

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
 Data File : BM026857.D
 Acq On : 23 Jul 2020 18:45
 Operator : JU/CG
 Sample : L3381-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 339-GRID-B3-B4

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BM070820.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.925	366	372	390	rBV	1492258	2130623	58.01%	9.031%
2	6.495	633	639	653	rBV	1408243	2188661	59.59%	9.277%
3	7.278	765	772	780	rBV	406381	607189	16.53%	2.574%
4	8.431	960	968	979	rBV	828101	1380578	37.59%	5.852%
5	10.030	1232	1240	1251	rBV	481457	801562	21.82%	3.398%
6	11.083	1412	1419	1424	rBV	40562	68112	1.85%	0.289%
7	12.483	1652	1657	1662	rBV	54114	77518	2.11%	0.329%
8	12.548	1662	1668	1676	rVB	1803064	2586797	70.43%	10.965%
9	12.783	1704	1708	1715	rVB3	48719	86455	2.35%	0.366%
10	12.866	1717	1722	1728	rVB3	42820	72681	1.98%	0.308%
11	13.354	1801	1805	1811	rVB4	36970	61695	1.68%	0.262%
12	13.424	1811	1817	1821	rBV	246522	365080	9.94%	1.548%
13	13.460	1821	1823	1827	rVB	85158	94223	2.57%	0.399%
14	13.883	1891	1895	1900	rVV	79169	100262	2.73%	0.425%
15	13.948	1900	1906	1915	rVB	691423	952387	25.93%	4.037%
16	14.513	1998	2002	2007	rVB2	50444	66358	1.81%	0.281%
17	14.854	2056	2060	2064	rVB	105833	126651	3.45%	0.537%
18	15.260	2122	2129	2134	rBV	98856	165289	4.50%	0.701%
19	15.460	2157	2163	2171	rVB	1562080	2067685	56.29%	8.765%
20	15.724	2202	2208	2209	rBV	134977	180297	4.91%	0.764%
21	15.742	2209	2211	2220	rVB	192643	257056	7.00%	1.090%
22	16.071	2263	2267	2273	rBV5	18773	38932	1.06%	0.165%
23	16.512	2338	2342	2347	rBV	111992	150221	4.09%	0.637%
24	16.565	2347	2351	2358	rVB	95110	134786	3.67%	0.571%
25	16.712	2371	2376	2381	rBV	774048	1023604	27.87%	4.339%
26	17.254	2464	2468	2472	rVV	101972	110896	3.02%	0.470%
27	17.653	2532	2536	2542	rBV	169915	230157	6.27%	0.976%
28	17.948	2582	2586	2591	rVB	71454	83552	2.27%	0.354%
29	18.401	2658	2663	2669	rBV	154507	207399	5.65%	0.879%
30	18.595	2693	2696	2699	rVB	55420	58688	1.60%	0.249%
31	18.789	2725	2729	2732	rBV	70737	102868	2.80%	0.436%
32	19.153	2788	2791	2794	rVV	59410	66775	1.82%	0.283%
33	19.383	2824	2830	2835	rBV	3103878	3673055	100.00%	15.569%
34	20.689	3048	3052	3062	rBV	552949	732515	19.94%	3.105%

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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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35	20.942	3090	3095	3099	rBV	1007300	1260104	34.31%	5.341%
36	23.000	3441	3445	3457	rVB	788320	1280646	34.87%	5.428%

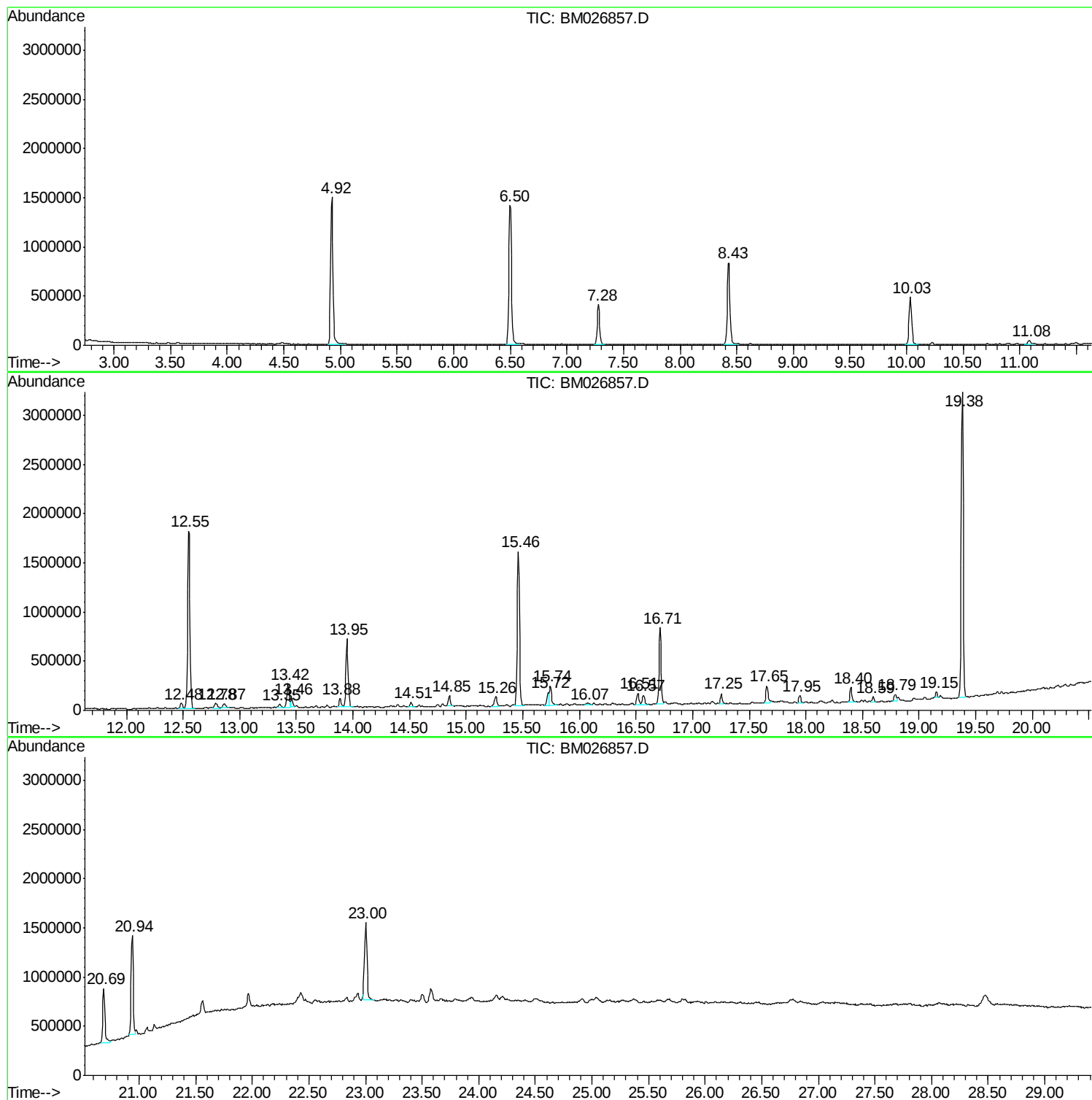
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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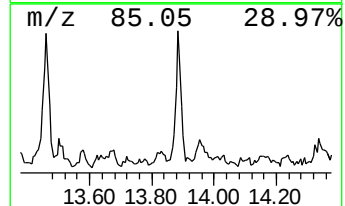
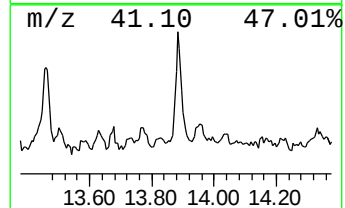
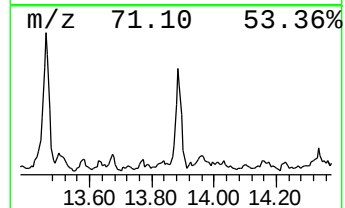
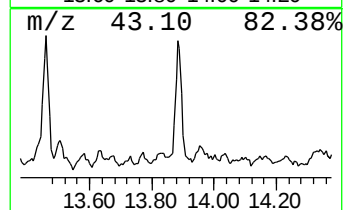
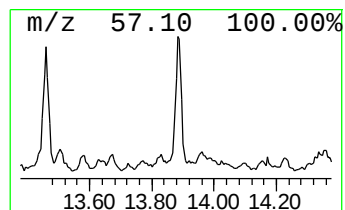
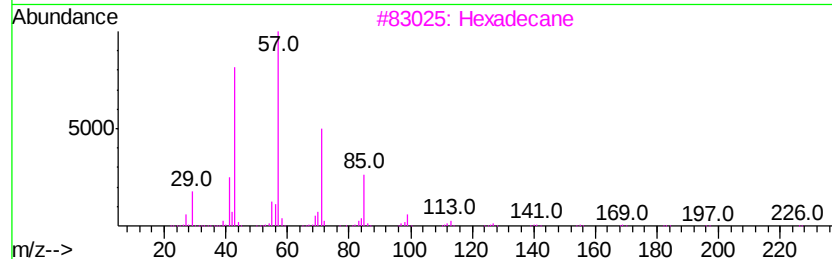
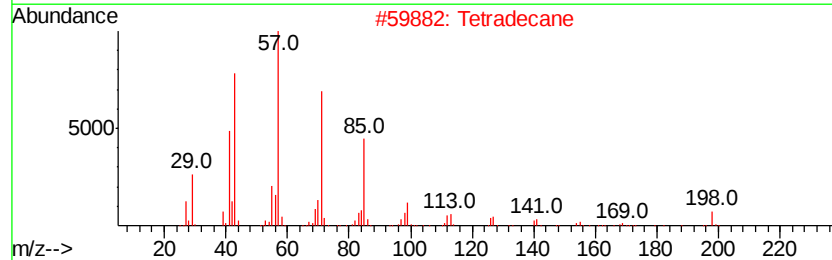
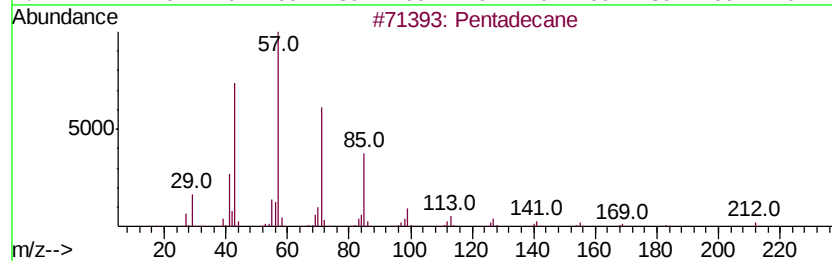
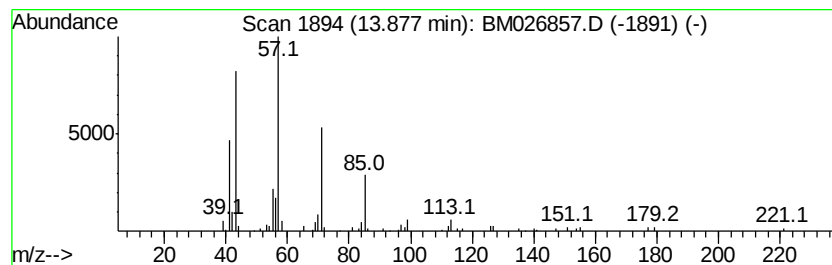
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Pentadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.88	2.11 ng	100262	Acenaphthene-d10		13.95
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	95
2	Tetradecane	198	C14H30	000629-59-4	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tridecane	184	C13H28	000629-50-5	90
5	Heptadecane	240	C17H36	000629-78-7	72



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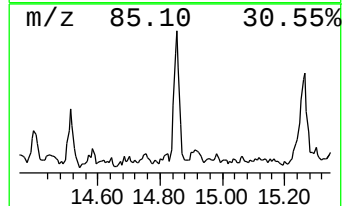
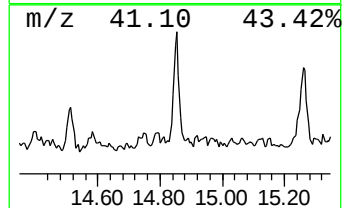
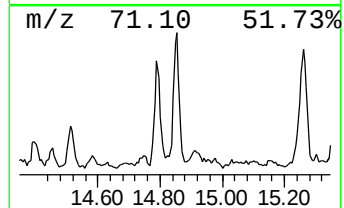
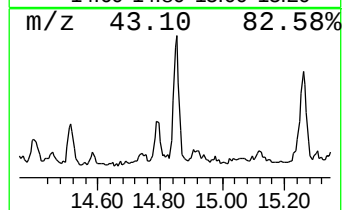
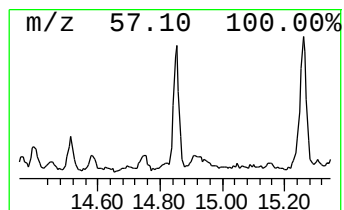
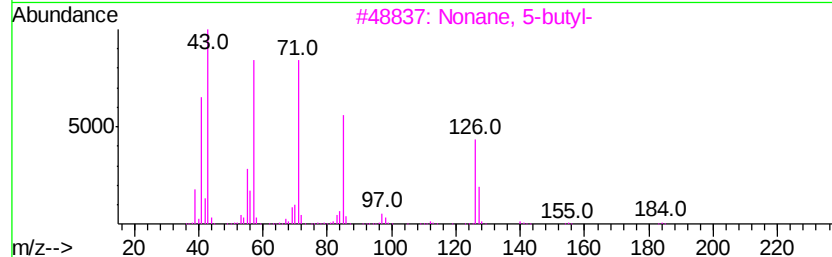
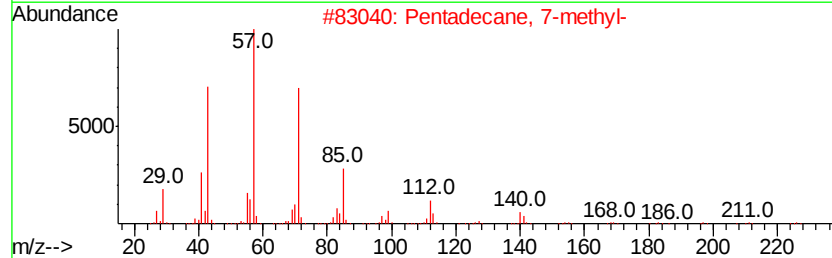
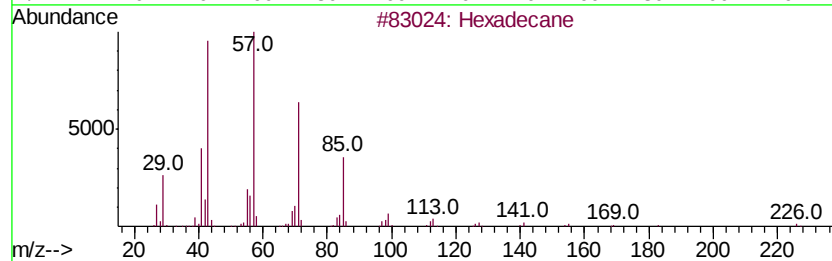
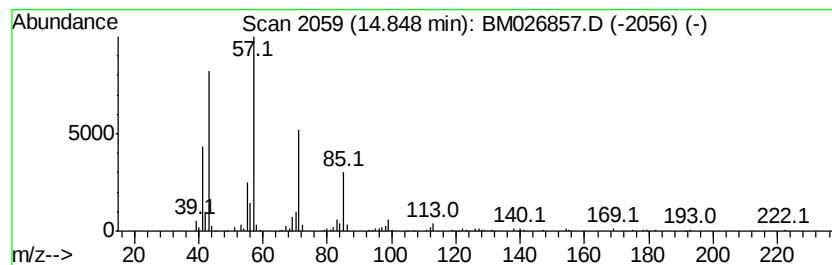
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.66 ng	126651	Acenaphthene-d10	13.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	93
2	Pentadecane, 7-methyl-	226 C16H34	006165-40-8	83
3	Nonane, 5-butyl-	184 C13H28	017312-63-9	72
4	Undecane, 6-ethyl-	184 C13H28	017312-60-6	59
5	Tridecane, 7-methyl-	198 C14H30	026730-14-3	50



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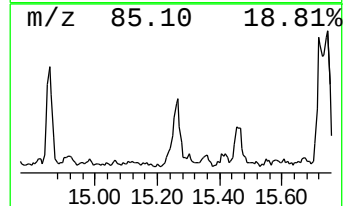
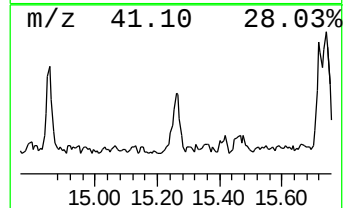
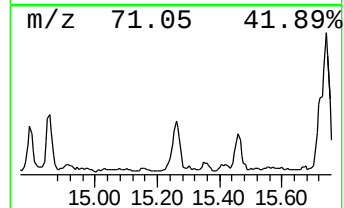
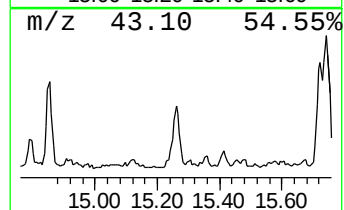
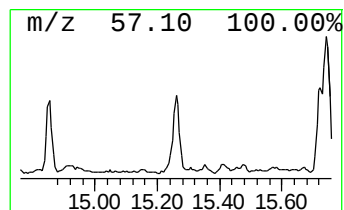
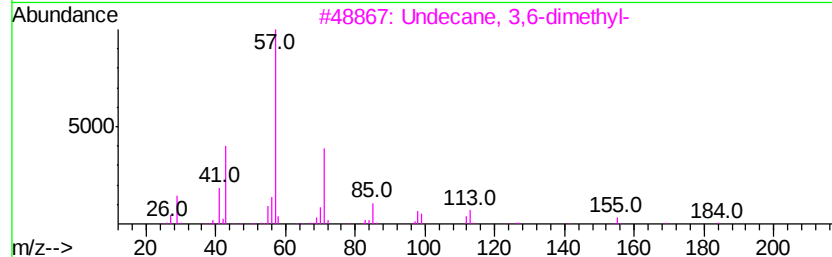
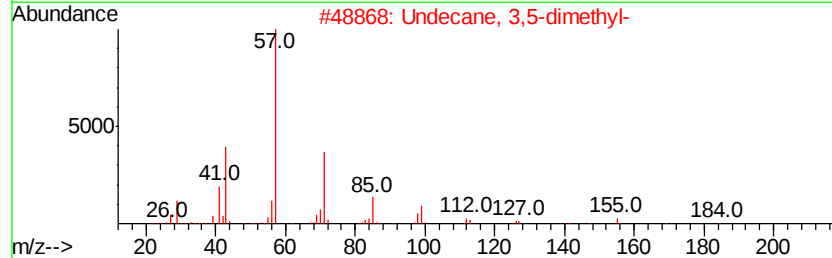
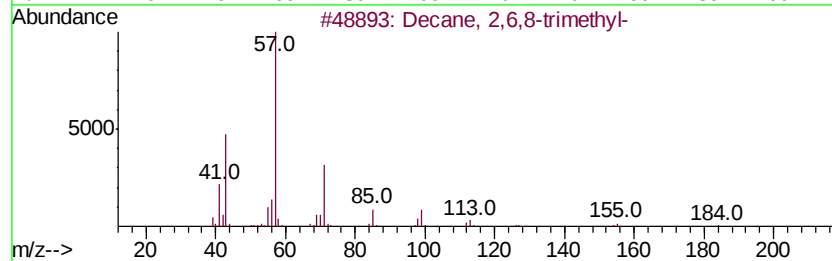
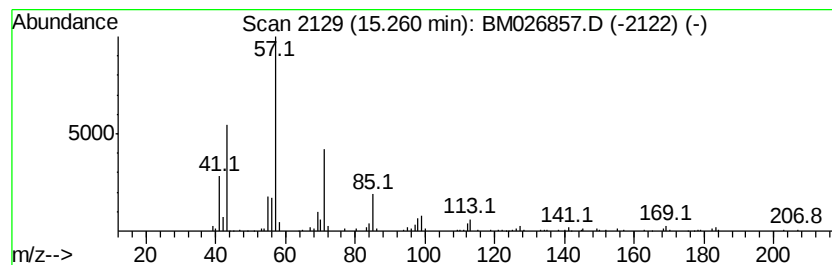
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Decane, 2,6,8-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	3.47 ng	165289	Acenaphthene-d10	13.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	87
2		Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	87
3		Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	83
4		Heptacosane	380	C27H56	000593-49-7	80
5		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	72



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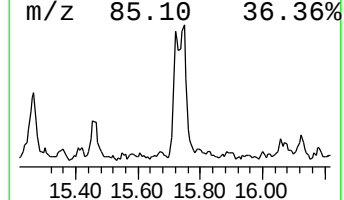
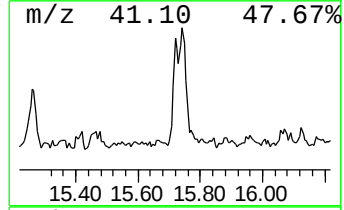
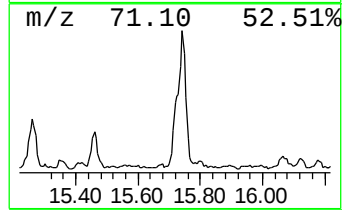
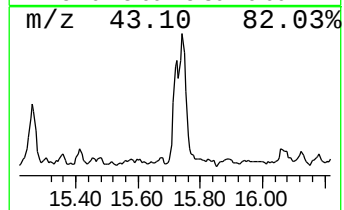
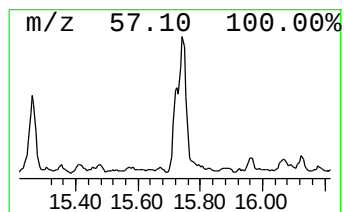
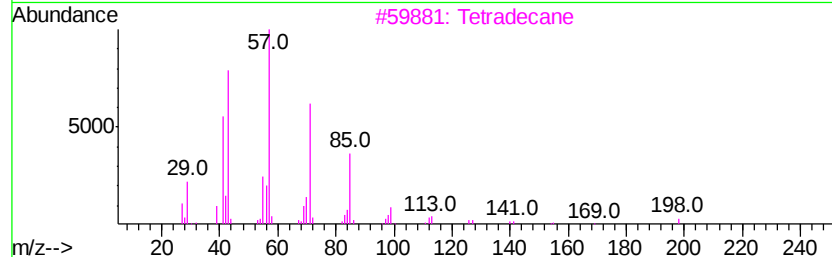
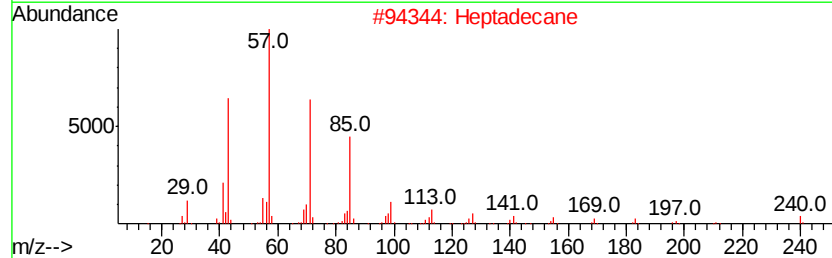
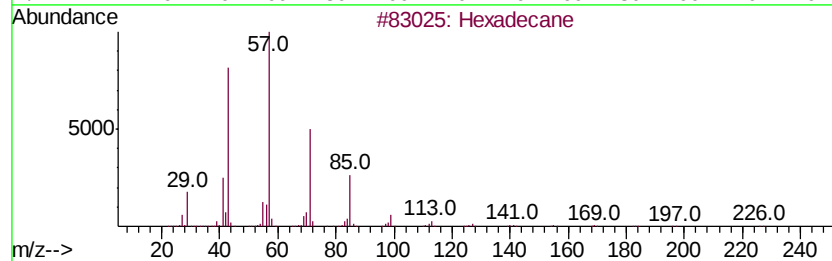
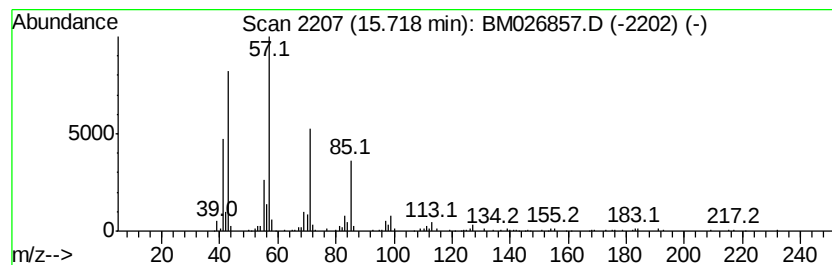
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	3.52 ng	180297	Phenanthrene-d10	16.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Heptadecane	240	C17H36	000629-78-7	90
3	Tetradecane	198	C14H30	000629-59-4	90
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	90
5	Heneicosane	296	C21H44	000629-94-7	90



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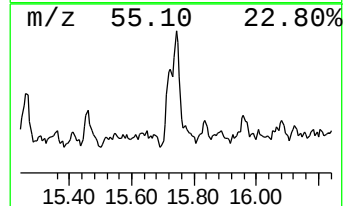
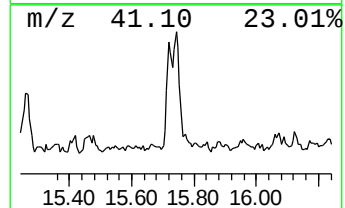
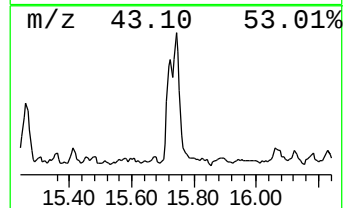
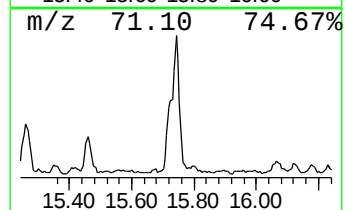
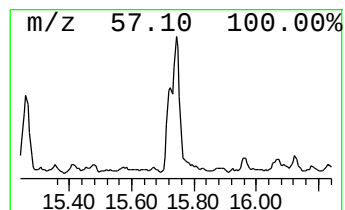
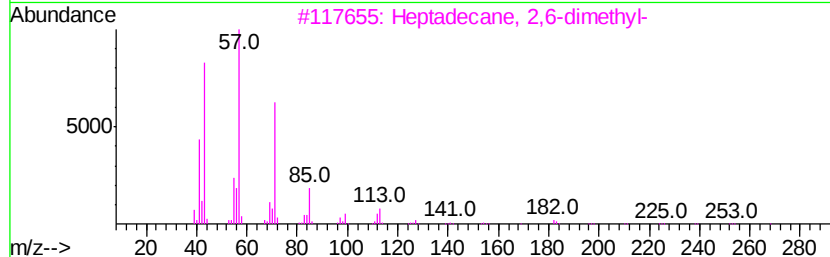
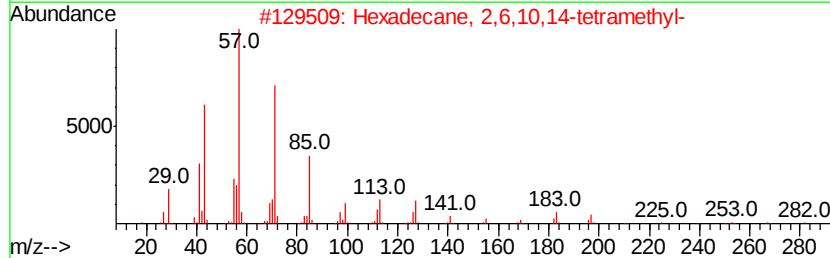
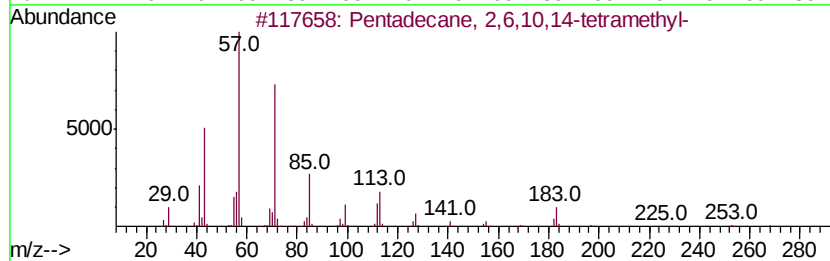
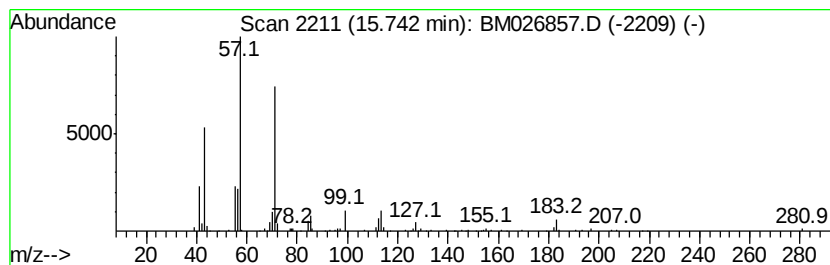
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Peak Number 5 Pentadecane, 2,6,10,14-tetr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.74	5.02 ng	257056	Phenanthrene-d10	16.71		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	83
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78
3		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	78
4		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
5		Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	72



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Operator : JU/CG
Sample : L3381-13
Misc :
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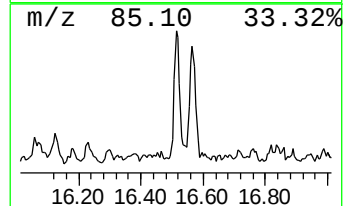
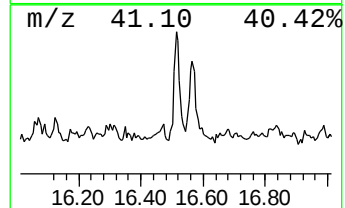
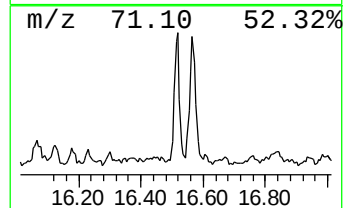
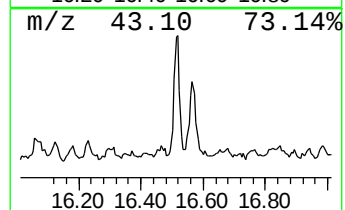
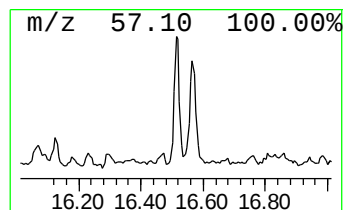
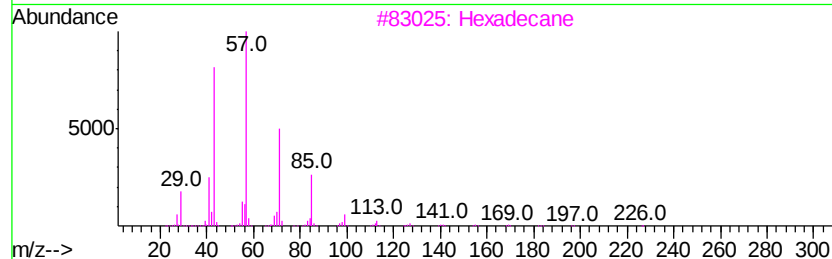
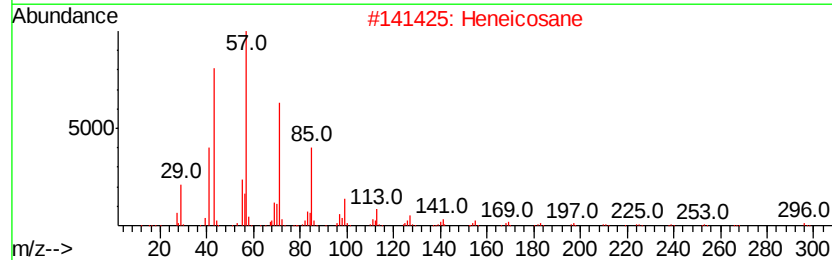
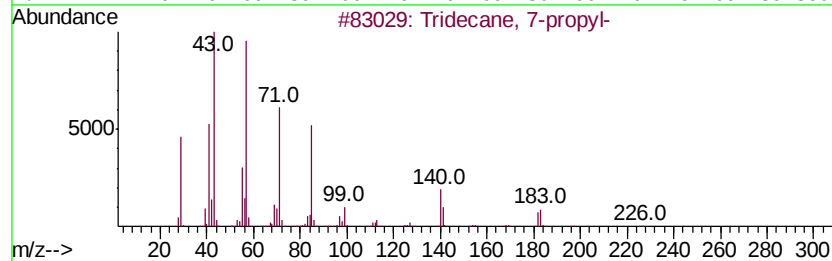
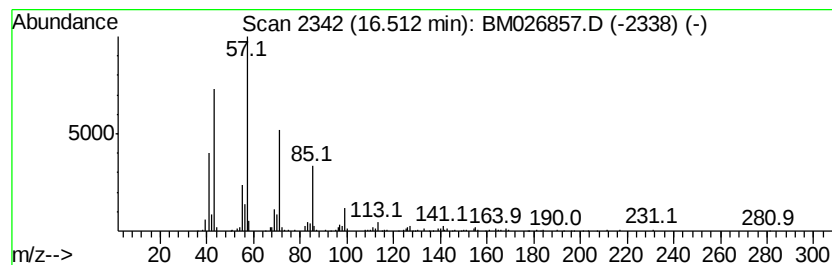
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecane, 7-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.51	2.94 ng	150221	Phenanthrene-d10	16.71	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 7-propyl-	226	C16H34	055045-09-5	86
2	Heneicosane	296	C21H44	000629-94-7	86
3	Hexadecane	226	C16H34	000544-76-3	80
4	Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	80
5	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
Data File : BM026857.D
Acq On : 23 Jul 2020 18:45
Operator : JU/CG
Sample : L3381-13
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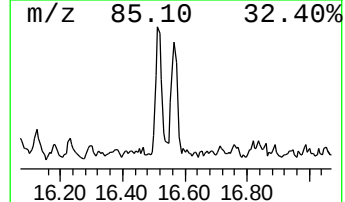
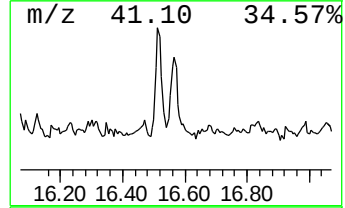
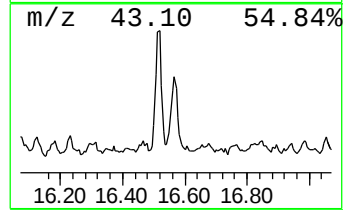
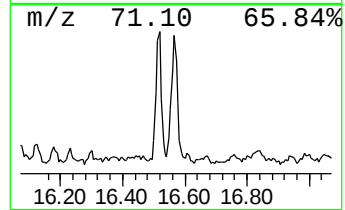
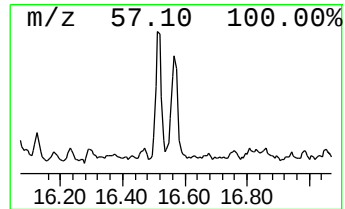
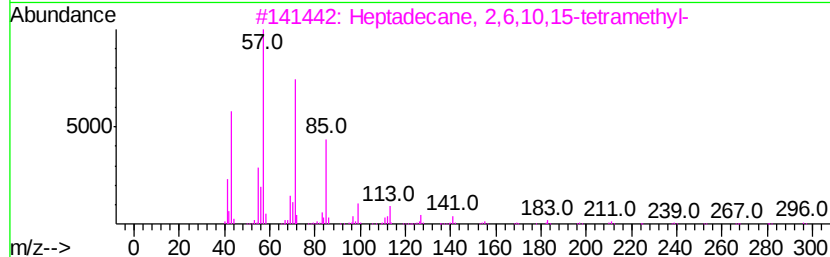
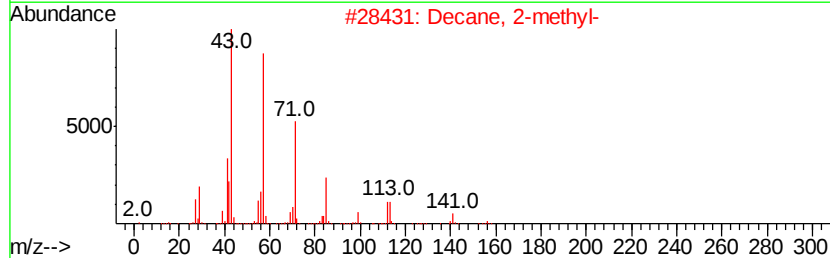
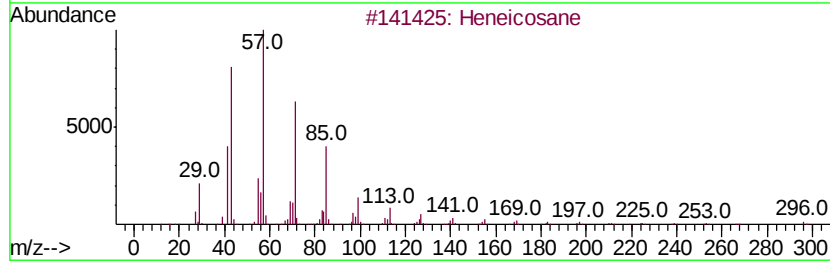
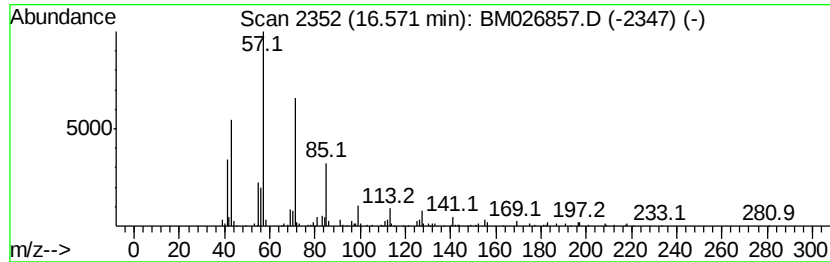
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.57	2.63 ng	134786	Phenanthrene-d10	16.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	86
2	Decane, 2-methyl-	156	C11H24	006975-98-0	80
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
4	Octane, 5-ethyl-2-methyl-	156	C11H24	062016-18-6	72
5	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
Data File : BM026857.D
Acq On : 23 Jul 2020 18:45
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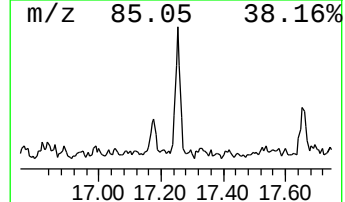
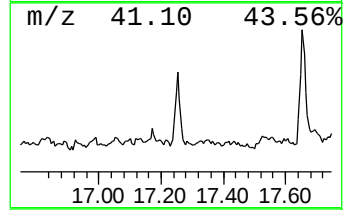
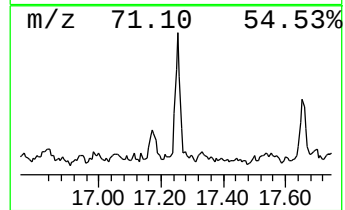
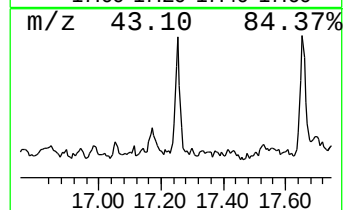
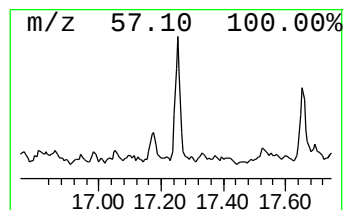
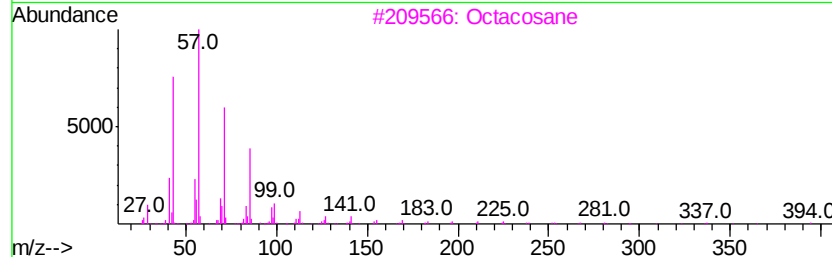
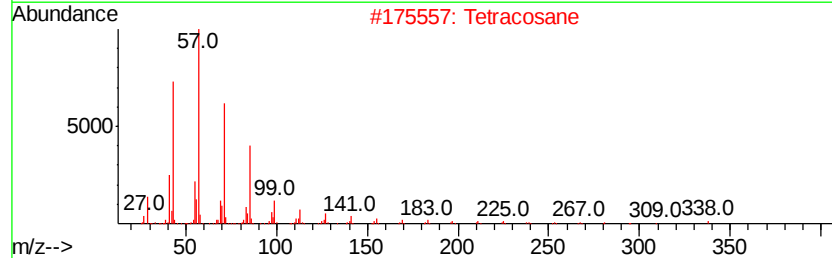
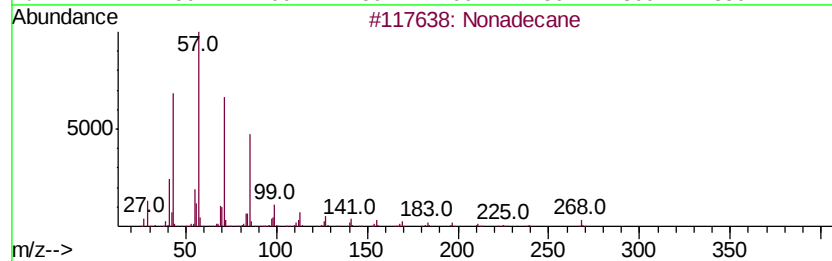
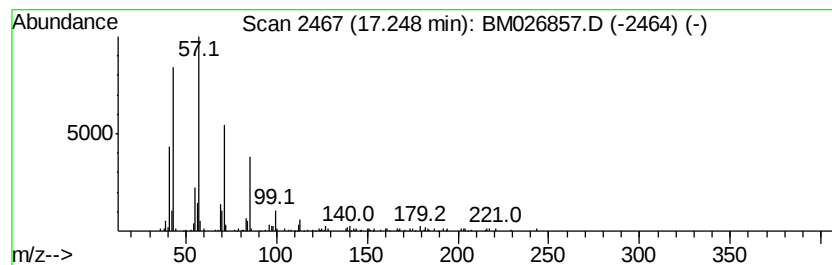
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.25	2.17 ng	110896	Phenanthrene-d10	16.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	94
2	Tetracosane		338	C24H50	000646-31-1	90
3	Octacosane		394	C28H58	000630-02-4	90
4	2-Bromo dodecane		248	C12H25Br	013187-99-0	86
5	Eicosane, 10-methyl-		296	C21H44	054833-23-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
Data File : BM026857.D
Acq On : 23 Jul 2020 18:45
Operator : JU/CG
Sample : L3381-13
Misc :
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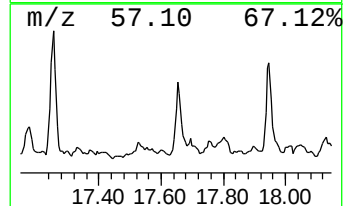
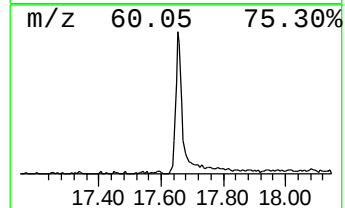
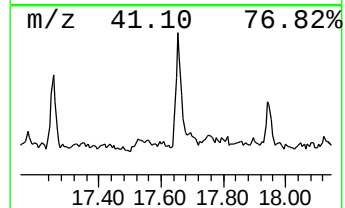
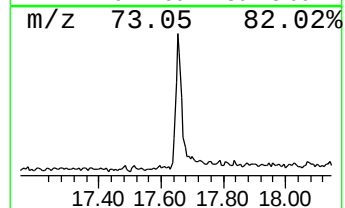
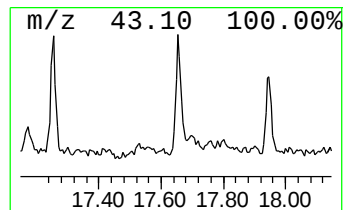
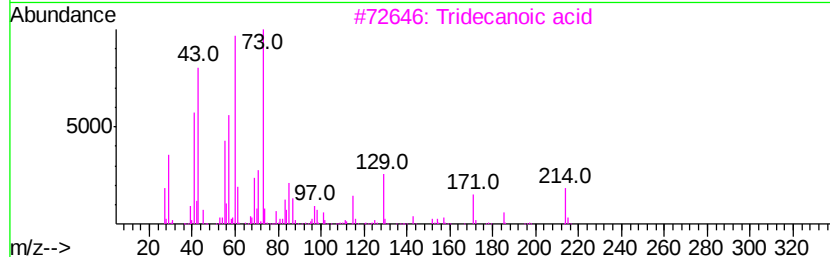
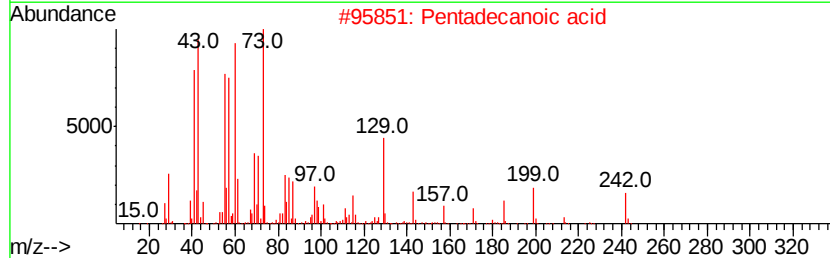
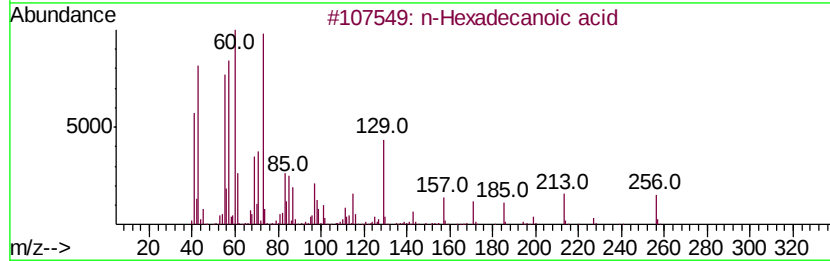
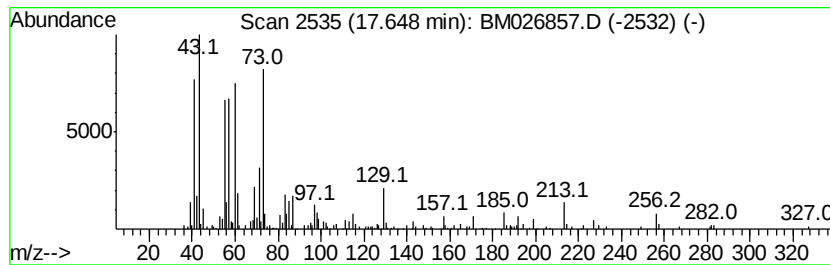
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.65	4.50 ng	230157	Phenanthrene-d10	16.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	Tridecanoic acid	214	C13H26O2	000638-53-9	87
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5	Undecanoic acid	186	C11H22O2	000112-37-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
Data File : BM026857.D
Acq On : 23 Jul 2020 18:45
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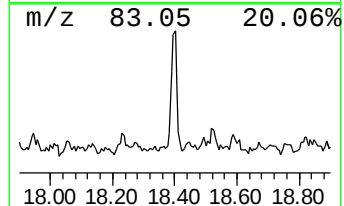
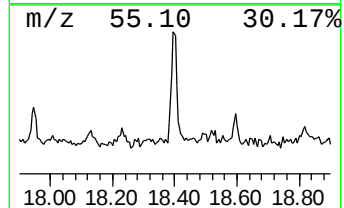
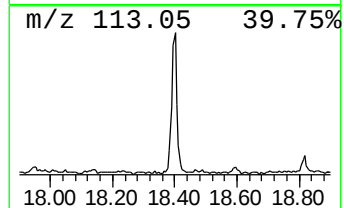
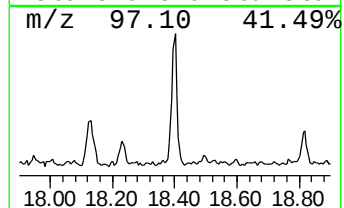
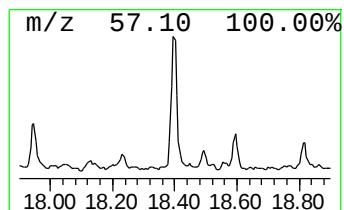
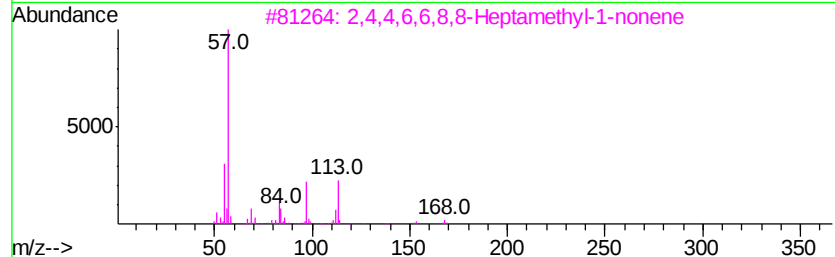
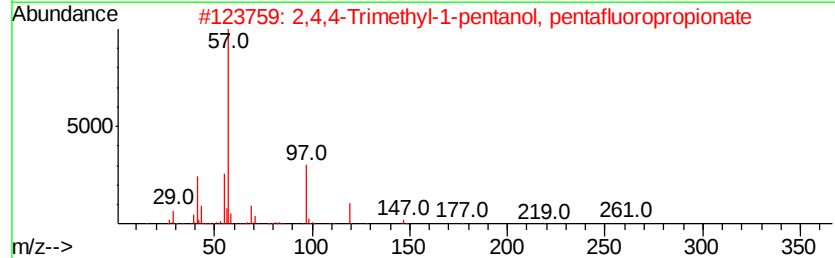
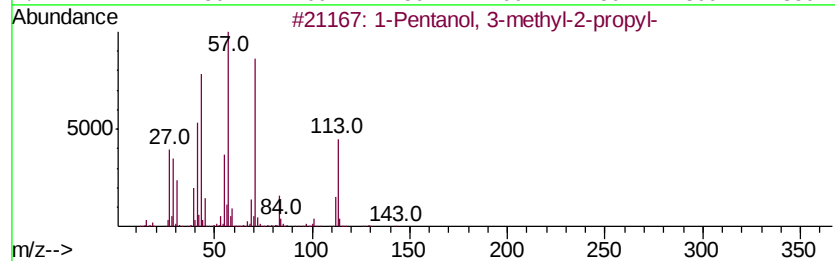
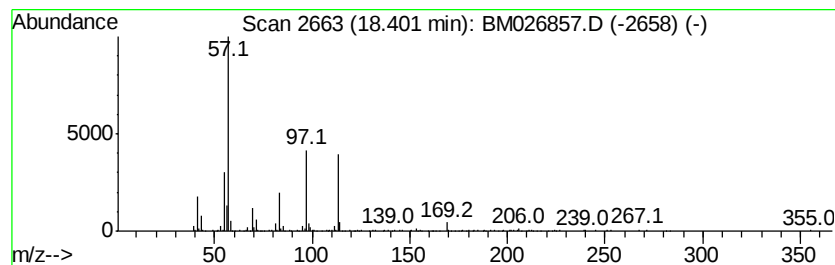
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown18.40 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	4.05 ng	207399	Phenanthrene-d10	16.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	40
2			2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	32
3			2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	32
4			2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	25
5			Formic acid, 2,4,4-trimethylpent...	158	C9H18O2	1000368-95-2	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072320\
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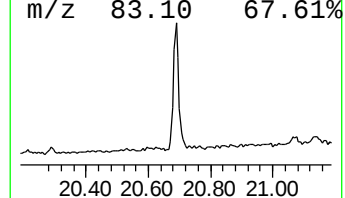
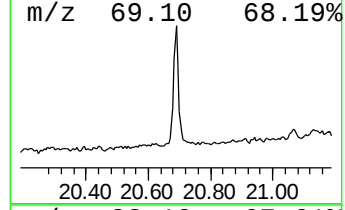
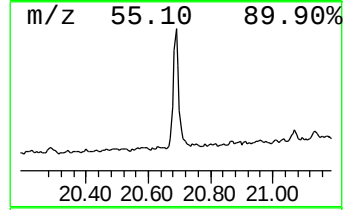
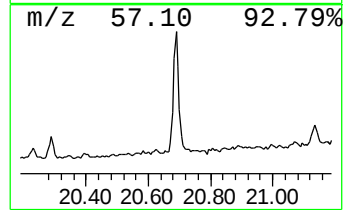
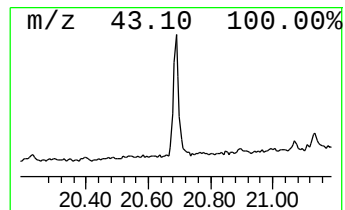
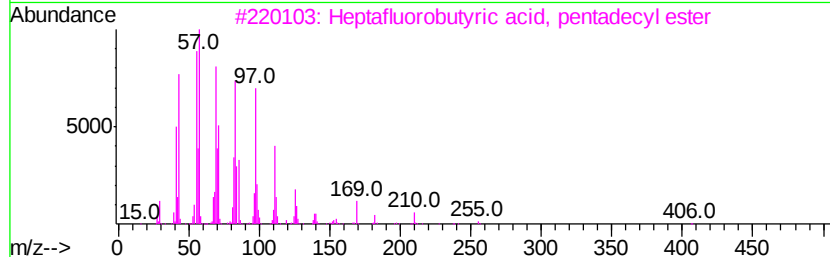
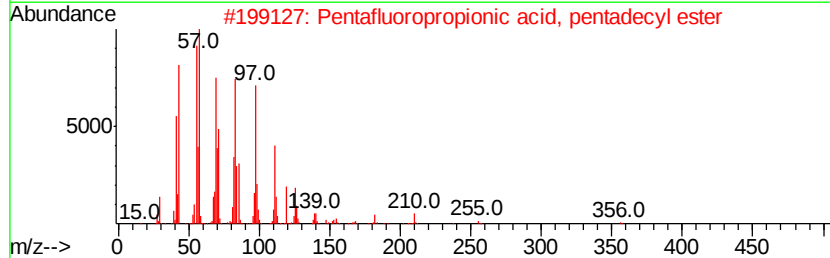
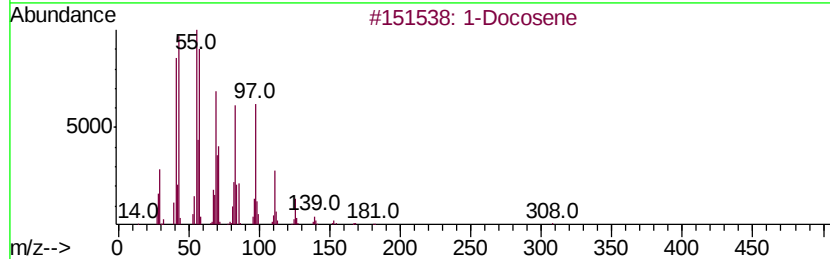
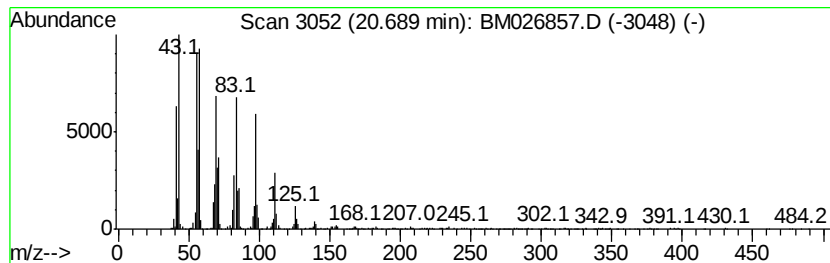
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Docosene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	11.63 ng	732515	Chrysene-d12	20.94
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene		308 C22H44	001599-67-3 95
2	Pentafluoropropionic acid, penta...		374 C18H31F5O2	959092-08-1 94
3	Heptafluorobutyric acid, pentade...		424 C19H31F7O2	959261-23-5 94
4	Heptadecyl trifluoroacetate		352 C19H35F3O2	1000351-87-0 94
5	1-Heneicosanol		312 C21H44O	015594-90-8 94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM072320\
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 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
Pentadecane	13.88	2.1 ng		100262	3	13.95	952387	20.0
Hexadecane	14.85	2.7 ng		126651	3	13.95	952387	20.0
Decane, 2,6,8-tri...	15.26	3.5 ng		165289	3	13.95	952387	20.0
Heptadecane	15.72	3.5 ng		180297	4	16.71	1023600	20.0
Pentadecane, 2,6,...	15.74	5.0 ng		257056	4	16.71	1023600	20.0
Tridecane, 7-propyl-	16.51	2.9 ng		150221	4	16.71	1023600	20.0
Heneicosane	16.57	2.6 ng		134786	4	16.71	1023600	20.0
Nonadecane	17.25	2.2 ng		110896	4	16.71	1023600	20.0
n-Hexadecanoic acid	17.65	4.5 ng		230157	4	16.71	1023600	20.0
unknown18.40	18.40	4.0 ng		207399	4	16.71	1023600	20.0
1-Docosene	20.69	11.6 ng		732515	5	20.94	1260100	20.0