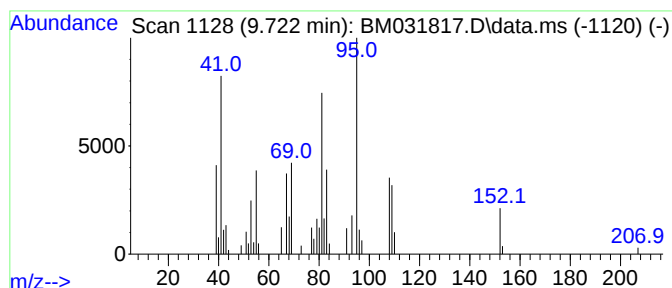
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (+)-2-Bornanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.722	2.95 ng	52412	Naphthalene-d8	10.298

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		(+)-2-Bornanone	152	C10H16O	000464-49-3	97
2		Camphor	152	C10H16O	000076-22-2	96
3		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	95
4		Carane, 4,5-epoxy-, trans	152	C10H16O	006909-20-2	46
5		5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	006752-80-3	43



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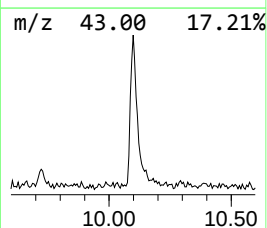
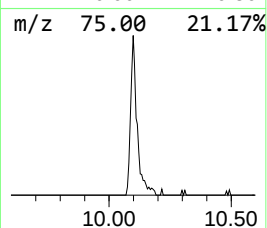
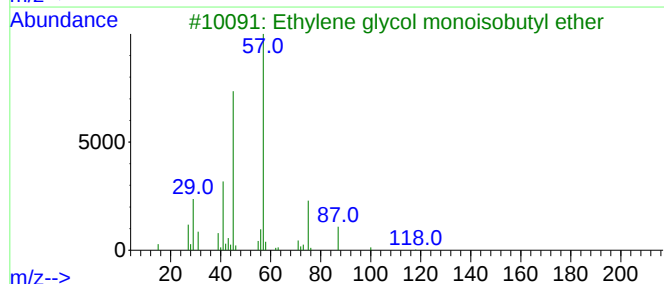
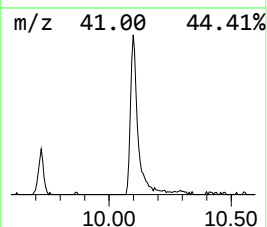
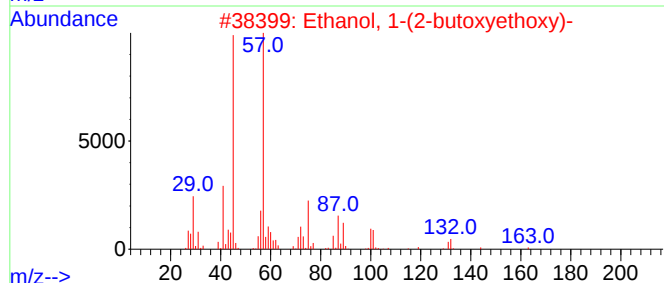
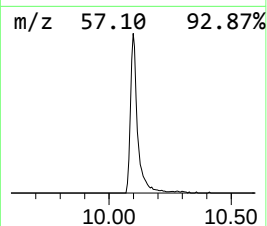
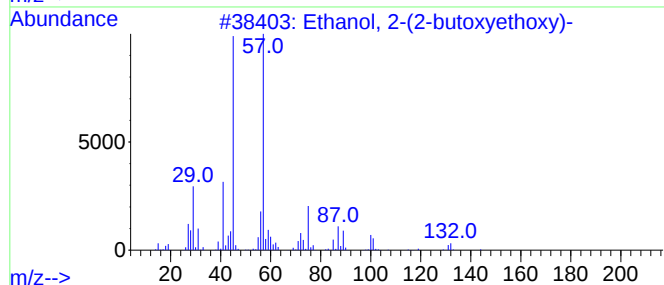
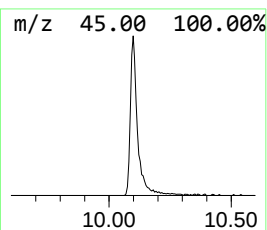
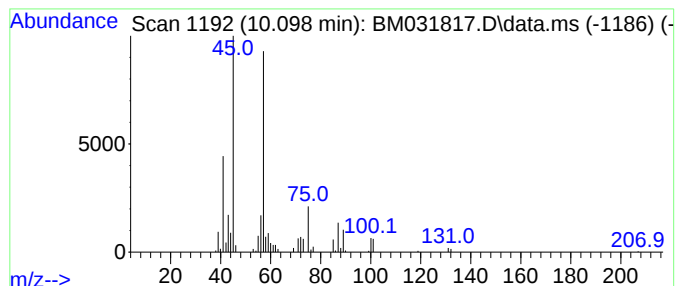
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.098	12.76 ng	227028	Naphthalene-d8	10.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	59
5			Di-sec-Butyl ether	130	C8H18O	006863-58-7	50



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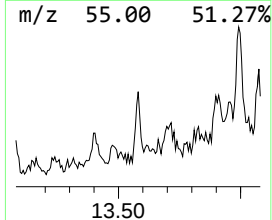
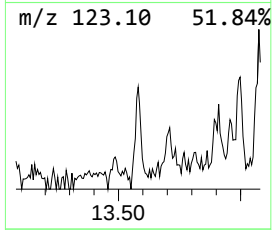
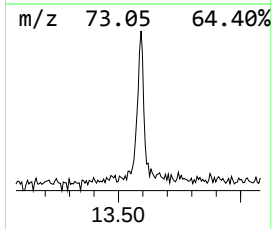
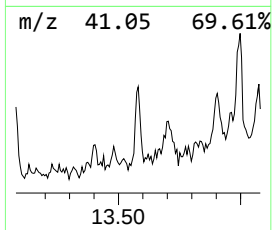
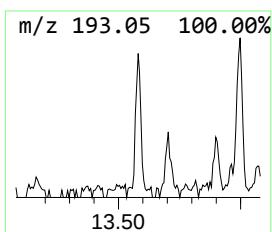
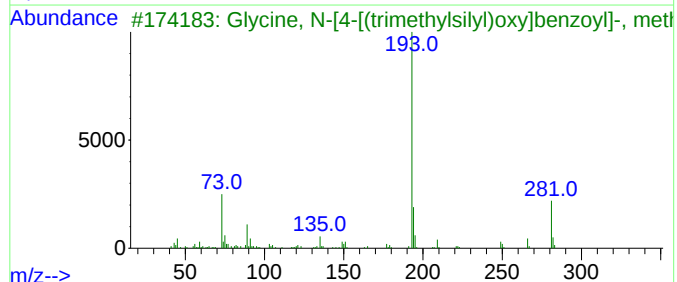
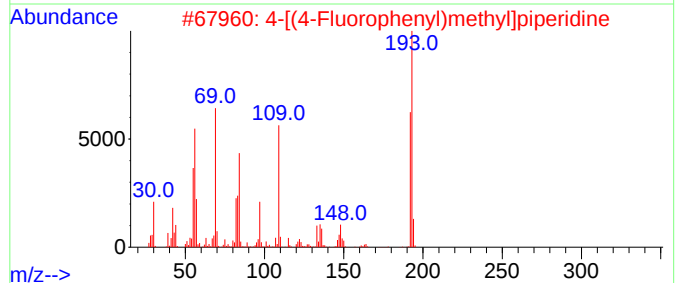
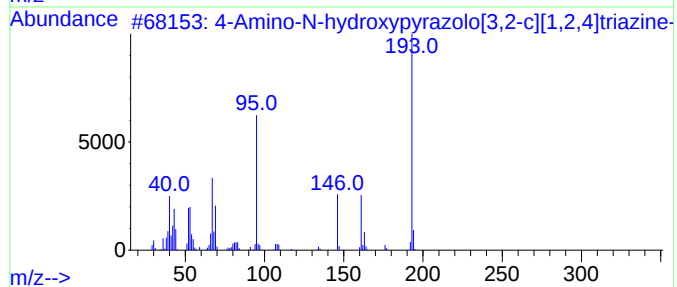
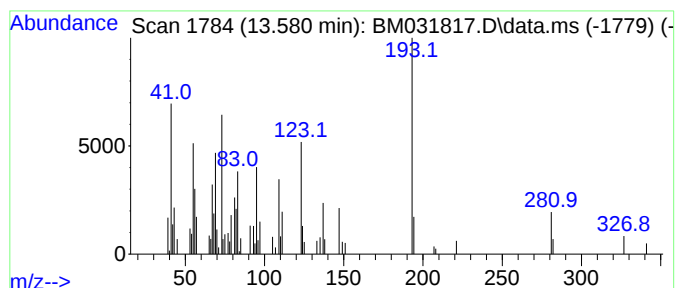
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 unknown13.580 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.580	2.35 ng	61444	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Amino-N-hydroxypyrazolo[3,2-c]...	193	C6H7N7O	1000443-50-2	27
2			4-[(4-Fluorophenyl)methyl]piperi...	193	C12H16FN	092822-02-1	27
3			Glycine, N-[4-[(trimethylsilyl)o...	281	C13H19NO4Si	055638-48-7	25
4			2-Diethylamino-4-phenylthiooct-2...	302	C18H26N2S	1000090-57-0	22
5			3-(3-Pyridyl)-D-alanine, 3TMS	382	C17H34N2O2Si3	1000512-77-7	22



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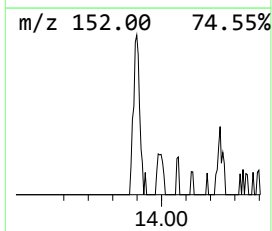
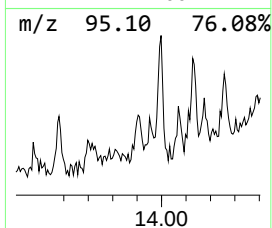
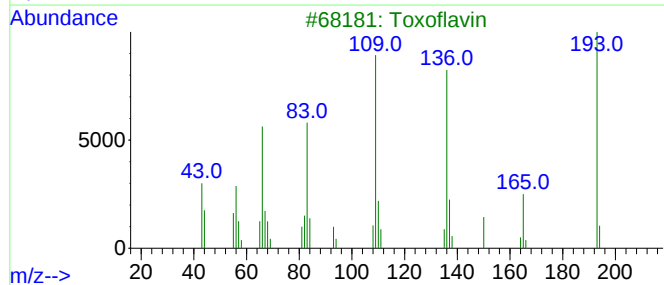
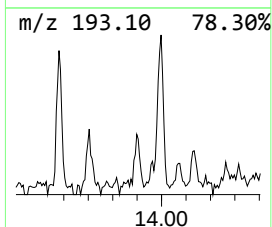
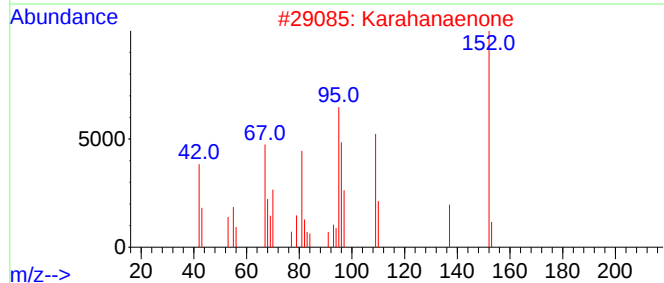
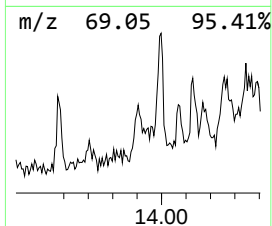
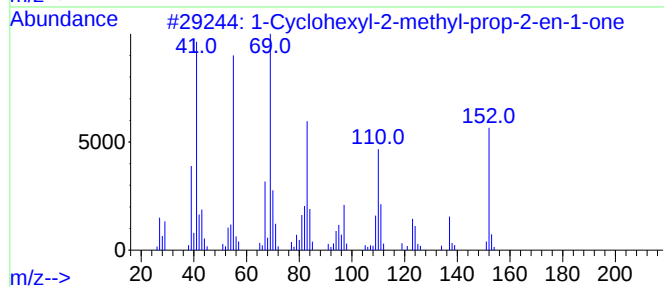
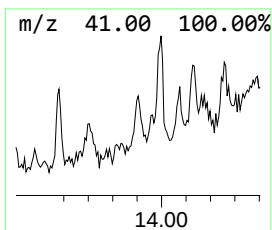
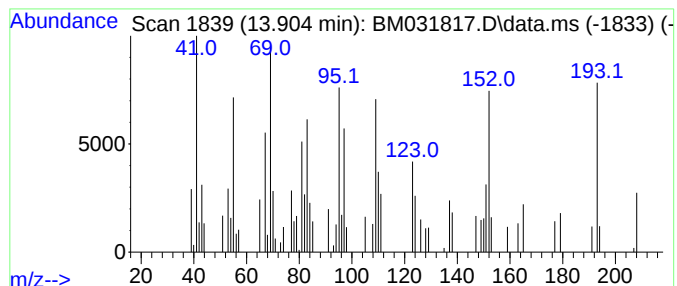
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 unknown13.904 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	2.16 ng	56432	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Cyclohexyl-2-methyl-prop-2-en-...	152	C10H16O	025183-82-8	41
2			Karahanaenone	152	C10H16O	1010427-29-1	41
3			Toxoflavin	193	C7H7N5O2	000084-82-2	35
4			Camphor	152	C10H16O	000076-22-2	30
5			(+)-2-Bornanone	152	C10H16O	000464-49-3	30



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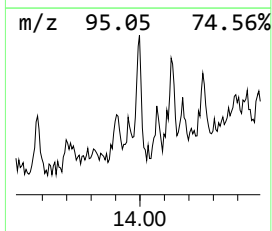
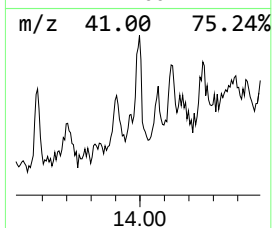
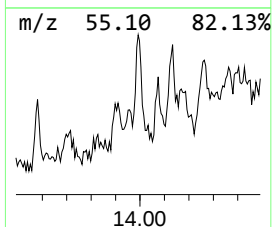
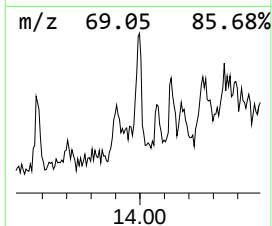
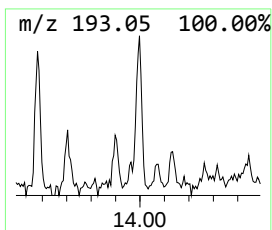
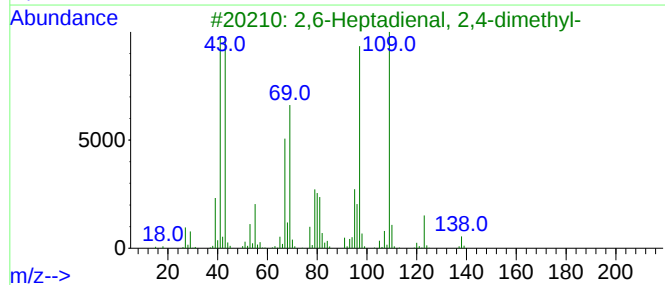
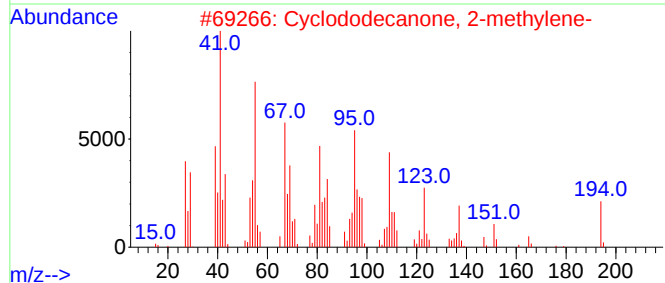
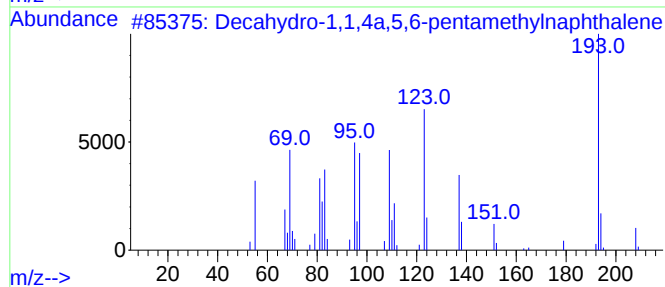
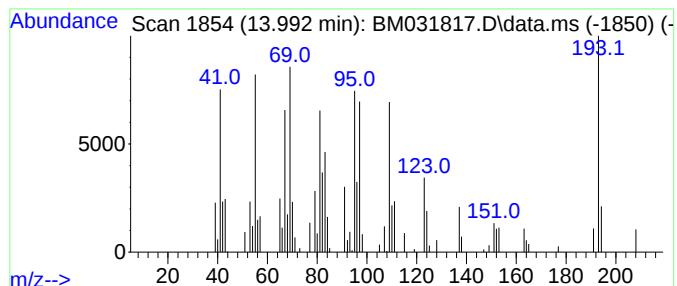
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Peak Number 5 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.992	2.85 ng	74362	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	70
2			Cyclododecanone, 2-methylene-	194	C13H22O	003045-76-9	46
3			2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	41
4			1,2-Dimethyl-4-oxocyclohex-2-ene...	152	C9H12O2	1000187-01-2	35
5			Phosphetane, 1-chloro-2,2,3,4,4-...	194	C8H16ClOP	029276-11-7	30



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Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
Operator : CG/JU
Sample : M3438-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

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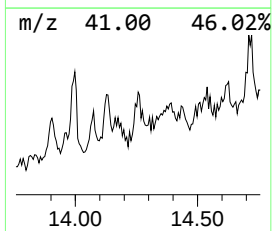
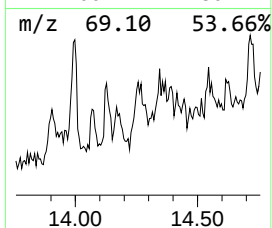
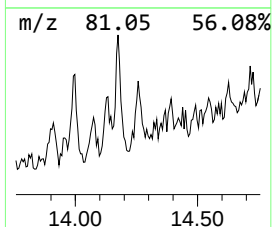
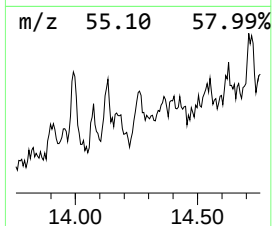
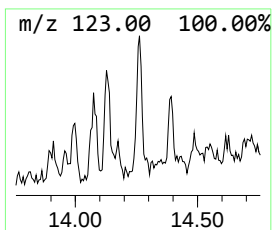
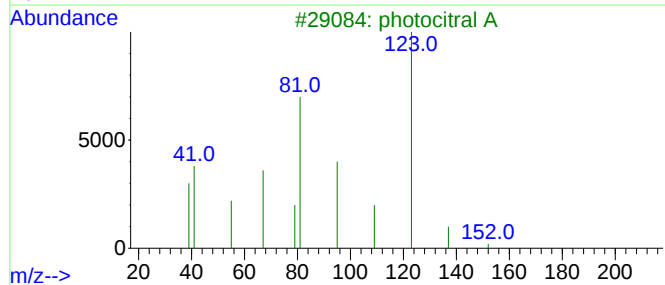
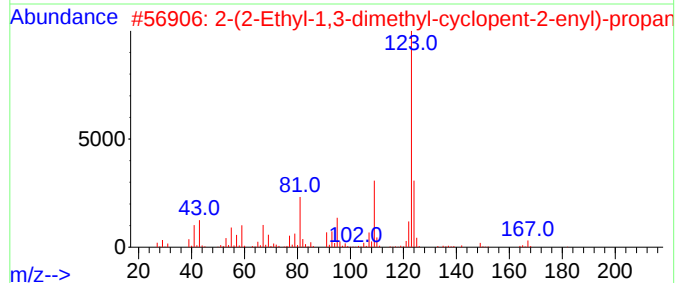
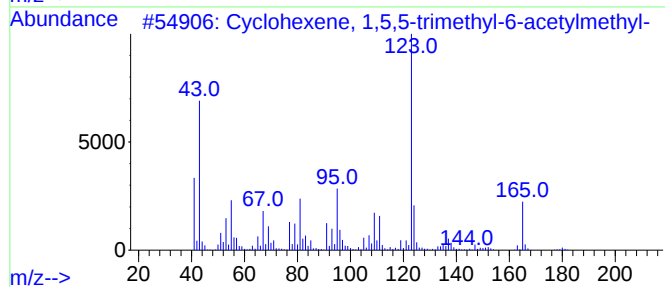
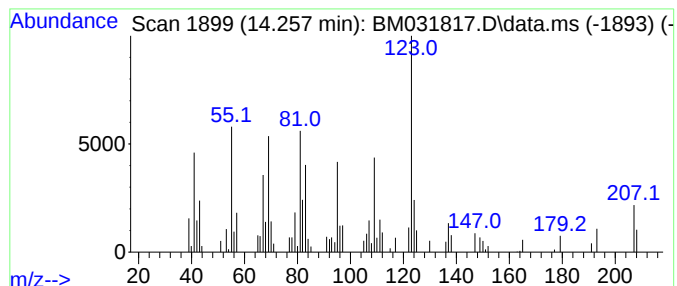
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown14.257 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.257	2.54 ng	66408	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 1,5,5-trimethyl-6-a...	180	C12H200	211563-96-1	43
2			2-(2-Ethyl-1,3-dimethyl-cyclopent...	182	C12H220	1000186-82-4	43
3			photocitral A	152	C10H160	055253-28-6	43
4			Benzamide, 3-fluoro-N-(2-phenyle...	299	C19H22FNO	1000451-30-2	38
5			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
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ClientSampleId :
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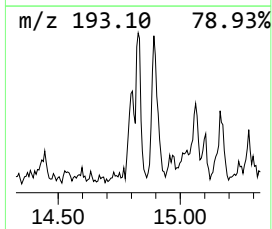
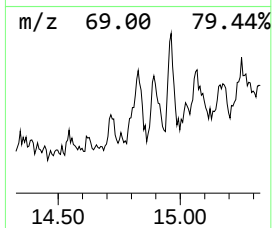
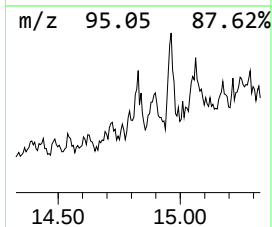
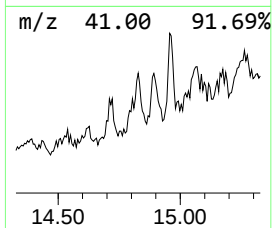
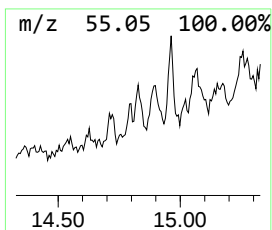
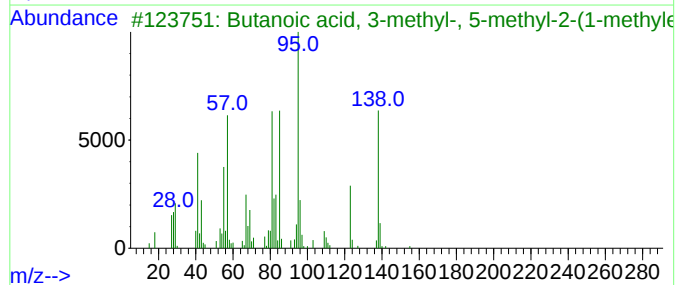
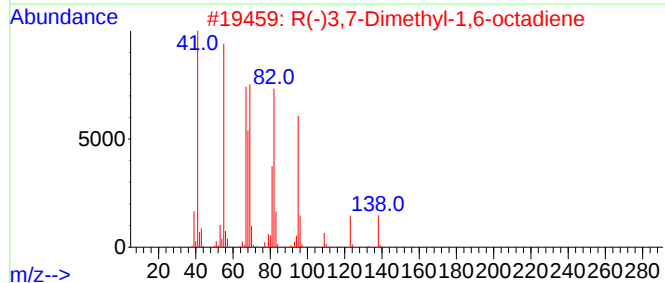
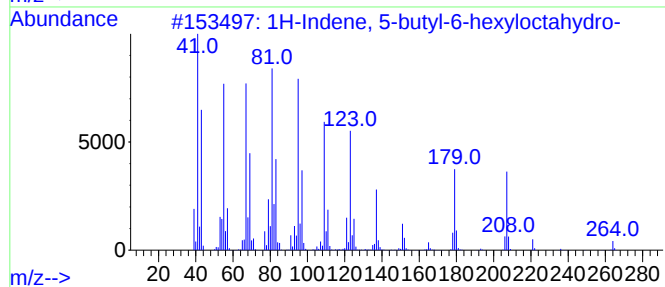
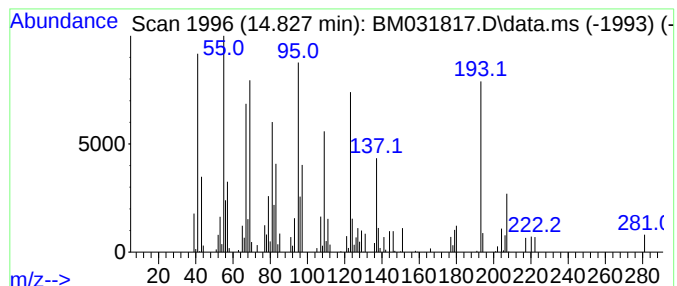
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown14.827 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.827	2.95 ng	77014	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 5-butyl-6-hexyloctahy...	264	C19H36	055044-36-5	43
2			R(-)3,7-Dimethyl-1,6-octadiene	138	C10H18	010281-56-8	27
3			Butanoic acid, 3-methyl-, 5-meth...	240	C15H28O2	000089-47-4	27
4			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	27
5			4-Decyne	138	C10H18	002384-86-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
Operator : CG/JU
Sample : M3438-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

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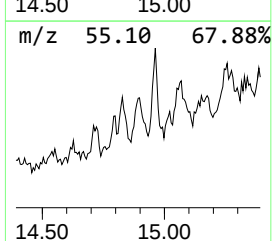
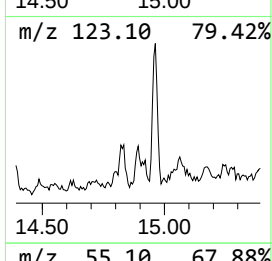
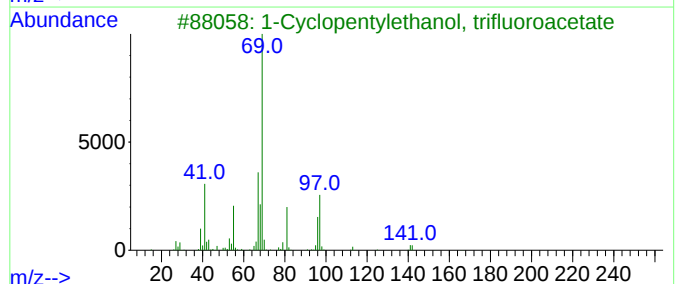
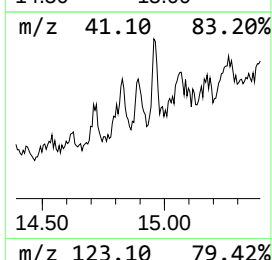
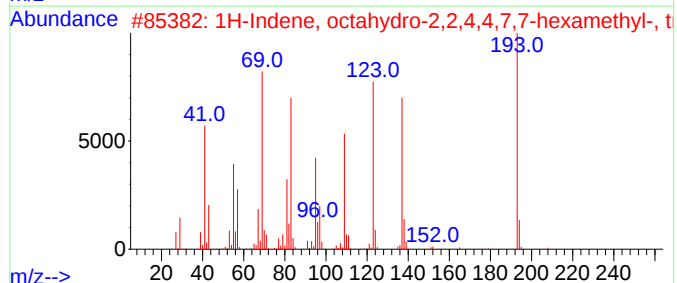
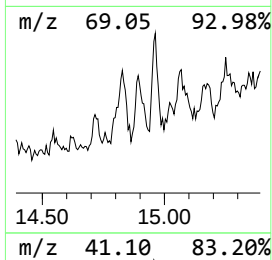
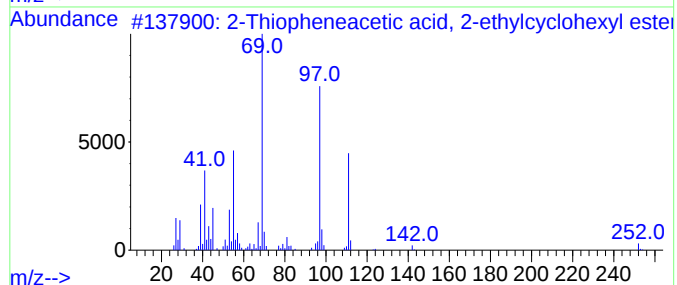
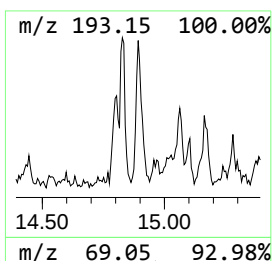
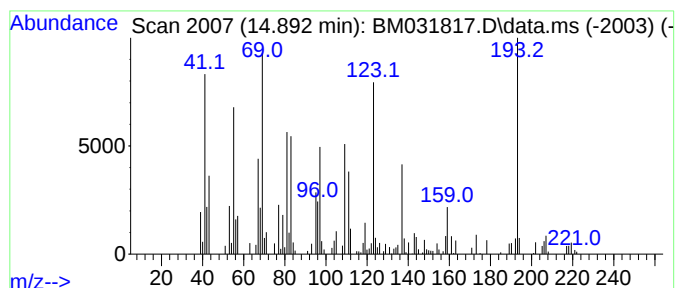
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.892 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.892	3.29 ng	86091	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Thiopheneacetic acid, 2-ethylc...	252	C14H20O2S	1000278-96-1	22
2			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	22
3			1-Cyclopentylethanol, trifluoroa...	210	C9H13F3O2	1000364-11-9	14
4			Benzamide, 4-fluoro-N-allyl-N-et...	207	C12H14FNO	1000446-42-7	11
5			2,4,6-Trimethyl-2-(4-methyl-pent...	206	C14H22O	1000193-58-6	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
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Misc :
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ClientSampleId :
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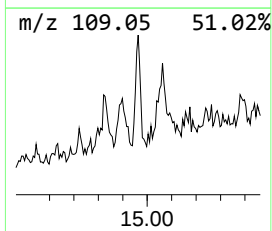
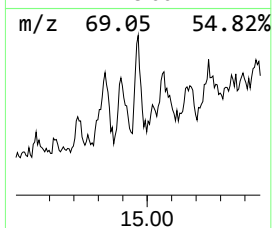
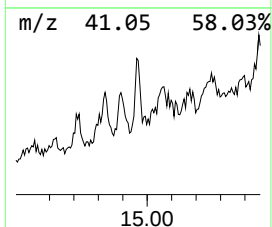
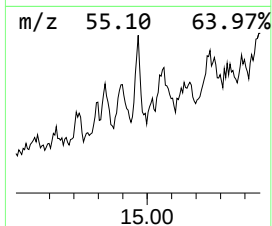
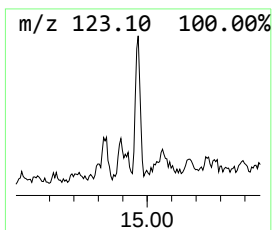
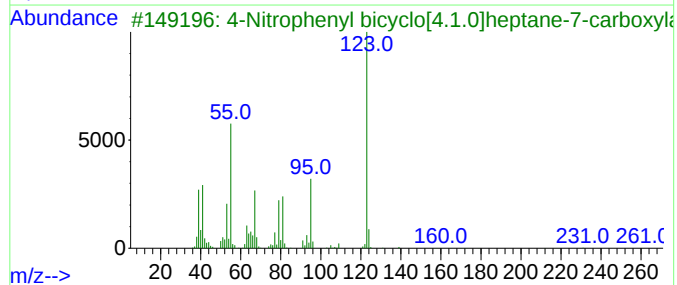
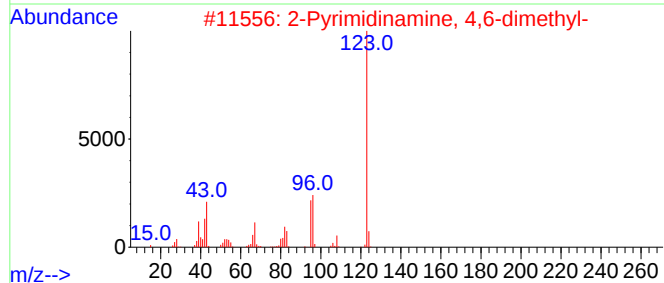
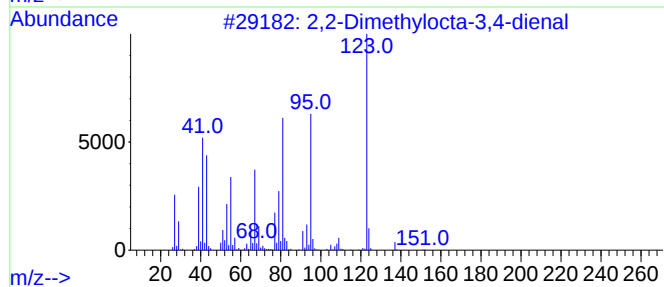
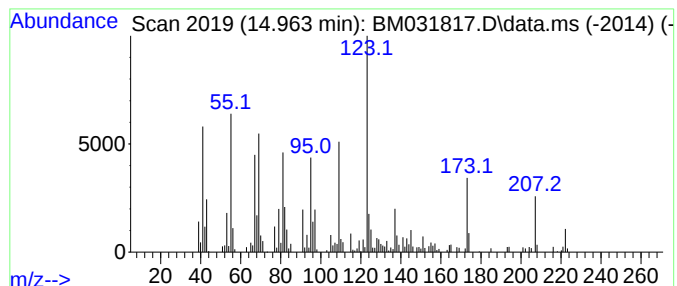
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown14.963 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.963	5.73 ng	149629	Acenaphthene-d10	14.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-Dimethylocta-3,4-dienal	152	C10H16O	000590-71-6	38
2			2-Pyrimidinamine, 4,6-dimethyl-	123	C6H9N3	000767-15-7	35
3			4-Nitrophenyl bicyclo[4.1.0]hept...	261	C14H15NO4	328938-65-4	35
4			Pyridine-3-carboxylic acid, 1-[...	290	C15H18N2O4	1010310-27-9	35
5			1,1-Dimethyl-4-methylenecyclohexane	124	C9H16	006007-96-1	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
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Sample : M3438-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

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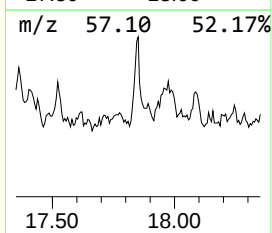
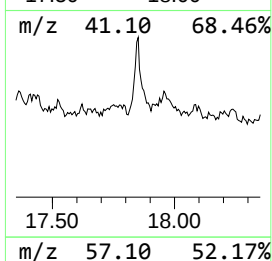
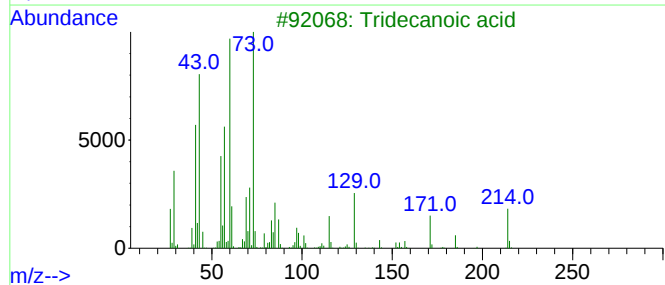
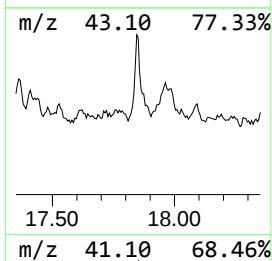
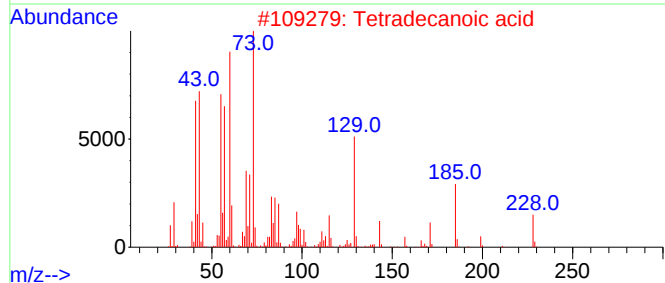
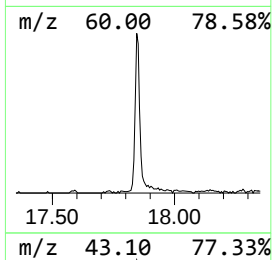
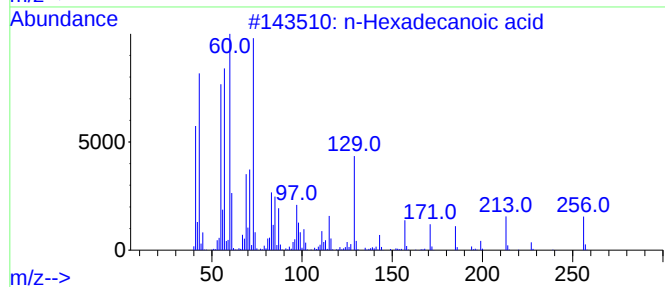
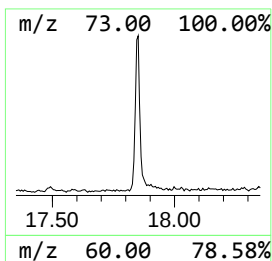
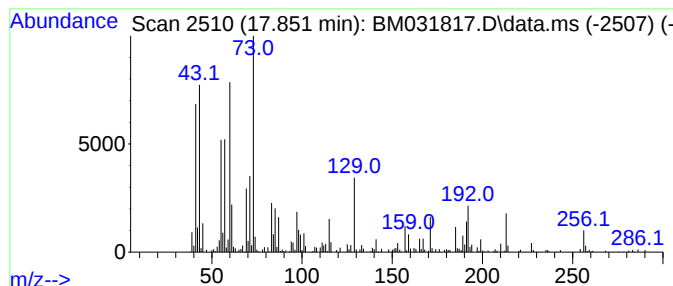
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.851	11.65 ng	452623	Phenanthrene-d10	16.933

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	95
3			Tridecanoic acid	214	C13H26O2	000638-53-9	93
4			Octadecanoic acid	284	C18H36O2	000057-11-4	76
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
Operator : CG/JU
Sample : M3438-03
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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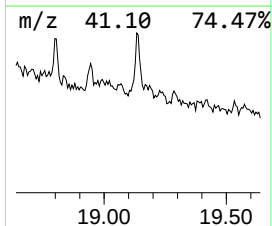
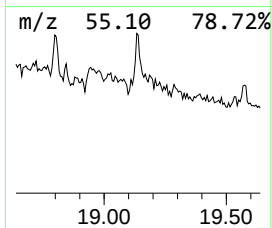
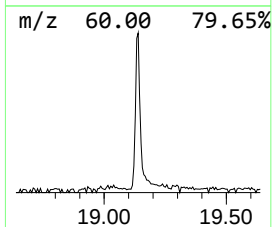
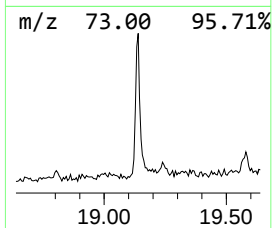
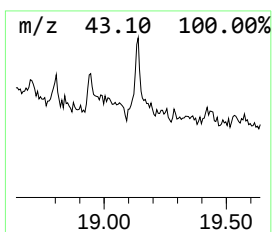
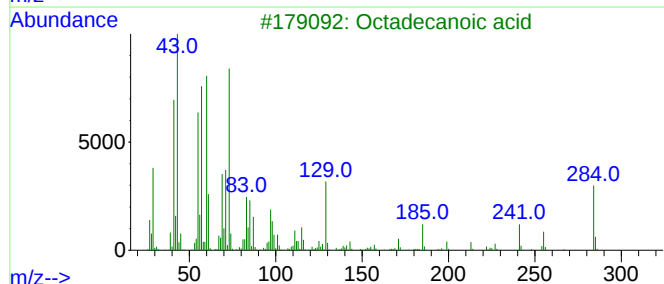
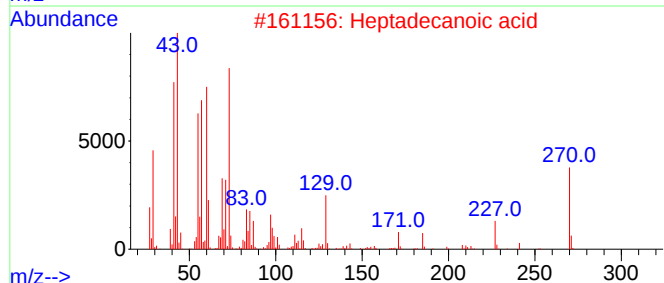
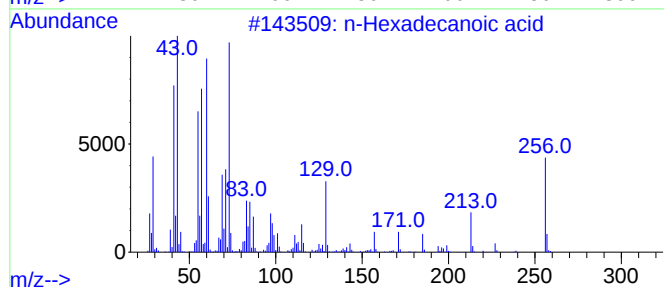
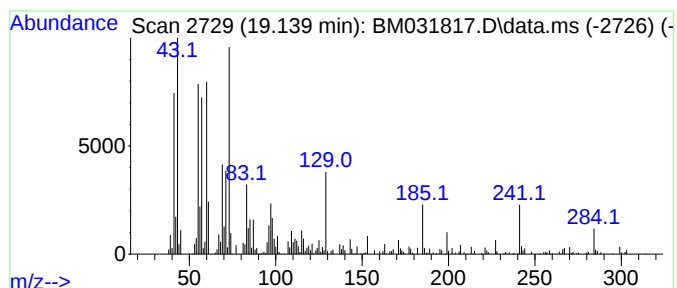
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TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.139	2.76 ng	132379	Chrysene-d12	21.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	89
2	Heptadecanoic acid			270	C17H34O2	000506-12-7	89
3	Octadecanoic acid			284	C18H36O2	000057-11-4	80
4	Tetradecanoic acid			228	C14H28O2	000544-63-8	72
5	n-Decanoic acid			172	C10H20O2	000334-48-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
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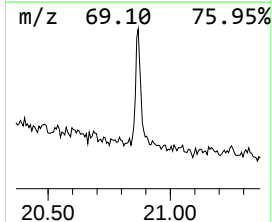
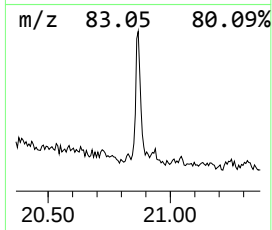
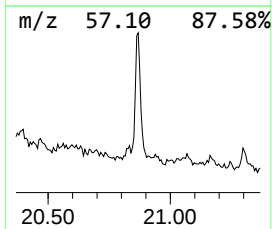
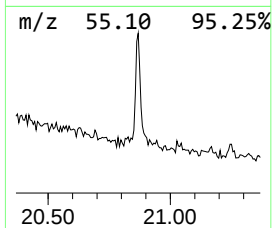
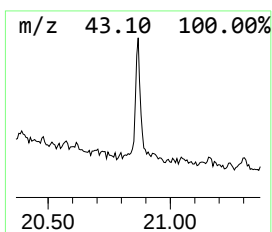
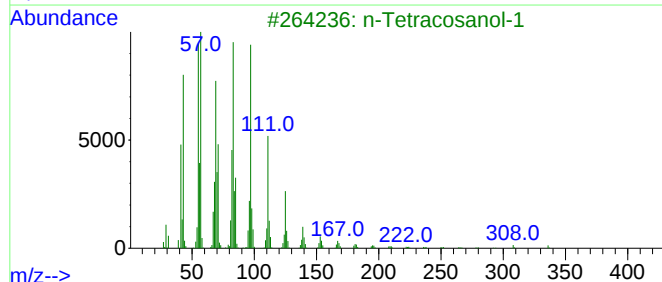
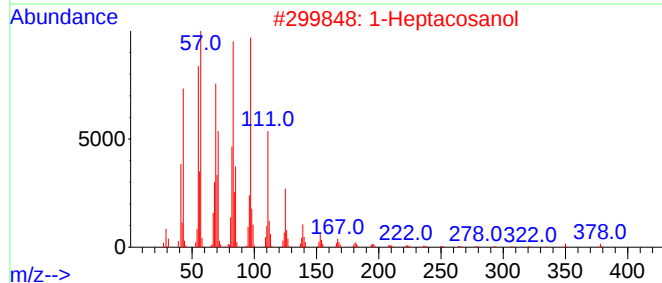
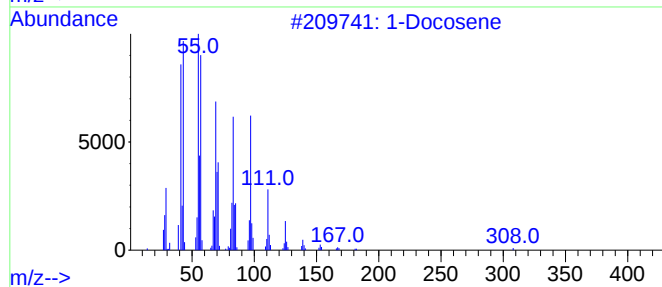
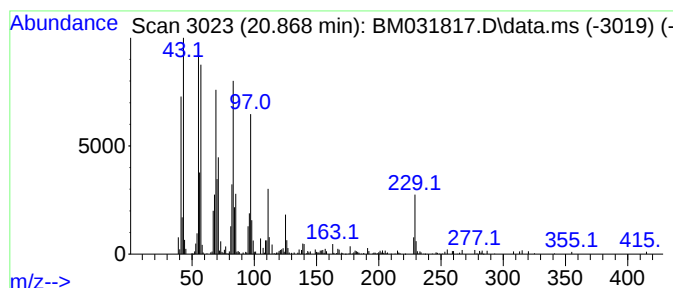
Instrument :
BNA_M
ClientSampleId :
CHRT-26855

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.868	4.17 ng	200266	Chrysene-d12	21.127	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	96
2	1-Heptacosanol	396	C27H56O	002004-39-9	93
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Octacosanol	410	C28H58O	000557-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
Operator : CG/JU
Sample : M3438-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CHRT-26855

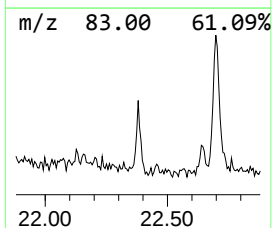
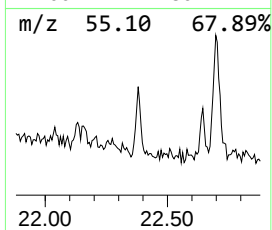
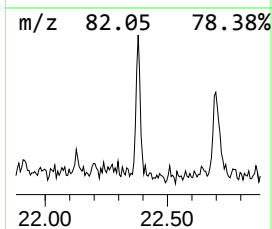
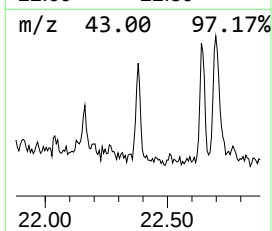
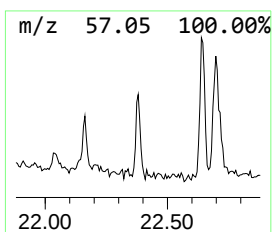
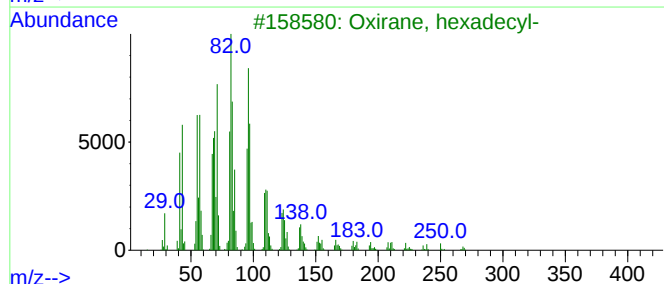
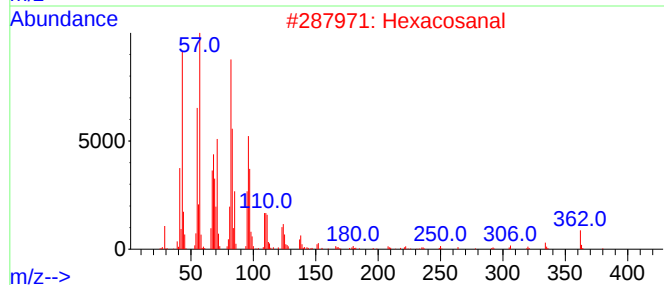
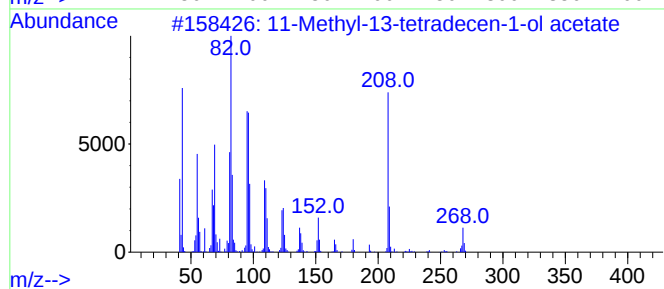
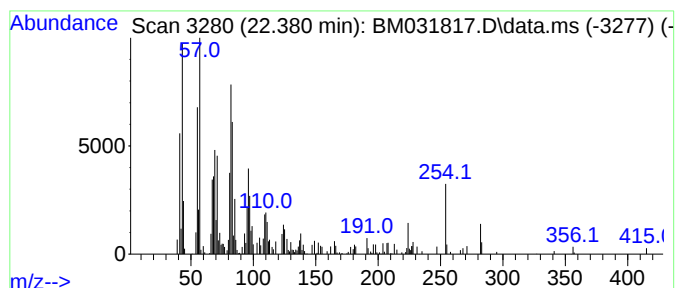
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 11-Methyl-13-tetradecen-1-o... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.380	2.74 ng	97122	Perylene-d12	23.274

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11-Methyl-13-tetradecen-1-ol ace...	268	C17H32O2	1000131-33-5	83
2			Hexacosanal	380	C26H52O	026627-85-0	81
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	78
4			Tetracosanal	352	C24H48O	057866-08-7	76
5			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
Operator : CG/JU
Sample : M3438-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CHRT-26855

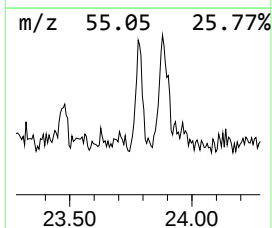
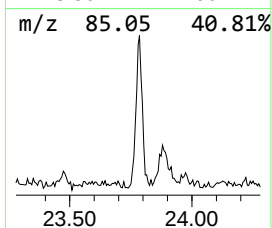
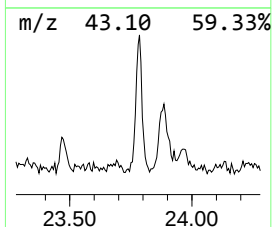
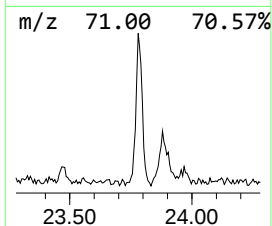
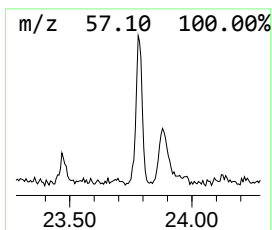
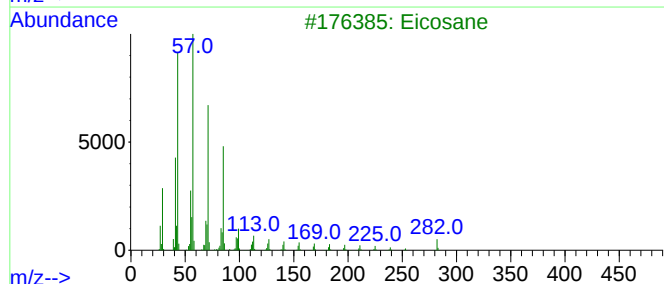
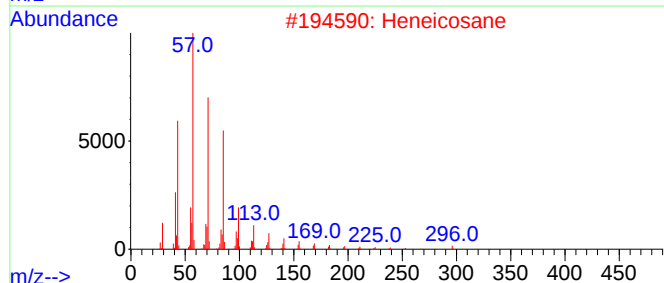
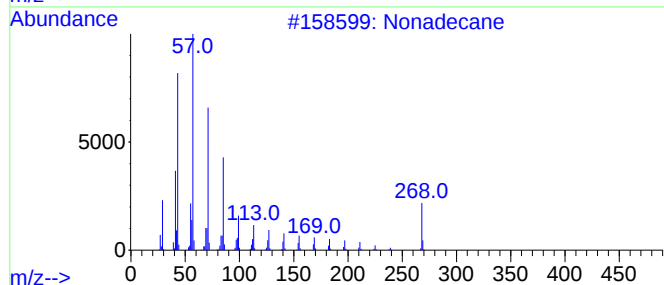
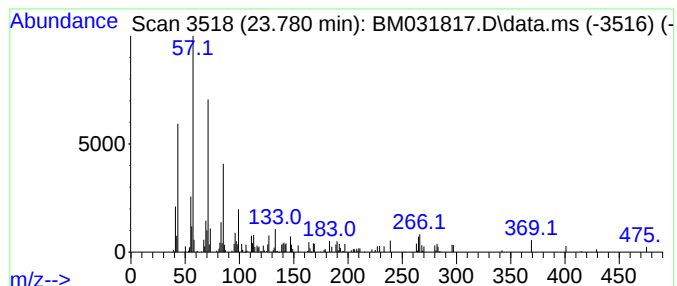
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.780	2.80 ng	99100	Perylene-d12	23.274

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	92
2			Heneicosane	296	C21H44	000629-94-7	90
3			Eicosane	282	C20H42	000112-95-8	86
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5			Hentriacontane	437	C31H64	000630-04-6	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
Data File : BM031817.D
Acq On : 18 Aug 2021 22:08
Operator : CG/JU
Sample : M3438-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CHRT-26855

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.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
(+)-2-Bornanone	9.722	3.0	ng	52412	2	10.298	355769	20.0
Ethanol, 2-(2-b...	10.098	12.8	ng	227028	2	10.298	355769	20.0
unknown13.580	13.580	2.4	ng	61444	3	14.175	522663	20.0
unknown13.904	13.904	2.2	ng	56432	3	14.175	522663	20.0
Decahydro-1,1,4...	13.992	2.9	ng	74362	3	14.175	522663	20.0
unknown14.257	14.257	2.5	ng	66408	3	14.175	522663	20.0
unknown14.827	14.827	3.0	ng	77014	3	14.175	522663	20.0
unknown14.892	14.892	3.3	ng	86091	3	14.175	522663	20.0
unknown14.963	14.963	5.7	ng	149629	3	14.175	522663	20.0
n-Hexadecanoic ...	17.851	11.7	ng	452623	4	16.933	777291	20.0
Heptadecanoic acid	19.139	2.8	ng	132379	5	21.127	959584	20.0
1-Docosene	20.868	4.2	ng	200266	5	21.127	959584	20.0
11-Methyl-13-te...	22.380	2.7	ng	97122	6	23.274	708793	20.0
Nonadecane	23.780	2.8	ng	99100	6	23.274	708793	20.0