

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
 Data File : BM038780.D
 Acq On : 17 Feb 2023 18:36
 Operator : CG/JU
 Sample : 01570-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 AU-6-021523

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-BM020123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM038780.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.969	292	303	310	rBV	257068	370922	13.45%	2.017%
2	5.440	375	383	388	rBV	778776	1087051	39.41%	5.911%
3	5.487	388	391	403	rVB	30699	65393	2.37%	0.356%
4	5.857	446	454	461	rBV3	24947	49316	1.79%	0.268%
5	6.840	617	621	626	rVB	24210	34700	1.26%	0.189%
6	6.951	635	640	644	rBV	25010	34574	1.25%	0.188%
7	7.010	644	650	653	rVV4	18120	33892	1.23%	0.184%
8	7.063	653	659	670	rVB	761682	1190022	43.14%	6.471%
9	7.522	732	737	743	rVB	46654	74847	2.71%	0.407%
10	7.881	793	798	805	rVB	287055	427811	15.51%	2.326%
11	8.422	883	890	899	rVB5	23505	56596	2.05%	0.308%
12	8.522	899	907	913	rBV5	23688	47326	1.72%	0.257%
13	8.986	980	986	991	rVB	29293	47816	1.73%	0.260%
14	9.063	991	999	1009	rBV	480257	798651	28.95%	4.343%
15	9.510	1069	1075	1080	rBV	20621	30784	1.12%	0.167%
16	10.086	1167	1173	1179	rBV4	19592	39356	1.43%	0.214%
17	10.645	1262	1268	1270	rBV3	19715	32302	1.17%	0.176%
18	10.692	1270	1276	1282	rVV	373498	632753	22.94%	3.441%
19	10.739	1282	1284	1290	rVB2	32435	50512	1.83%	0.275%
20	11.880	1469	1478	1481	rBV7	15353	32793	1.19%	0.178%
21	12.098	1509	1515	1520	rBV3	30768	44325	1.61%	0.241%
22	12.357	1555	1559	1567	rVB2	20554	33009	1.20%	0.179%
23	13.157	1689	1695	1702	rVB	1602618	2179569	79.02%	11.852%
24	13.339	1715	1726	1734	rVB	45340	80743	2.93%	0.439%
25	13.998	1834	1838	1841	rVB5	26854	34776	1.26%	0.189%
26	14.045	1841	1846	1850	rVV6	18899	34127	1.24%	0.186%
27	14.116	1850	1858	1864	rVB9	18573	49591	1.80%	0.270%
28	14.415	1904	1909	1913	rVB	62969	76191	2.76%	0.414%
29	14.533	1923	1929	1935	rVB	570367	830333	30.10%	4.515%
30	14.927	1991	1996	2005	rVB9	16841	46740	1.69%	0.254%
31	15.362	2066	2070	2075	rVB	70062	92049	3.34%	0.501%
32	15.768	2132	2139	2145	rBV2	19007	43845	1.59%	0.238%
33	15.898	2156	2161	2167	rVB	379379	490241	17.77%	2.666%
34	16.027	2178	2183	2189	rVB	1186968	1632169	59.17%	8.876%
35	16.227	2213	2217	2225	rBV	71153	129923	4.71%	0.707%
36	16.457	2252	2256	2263	rBV3	23078	35530	1.29%	0.193%

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37	16.562	2271	2274	2281	rBV8	14423	29416	1.07%	0.160%
38	17.021	2348	2352	2356	rBV	62192	70910	2.57%	0.386%
39	17.068	2356	2360	2364	rVV	21106	30625	1.11%	0.167%
40	17.286	2391	2397	2401	rBV	657782	910997	33.03%	4.954%
41	17.327	2401	2404	2411	rVB	70501	92595	3.36%	0.504%
42	17.762	2474	2478	2482	rBV	57102	65641	2.38%	0.357%
43	17.939	2505	2508	2512	rVB	47300	57701	2.09%	0.314%
44	18.051	2522	2527	2531	rBV7	20899	41141	1.49%	0.224%
45	18.174	2543	2548	2552	rBV	300217	419328	15.20%	2.280%
46	18.356	2574	2579	2582	rBV5	24162	43760	1.59%	0.238%
47	18.456	2592	2596	2600	rBV	52202	71881	2.61%	0.391%
48	19.086	2699	2703	2706	rBV2	183284	212316	7.70%	1.155%
49	19.115	2706	2708	2715	rVB4	84660	99511	3.61%	0.541%
50	19.339	2742	2746	2751	rVB	127958	169996	6.16%	0.924%
51	19.386	2751	2754	2759	rVB2	68359	82437	2.99%	0.448%
52	19.445	2761	2764	2775	rBV2	143755	217490	7.88%	1.183%
53	19.668	2799	2802	2805	rBV3	55722	66888	2.42%	0.364%
54	19.703	2805	2808	2816	rVB	119595	173696	6.30%	0.945%
55	19.903	2837	2842	2847	rBV	2353143	2758340	100.00%	15.000%
56	21.162	3053	3056	3065	rVB2	195739	276386	10.02%	1.503%
57	21.468	3103	3108	3112	rBV	636520	850381	30.83%	4.624%
58	23.856	3509	3514	3520	rBV2	362704	677474	24.56%	3.684%

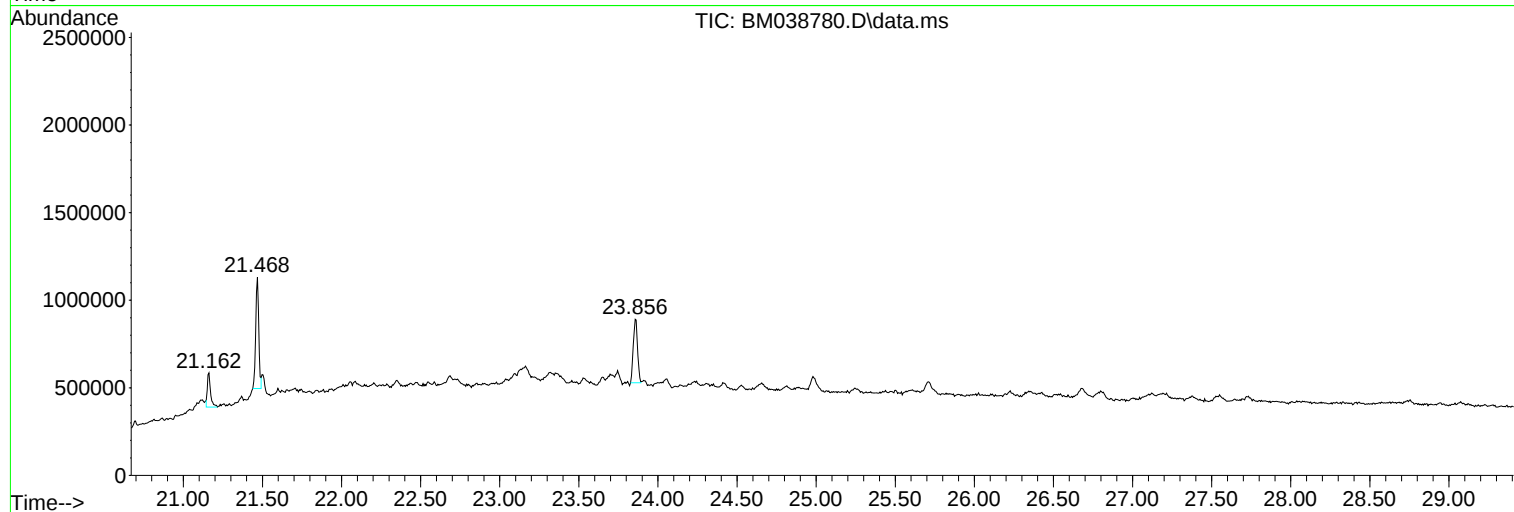
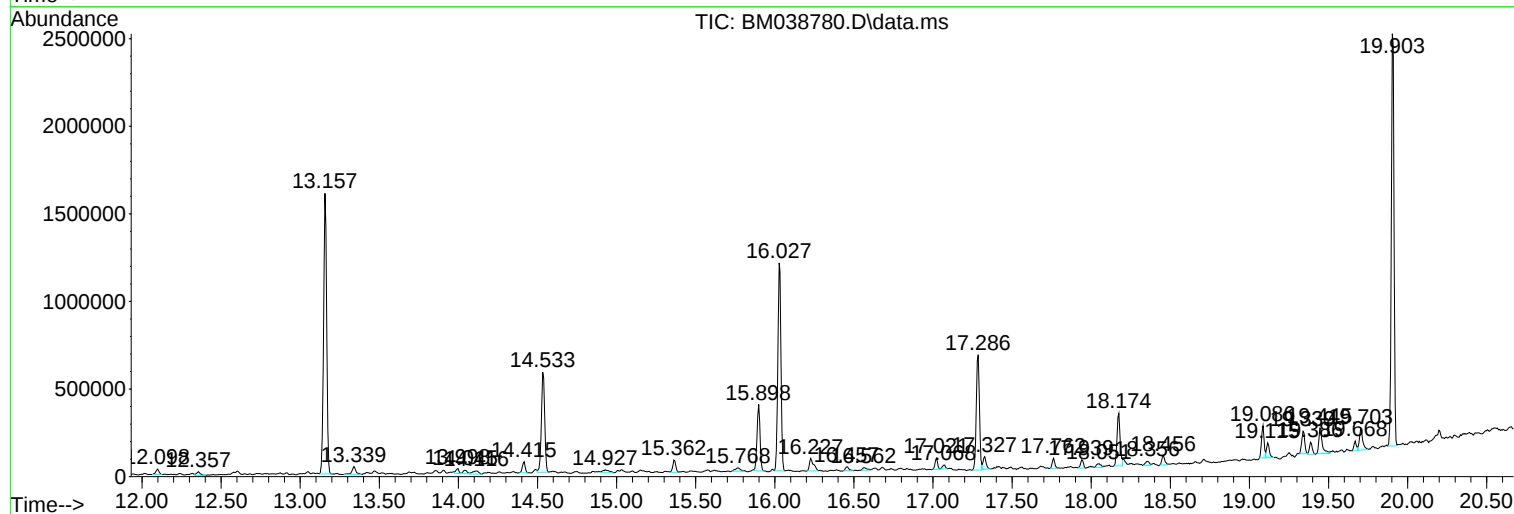
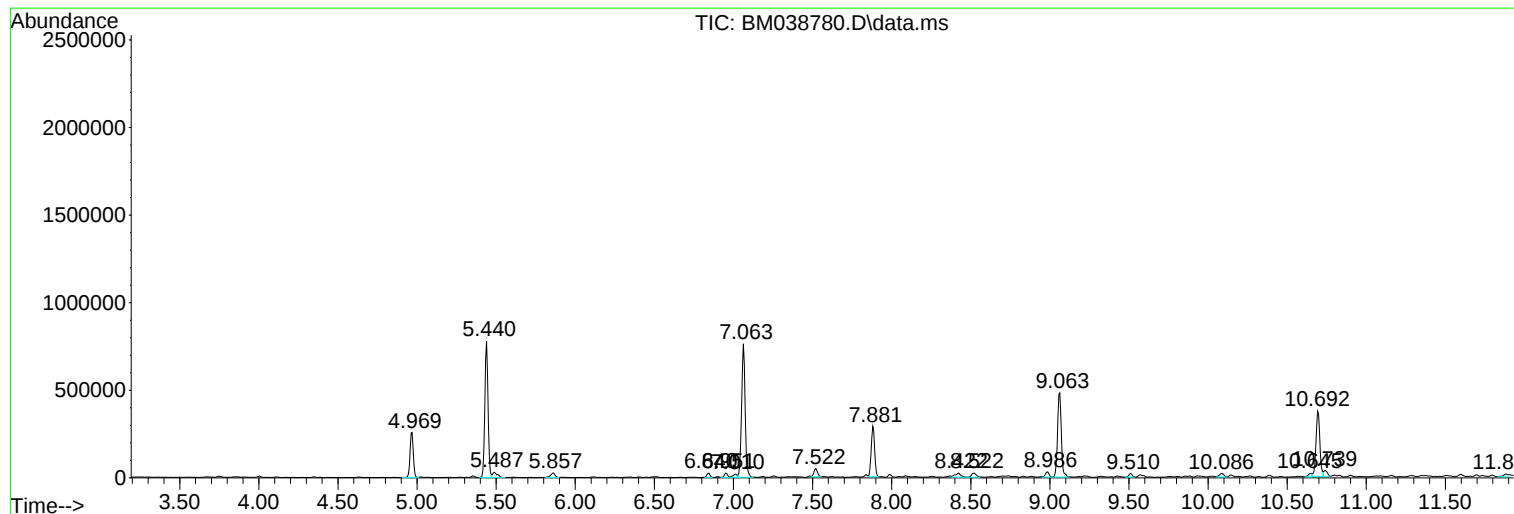
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020123.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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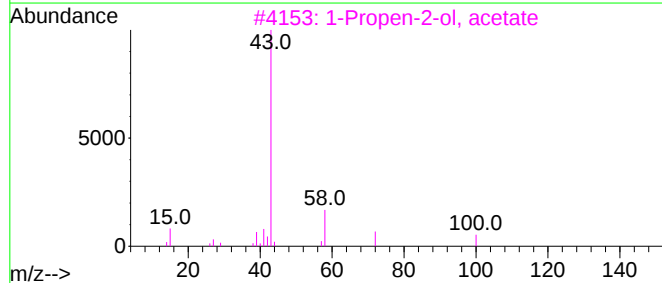
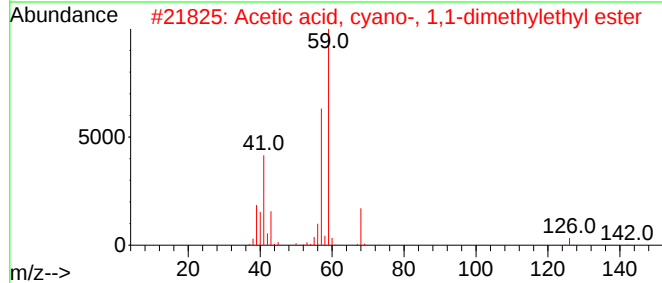
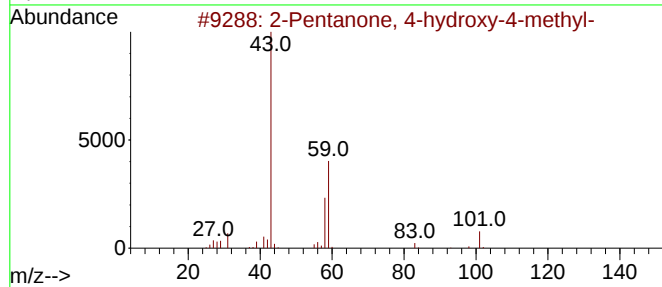
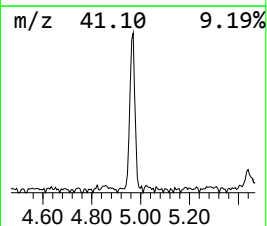
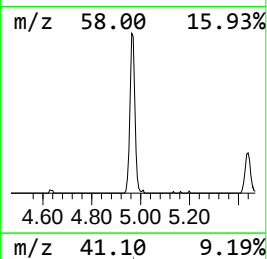
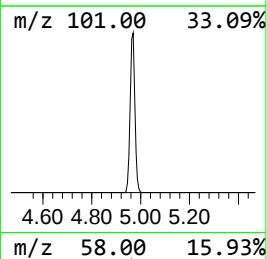
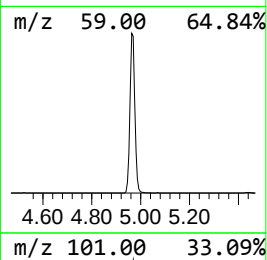
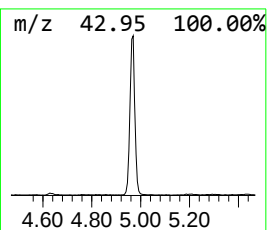
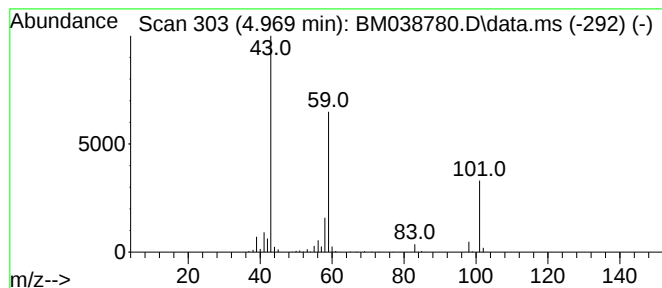
Instrument :
BNA_M
ClientSampleId :
AU-6-021523

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.969	17.34 ng	370922	1,4-Dichlorobenzene-d4			7.881
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	59
2	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
3	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	10
4	Propane, 1-ethoxy-2-methyl-		102	C6H14O	000627-02-1	9
5	Morpholine, 4-methyl-		101	C5H11NO	000109-02-4	9



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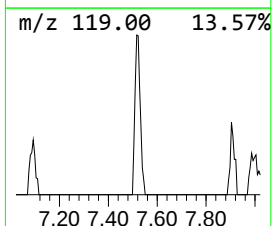
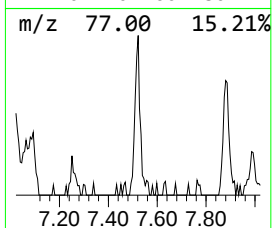
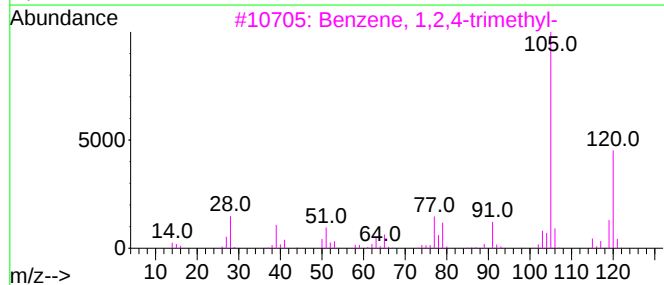
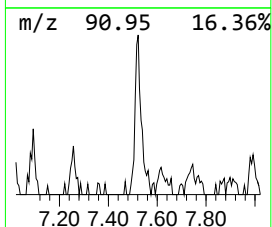
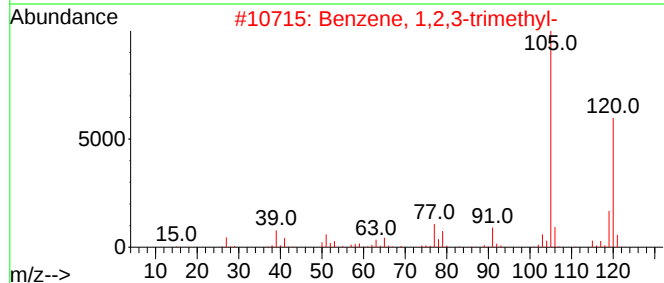
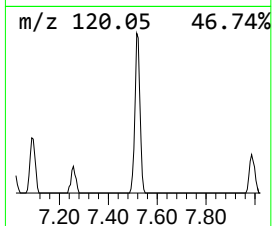
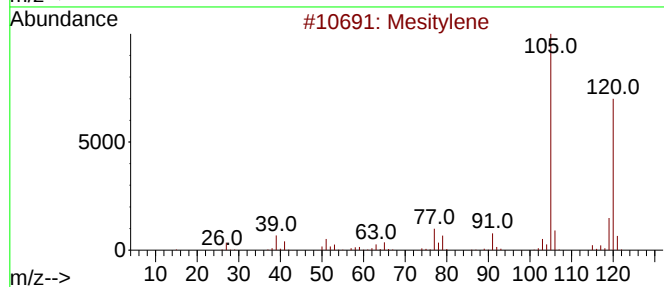
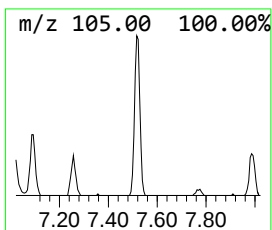
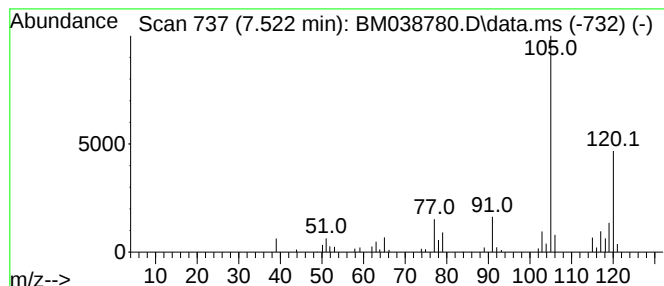
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.522	3.50 ng	74847	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	90
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87
4			1,3-Cyclopentadiene, 5-(1-methyl...	120	C9H12	003141-02-4	83
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	81



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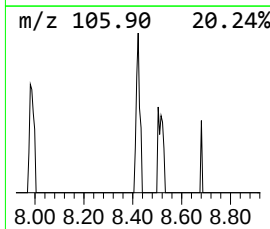
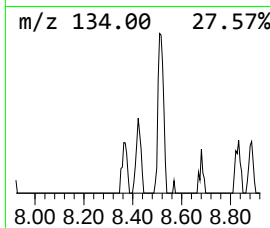
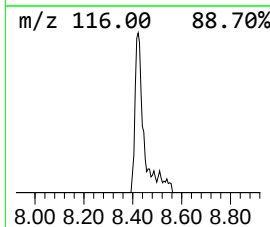
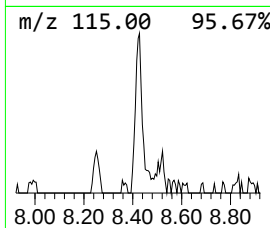
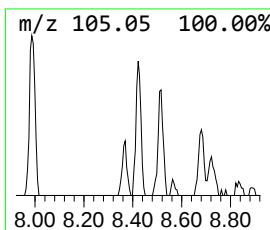
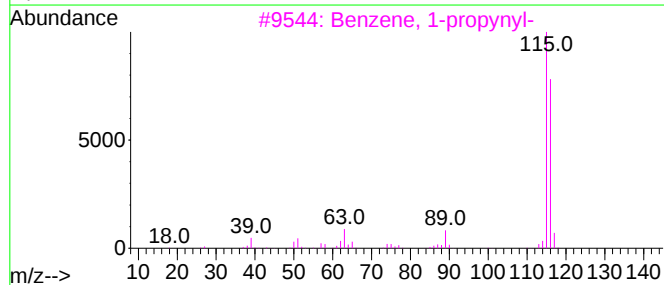
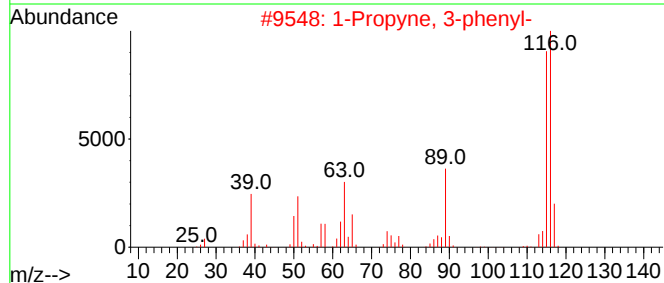
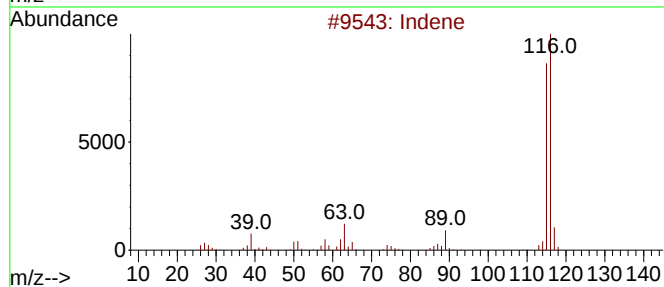
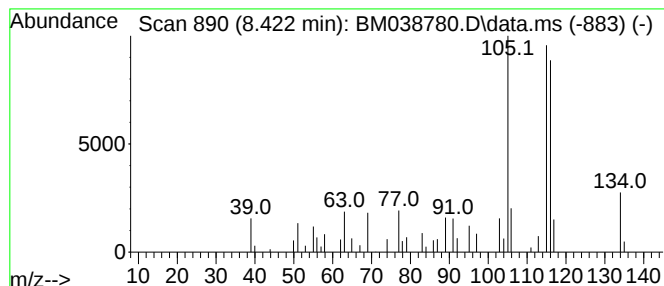
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.422	2.65 ng	56596	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	64
2			1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	58
3			Benzene, 1-propynyl-	116	C9H8	000673-32-5	47
4			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	47
5			Benzonitrile, 3,5-dimethyl-	131	C9H9N	022445-42-7	35



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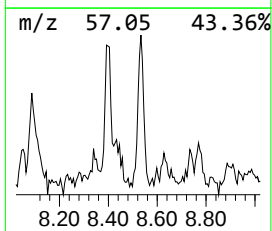
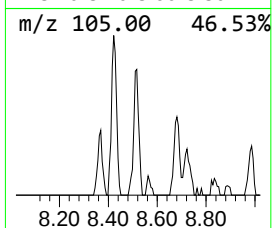
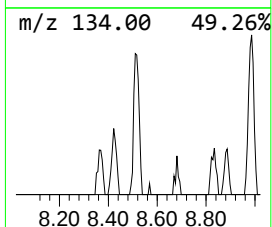
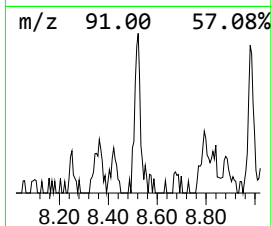
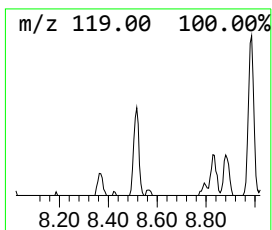
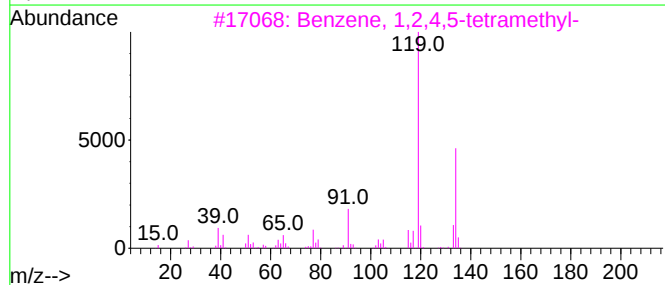
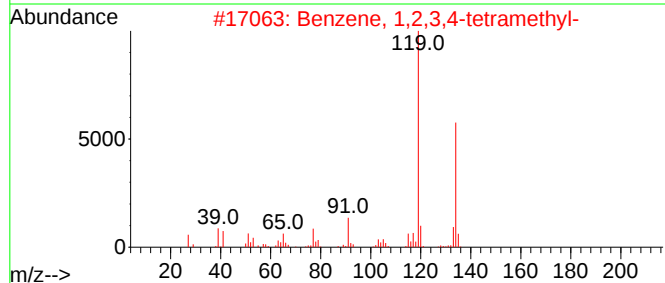
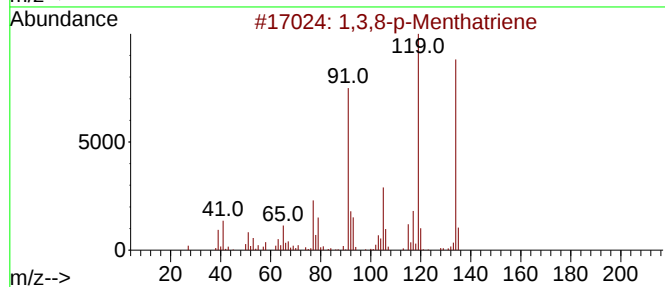
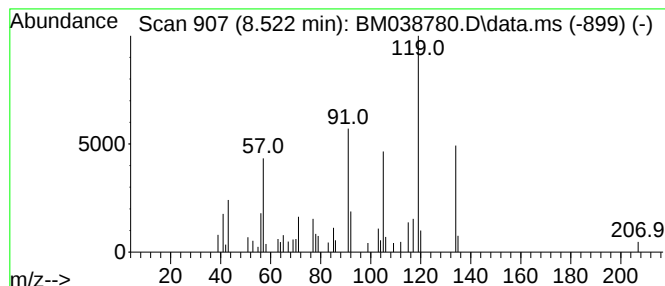
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,3,8-p-Menthatriene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.522	2.21 ng	47326	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	90
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	87
4			2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	87
5			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	87



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AU-6-021523

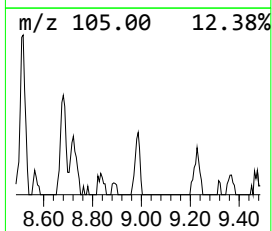
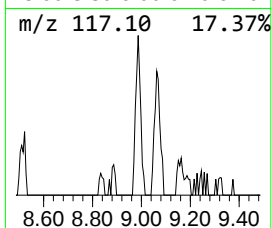
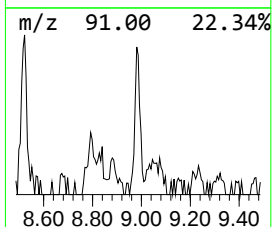
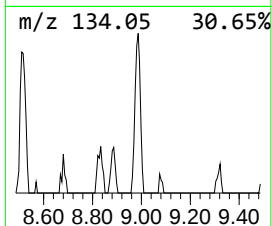
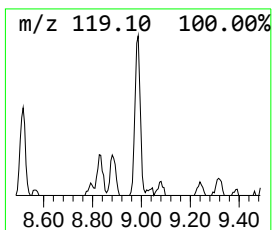
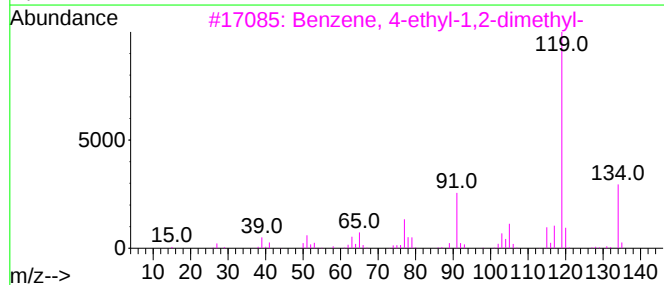
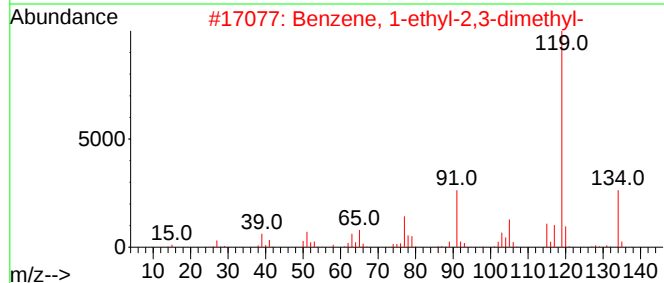
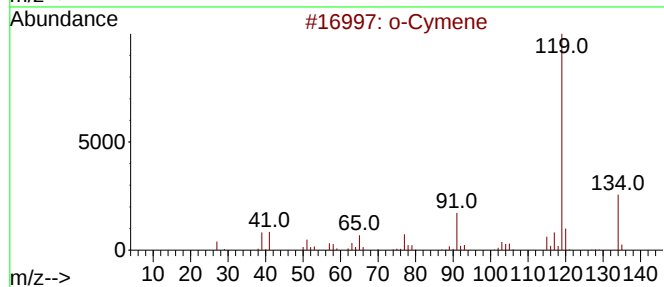
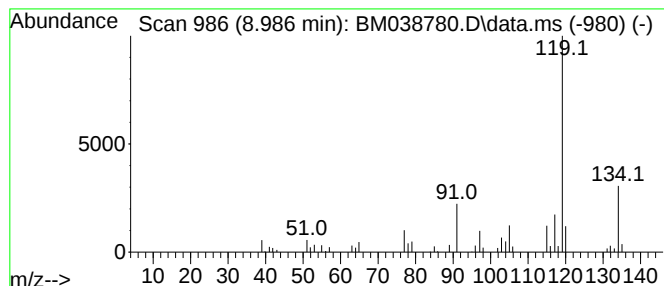
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 o-Cymene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.986	2.24 ng	47816	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	94
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
Operator : CG/JU
Sample : 01570-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AU-6-021523

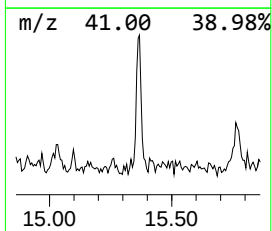
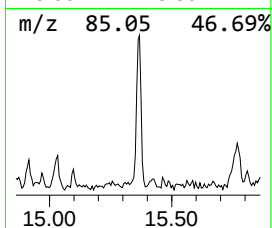
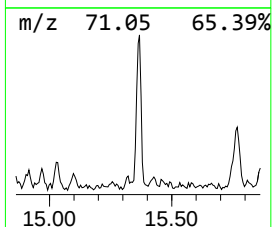
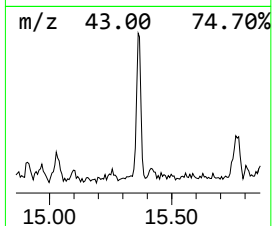
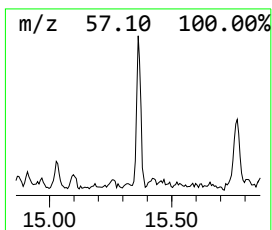
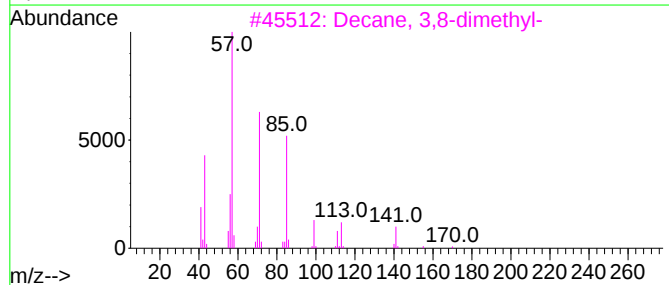
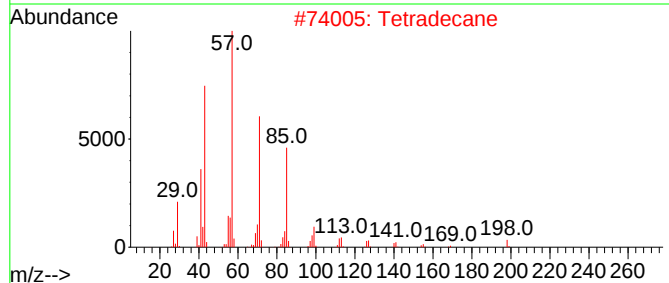
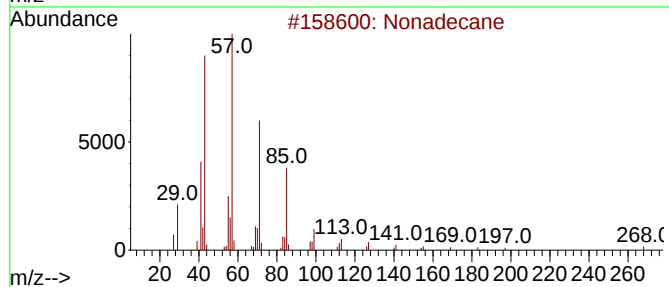
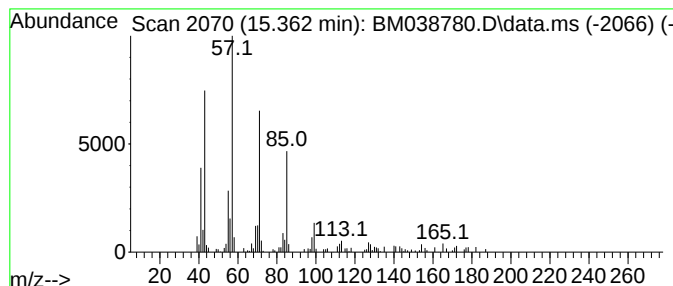
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.363	2.22 ng	92049	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	83
2			Tetradecane	198	C14H30	000629-59-4	83
3			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	81
4			Octacosane	394	C28H58	000630-02-4	80
5			Pentadecane	212	C15H32	000629-62-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
Operator : CG/JU
Sample : 01570-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AU-6-021523

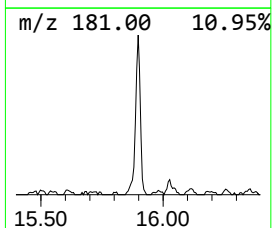
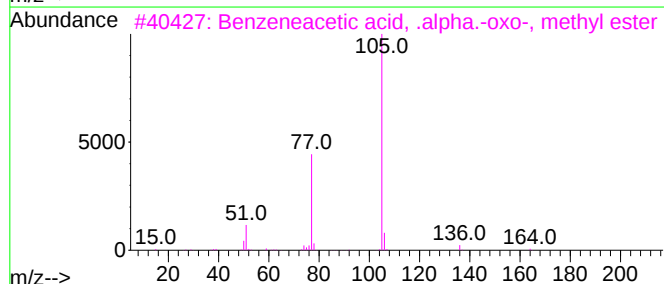
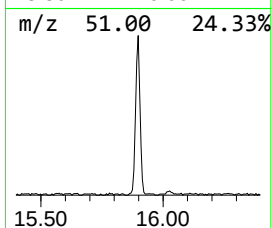
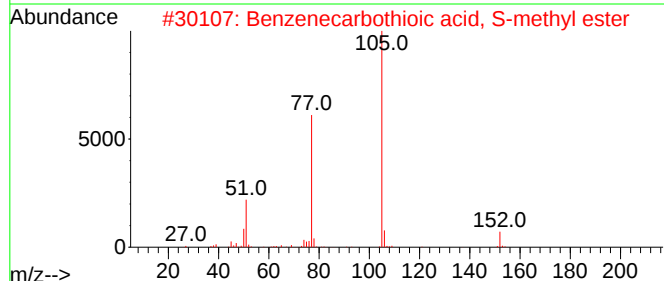
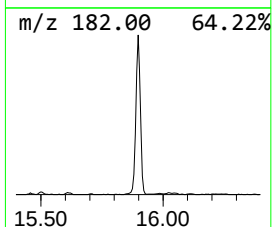
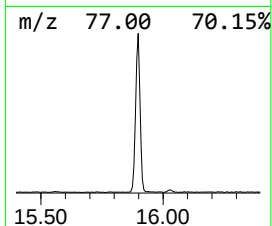
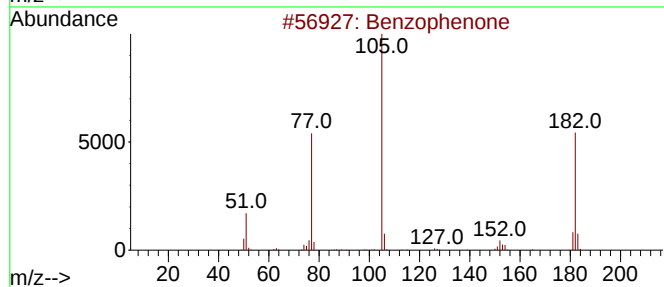
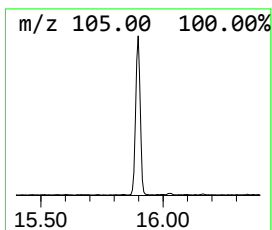
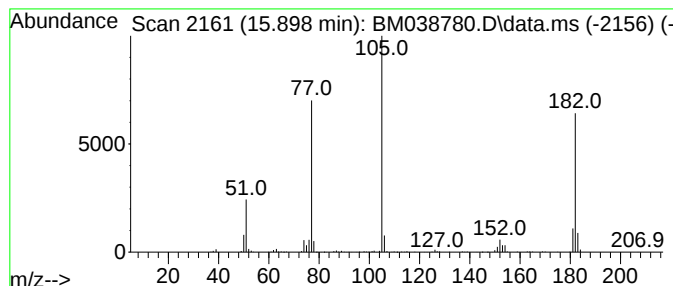
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.898	11.81 ng	490241	Acenaphthene-d10	14.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	49
3			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	47
4			Ethanone, 2,2,2-trifluoro-1-phenyl-	174	C8H5F3O	000434-45-7	47
5			N-Methylbenzamide, N-pentafluoro...	281	C11H8F5NO2	1000446-98-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
Operator : CG/JU
Sample : 01570-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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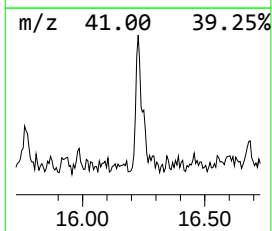
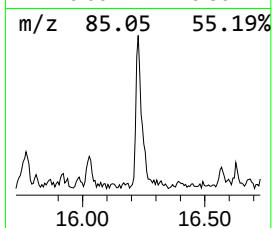
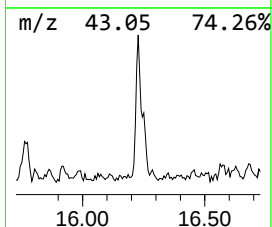
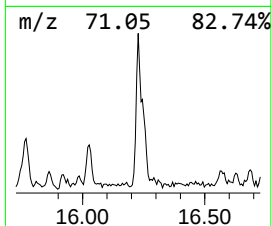
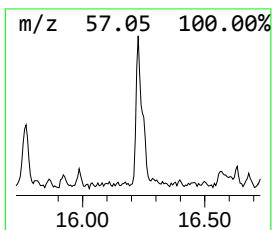
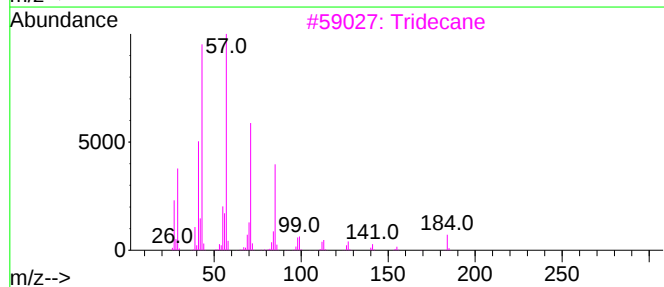
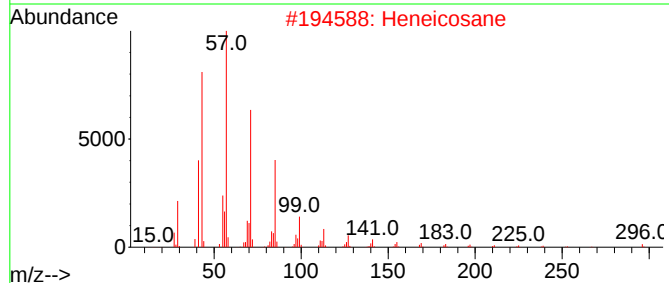
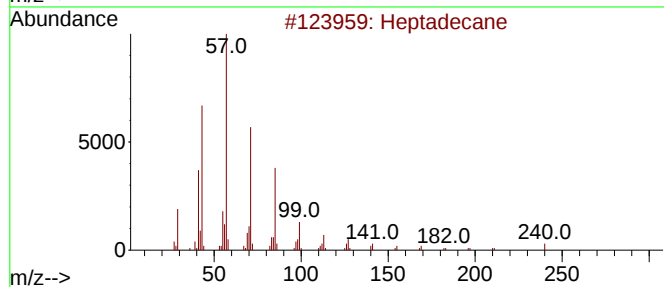
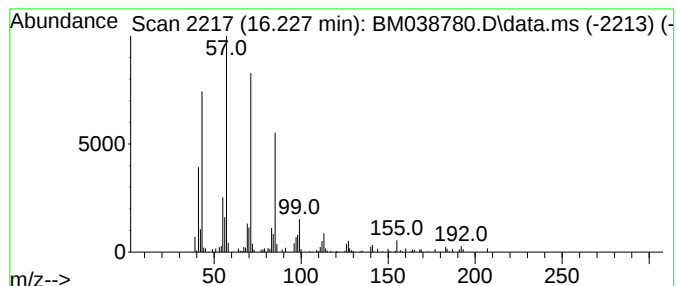
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.227	2.85 ng	129923	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	91
2			Heneicosane	296	C21H44	000629-94-7	90
3			Tridecane	184	C13H28	000629-50-5	90
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	90
5			Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
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ALS Vial : 9 Sample Multiplier: 1

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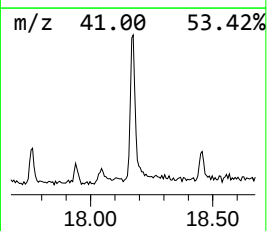
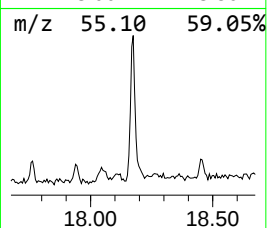
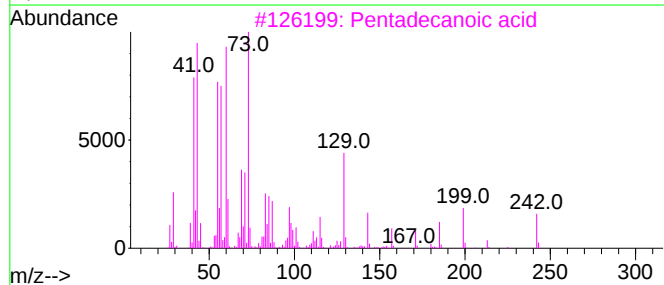
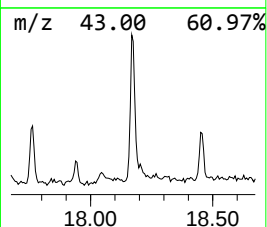
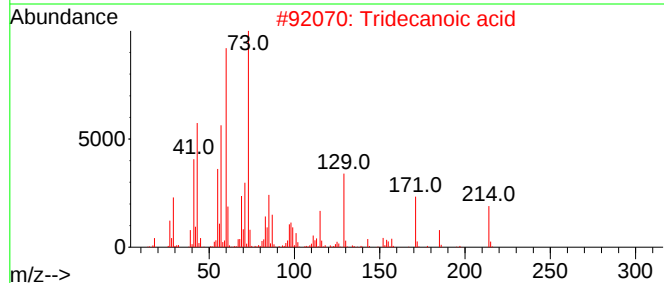
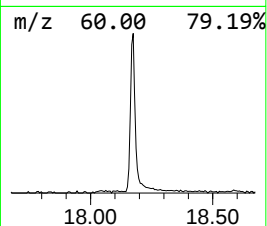
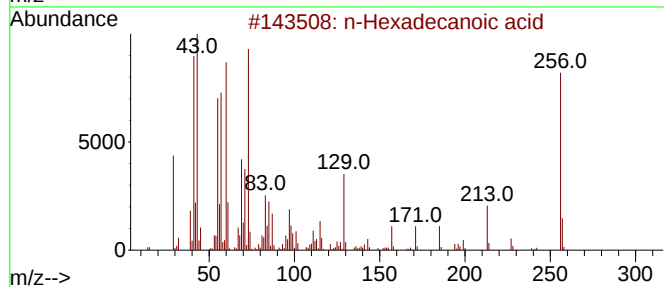
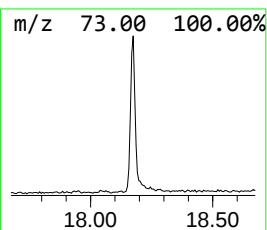
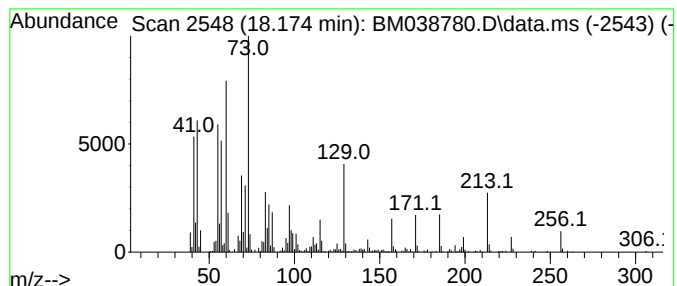
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.174	9.21 ng	419328	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tridecanoic acid	214	C13H26O2	000638-53-9	90
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	68
5			n-Decanoic acid	172	C10H20O2	000334-48-5	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
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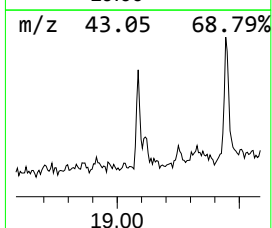
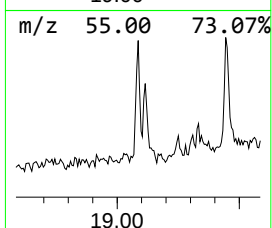
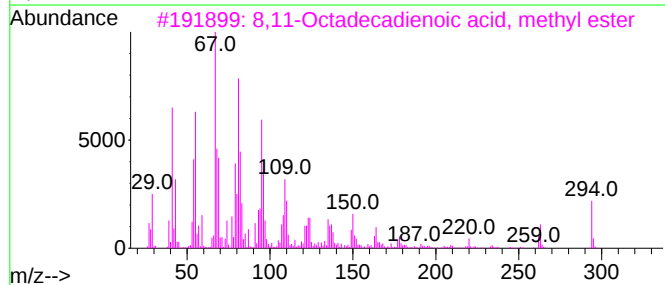
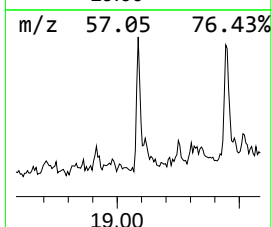
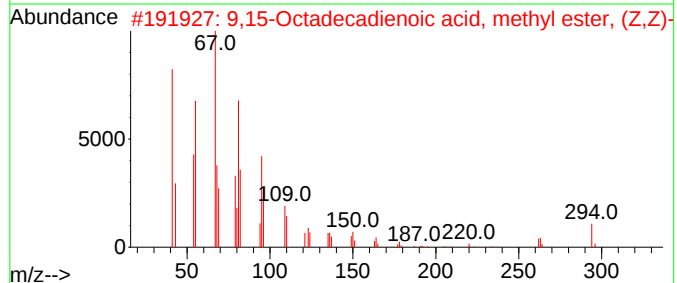
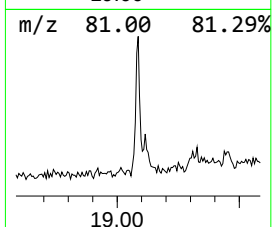
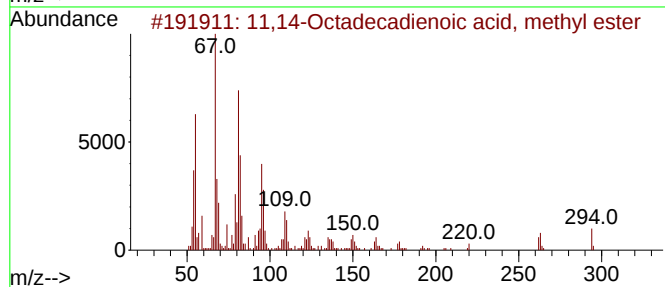
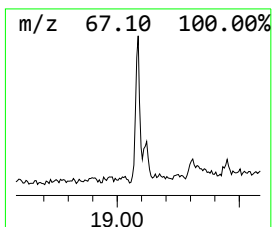
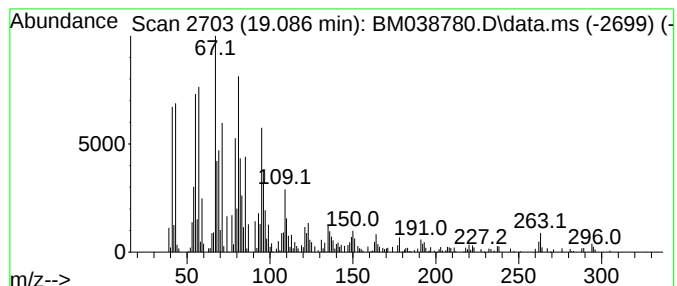
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11,14-Octadecadienoic acid,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.086	4.66 ng	212316	Phenanthrene-d10	17.286	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	97
2	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	96
3	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	96
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	95
5	9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
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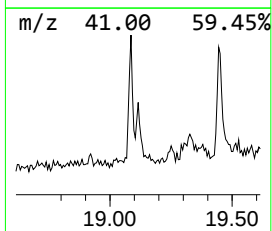
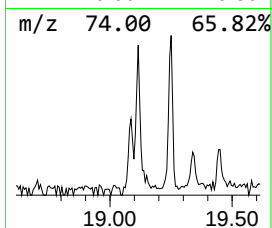
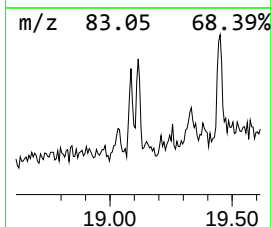
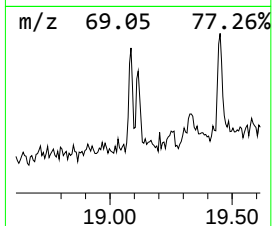
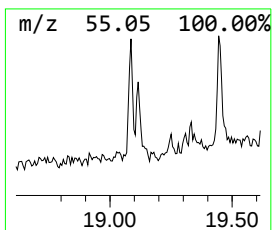
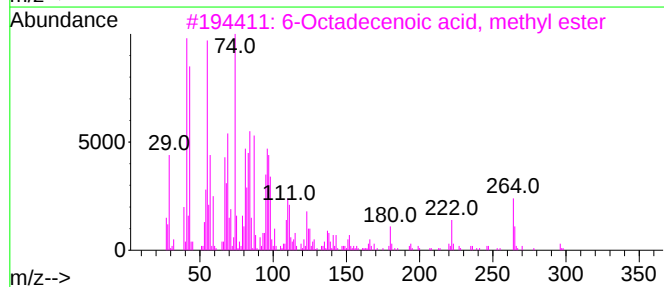
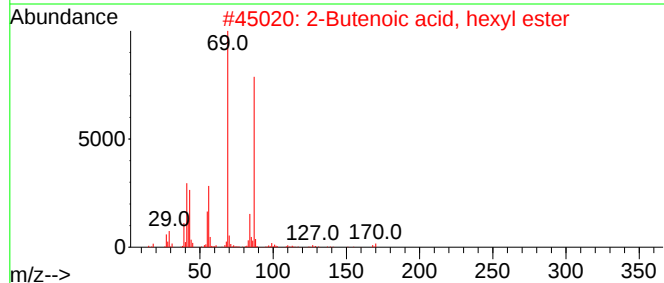
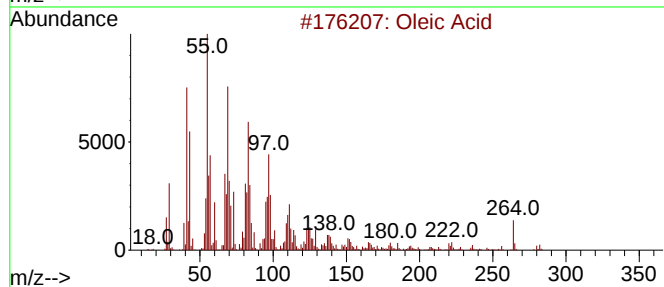
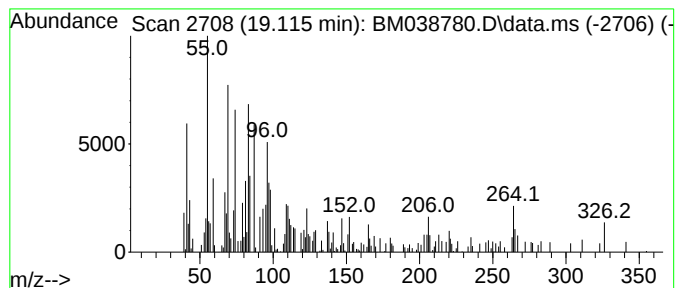
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown19.115 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.115	2.18 ng	99511	Phenanthrene-d10	17.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleic Acid			282	C18H34O2	000112-80-1	41
2	2-Butenoic acid, hexyl ester			170	C10H18O2	019089-92-0	35
3	6-Octadecenoic acid, methyl ester			296	C19H36O2	052355-31-4	35
4	Methyl myristoleate			240	C15H28O2	056219-06-8	30
5	3-Ethyl-4-octanol			158	C10H22O	063126-48-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
Operator : CG/JU
Sample : 01570-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AU-6-021523

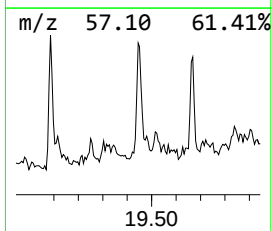
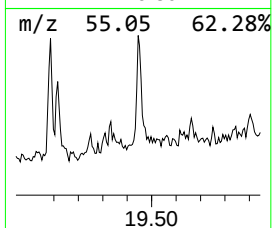
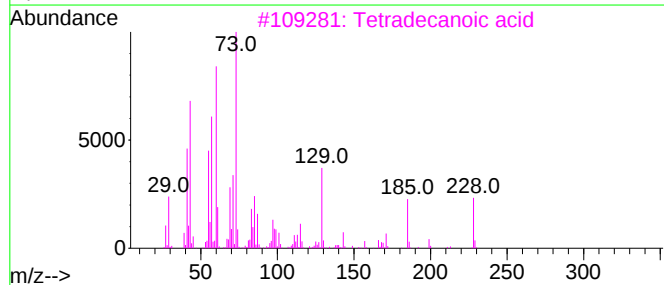
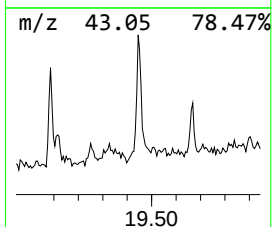
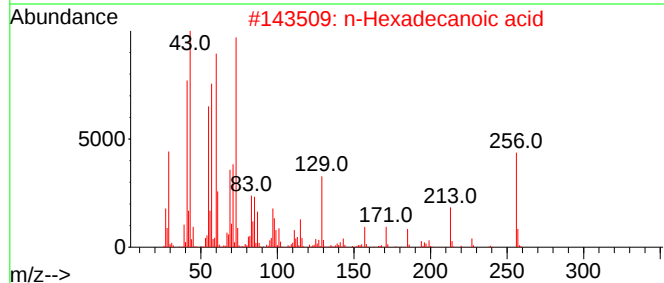
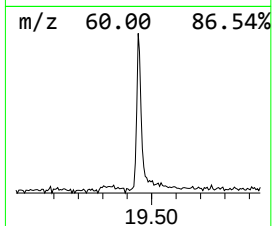
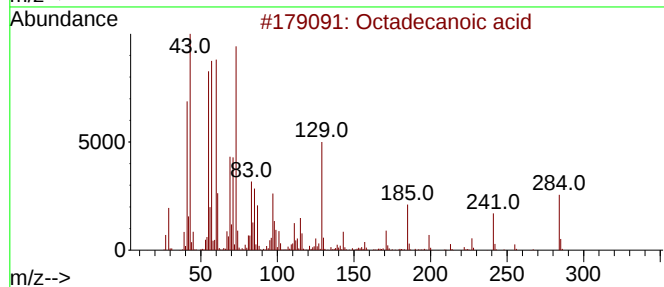
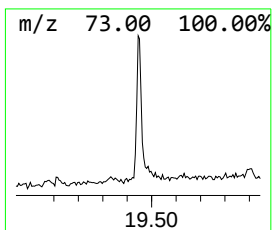
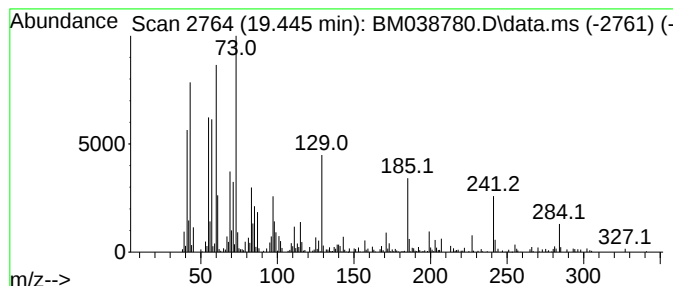
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	5.12 ng	217490	Chrysene-d12	21.468

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	92
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64
5			Nonadecanoic acid	298	C19H38O2	000646-30-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
Operator : CG/JU
Sample : 01570-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AU-6-021523

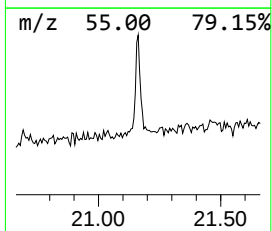
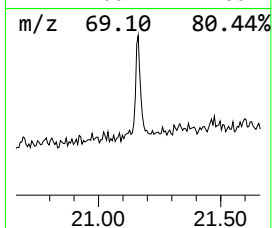
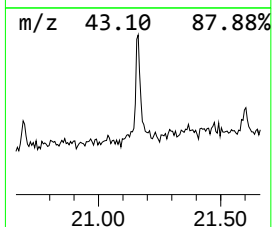
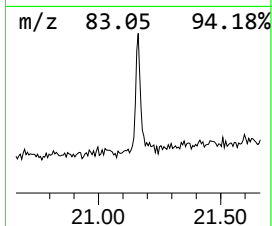
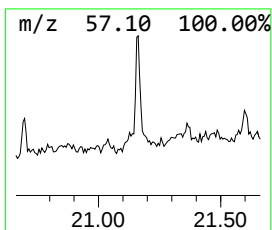
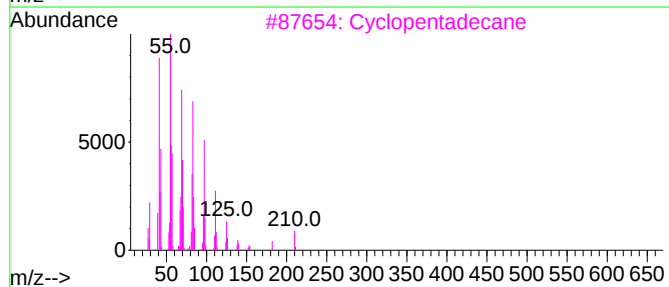
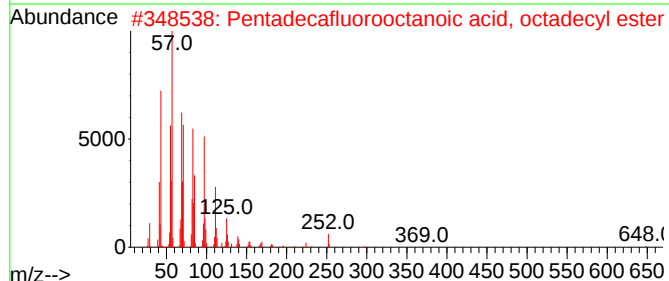
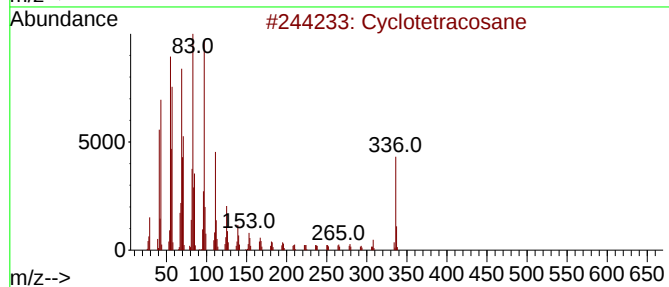
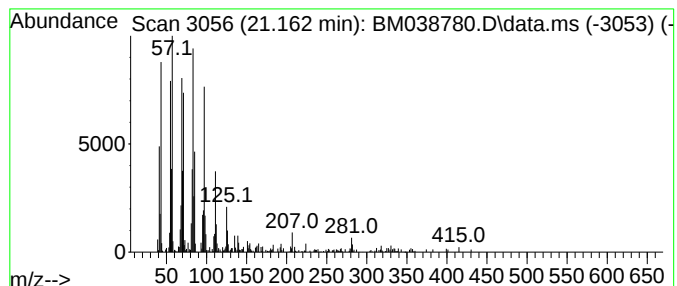
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Cyclotetracosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.162	6.50 ng	276386	Chrysene-d12	21.468

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	98
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3			Cyclopentadecane	210	C15H30	000295-48-7	93
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021723\
Data File : BM038780.D
Acq On : 17 Feb 2023 18:36
Operator : CG/JU
Sample : 01570-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
AU-6-021523

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020123.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-Pentanone, 4-...	4.969	17.3	ng	370922	1	7.881	427811	20.0
Mesitylene	7.522	3.5	ng	74847	1	7.881	427811	20.0
Indene	8.422	2.6	ng	56596	1	7.881	427811	20.0
1,3,8-p-Menthat...	8.522	2.2	ng	47326	1	7.881	427811	20.0
o-Cymene	8.986	2.2	ng	47816	1	7.881	427811	20.0
Nonadecane	15.363	2.2	ng	92049	3	14.533	830333	20.0
Benzophenone	15.898	11.8	ng	490241	3	14.533	830333	20.0
Heptadecane	16.227	2.9	ng	129923	4	17.286	910997	20.0
n-Hexadecanoic ...	18.174	9.2	ng	419328	4	17.286	910997	20.0
11,14-Octadecad...	19.086	4.7	ng	212316	4	17.286	910997	20.0
unknown19.115	19.115	2.2	ng	99511	4	17.286	910997	20.0
Octadecanoic acid	19.445	5.1	ng	217490	5	21.468	850381	20.0
Cyclotetracosane	21.162	6.5	ng	276386	5	21.468	850381	20.0