

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
 Data File : BM026736.D
 Acq On : 10 Jul 2020 16:24
 Operator : JU/CG
 Sample : L3217-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E3-41-BOILER-ROOM

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	4.963	346	353	369	rBV	2510749	3440766	49.64%	7.955%
2	6.534	611	620	633	rBV	2242106	3368247	48.60%	7.788%
3	6.963	688	693	708	rBV	92874	174750	2.52%	0.404%
4	7.322	748	754	760	rBV	502138	742736	10.72%	1.717%
5	7.633	800	807	817	rBV	2135362	3187522	45.99%	7.370%
6	8.475	942	950	961	rBV	1652573	2710827	39.11%	6.268%
7	10.074	1216	1222	1229	rBV	584583	918243	13.25%	2.123%
8	12.592	1643	1650	1665	rVB	3548253	5089897	73.43%	11.768%
9	12.898	1696	1702	1709	rBV4	27775	72349	1.04%	0.167%
10	13.457	1792	1797	1801	rBV	553338	761839	10.99%	1.761%
11	13.498	1801	1804	1808	rVB	93933	122030	1.76%	0.282%
12	13.857	1861	1865	1870	rBV6	37211	73186	1.06%	0.169%
13	13.980	1881	1886	1892	rVB2	735608	1047008	15.11%	2.421%
14	14.433	1960	1963	1967	rBV4	41588	79491	1.15%	0.184%
15	14.545	1979	1982	1988	rVB4	84679	120055	1.73%	0.278%
16	14.715	2008	2011	2014	rBV4	50096	79542	1.15%	0.184%
17	14.786	2020	2023	2027	rVB3	68096	82902	1.20%	0.192%
18	15.298	2106	2110	2114	rVB	276245	403418	5.82%	0.933%
19	15.492	2138	2143	2150	rVB	2788400	3922934	56.60%	9.070%
20	15.780	2187	2192	2199	rVB	1465123	2041297	29.45%	4.720%
21	15.874	2205	2208	2211	rVB	154203	183172	2.64%	0.424%
22	16.115	2246	2249	2253	rVB2	189626	247887	3.58%	0.573%
23	16.598	2327	2331	2336	rVB	1034460	1513152	21.83%	3.499%
24	16.745	2352	2356	2361	rVB	844502	1074657	15.50%	2.485%
25	17.203	2431	2434	2439	rVB	331130	459103	6.62%	1.061%
26	17.686	2513	2516	2521	rVB	461931	589260	8.50%	1.362%
27	19.280	2784	2787	2793	rVB	231858	325749	4.70%	0.753%
28	19.409	2804	2809	2820	rVB	5689317	6931245	100.00%	16.026%
29	20.191	2940	2942	2945	rVB	152144	141419	2.04%	0.327%
30	20.715	3027	3031	3037	rVB	602450	696518	10.05%	1.610%
31	20.968	3069	3074	3081	rVB	1046672	1301064	18.77%	3.008%
32	23.032	3420	3425	3434	rVB2	799794	1348636	19.46%	3.118%

LSC Area Percent Report

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ClientSampleId :
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM070820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

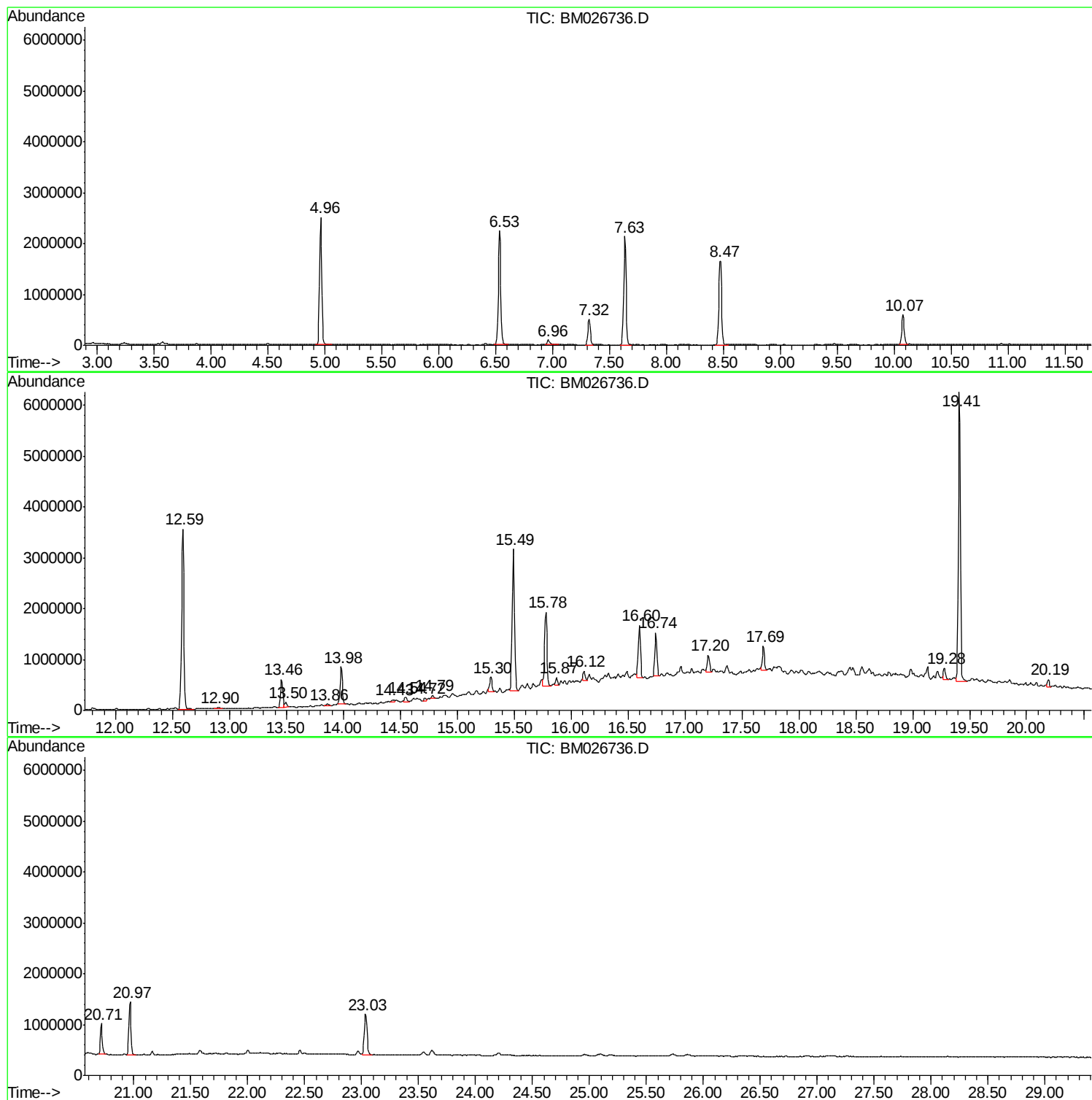
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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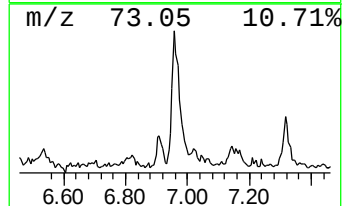
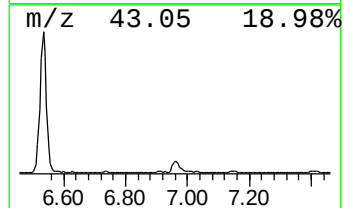
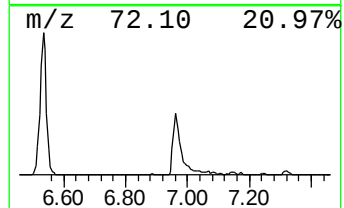
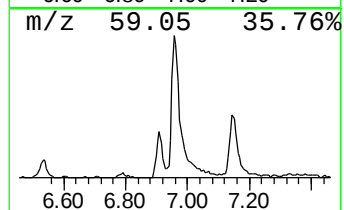
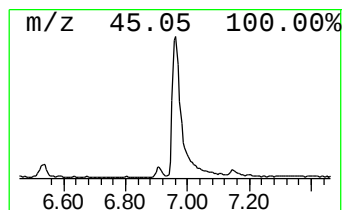
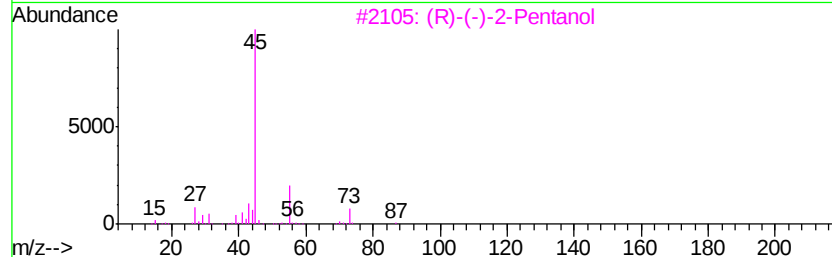
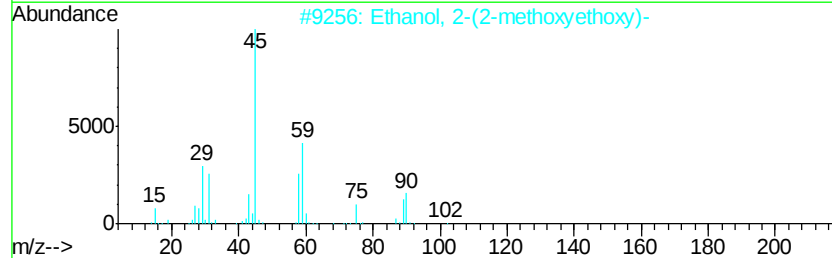
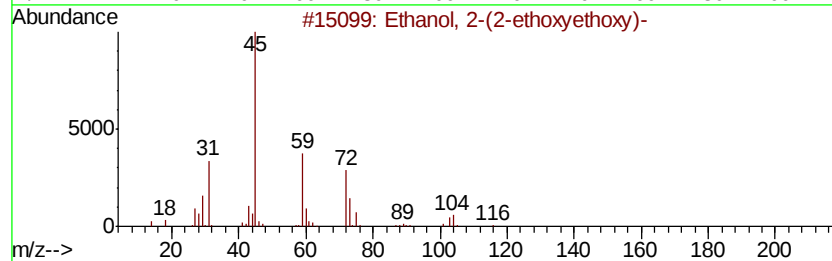
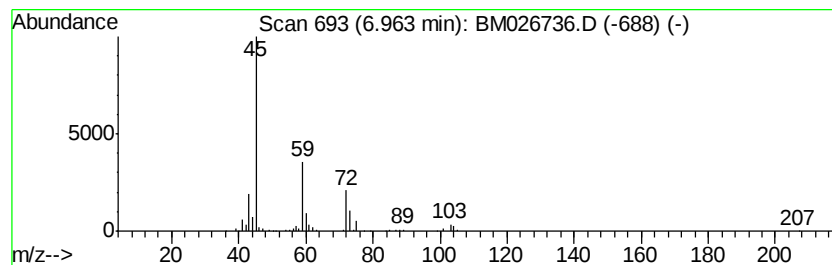
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.96	4.71 ng	174750	1,4-Dichlorobenzene-d4	7.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2	Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	42
3	(R)-(-)-2-Pentanol	88	C5H12O	031087-44-2	38
4	.beta.-Methoxyethoxymethyl chloride	124	C4H9ClO2	003970-21-6	36
5	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	36



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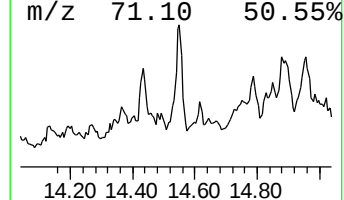
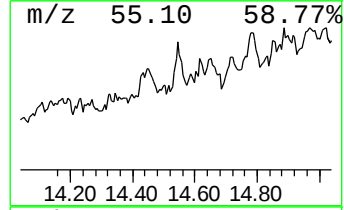
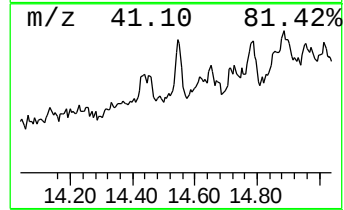
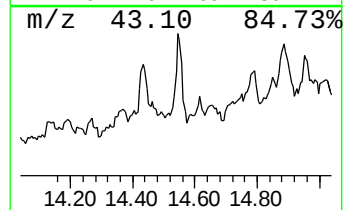
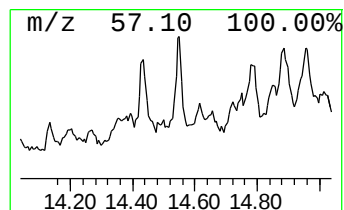
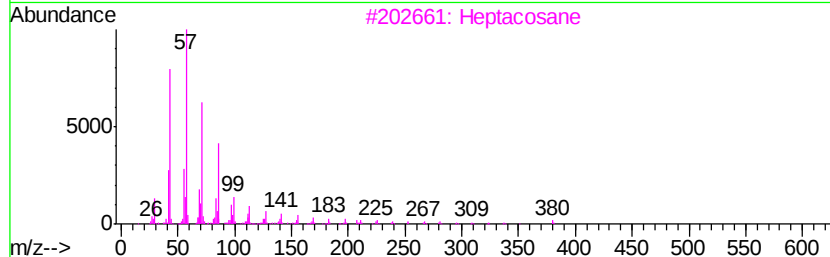
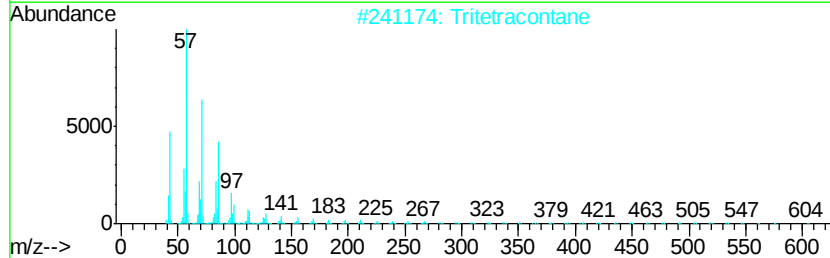
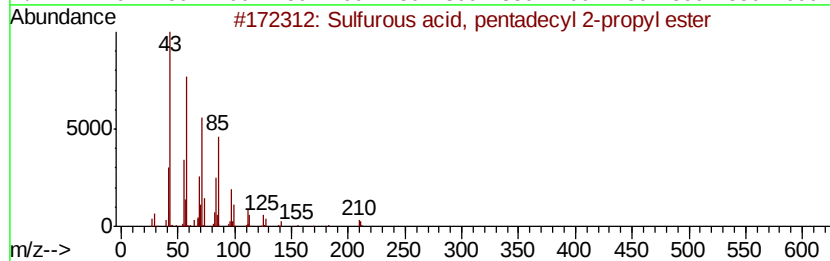
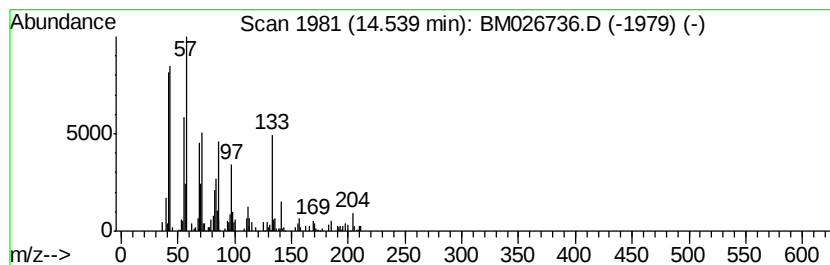
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Sulfurous acid, pentadecyl ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	2.29 ng	120055	Acenaphthene-d10	13.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Sulfurous acid, pentadecyl 2-pro...	334 C18H38O3S	1000309-12-6 86
2		Tritetracontane	605 C43H88	007098-21-7 80
3		Heptacosane	380 C27H56	000593-49-7 72
4		Sulfurous acid, butyl heptadecyl...	376 C21H44O3S	1000309-18-4 72
5		Hexacosane	366 C26H54	000630-01-3 64



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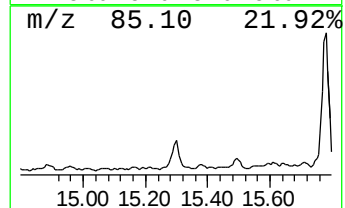
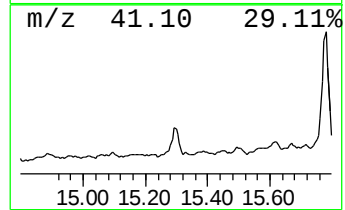
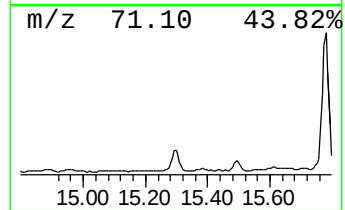
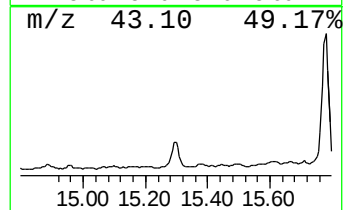
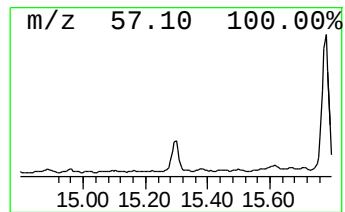
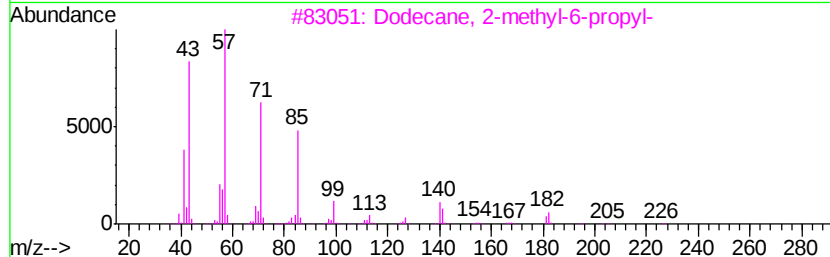
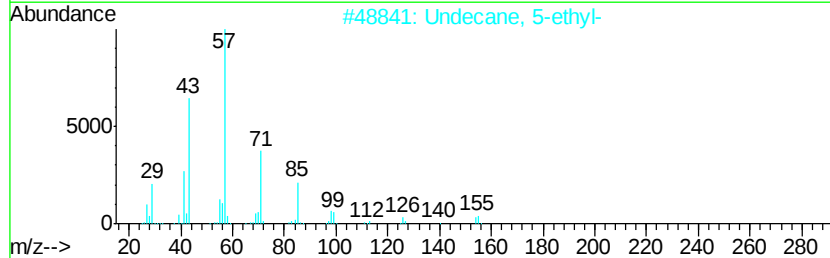
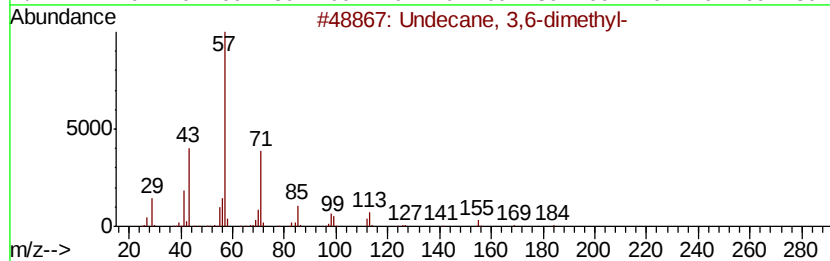
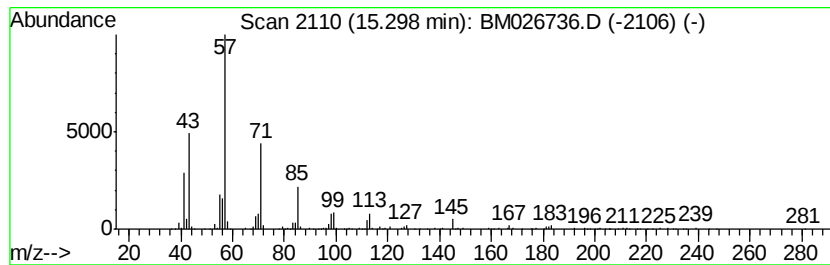
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Undecane, 3,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.30	7.71 ng	403418	Acenaphthene-d10	13.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	72
2	Undecane, 5-ethyl-	184	C13H28	017453-94-0	72
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	59
4	Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	59
5	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	59



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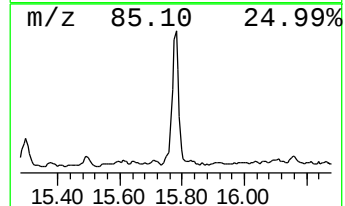
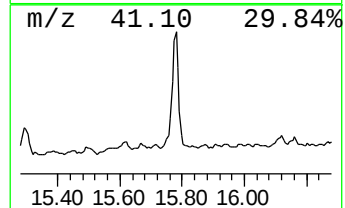
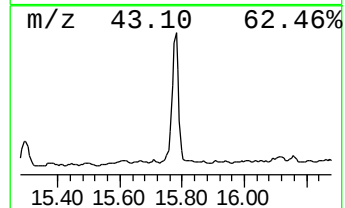
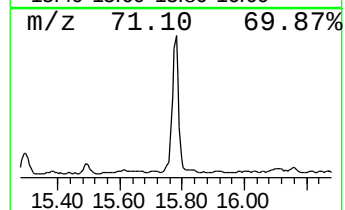
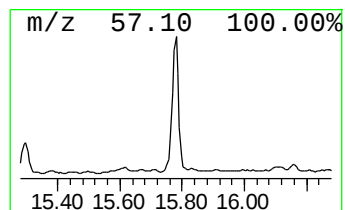
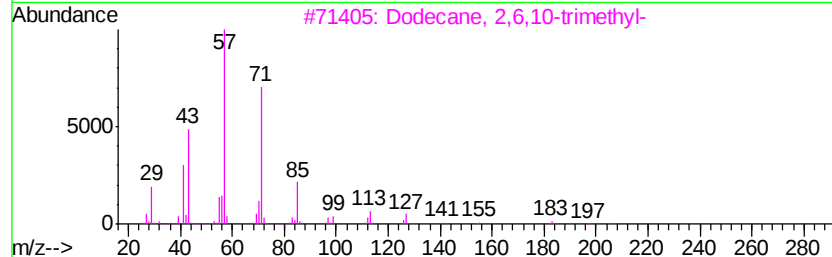
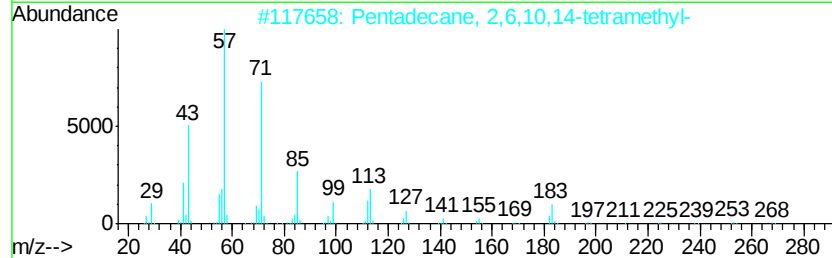
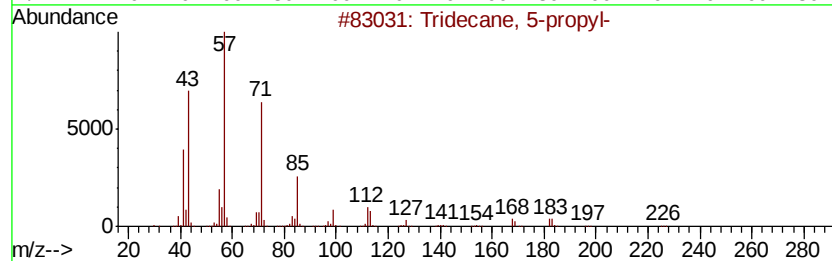
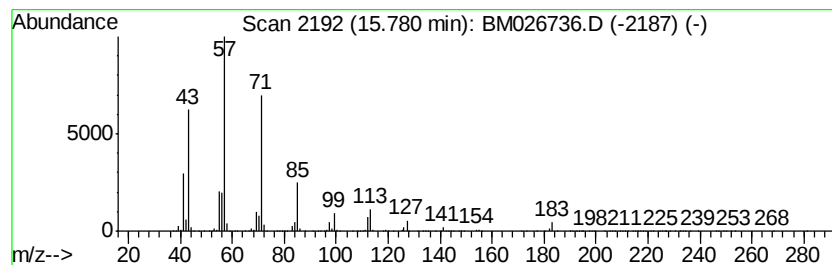
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tridecane, 5-propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.78	37.99 ng	2041300	Phenanthrene-d10	16.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
3		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	83
4		Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	81
5		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	72



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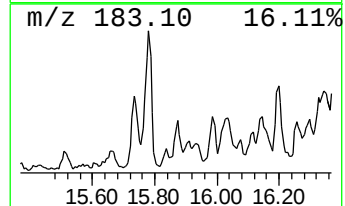
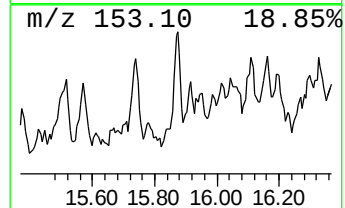
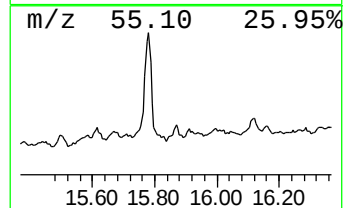
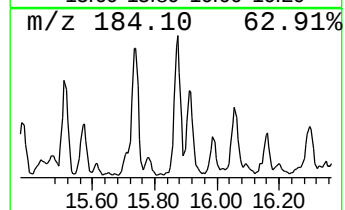
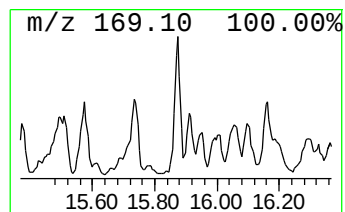
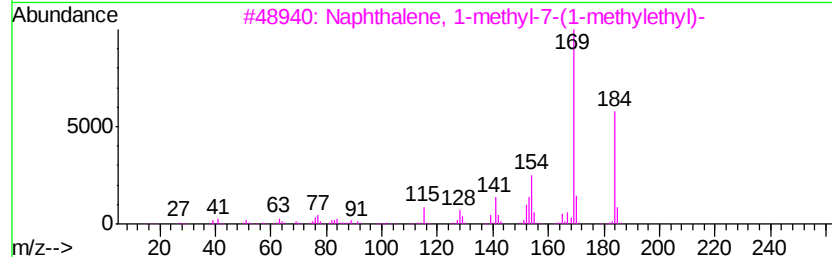
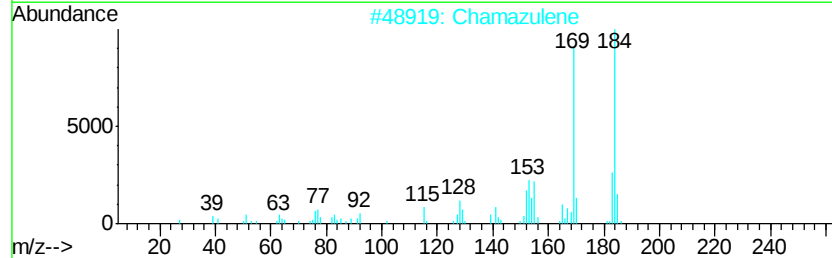
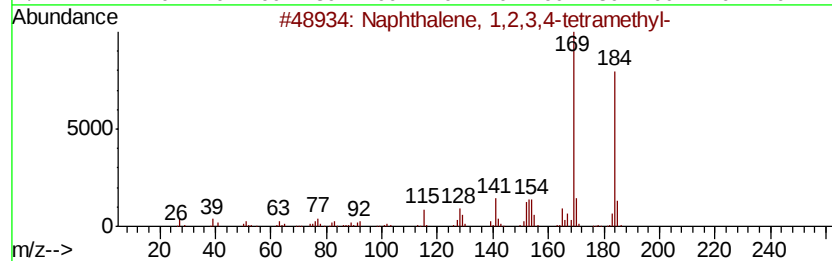
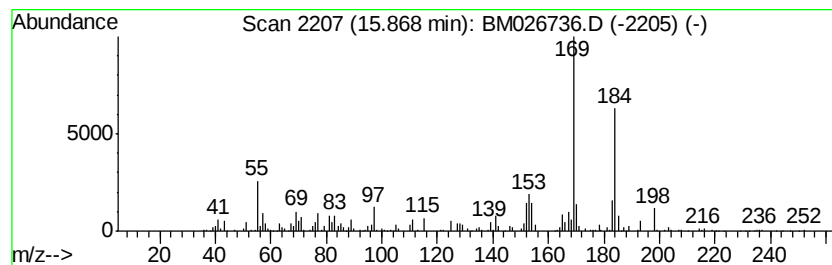
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Peak Number 6 Naphthalene, 1,2,3,4-tetram... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.87	3.41 ng	183172	Phenanthrene-d10	16.74	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	93
2	Chamazulene	184	C14H16	000529-05-5	93
3	Naphthalene, 1-methyl-7-(1-methyl-2-propenyl)-	184	C14H16	000490-65-3	93
4	Ethanone, 1-(5-chloro-2-hydroxy-2-propyl)-	184	C9H9ClO2	028480-70-8	58
5	Naphthalene, 1-(1,1-dimethylethyl)-	184	C14H16	017085-91-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
Data File : BM026736.D
Acq On : 10 Jul 2020 16:24
Operator : JU/CG
Sample : L3217-14
Misc :
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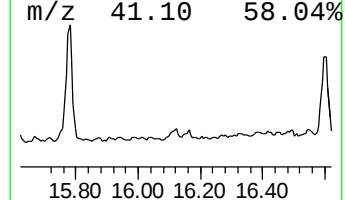
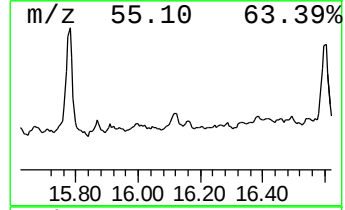
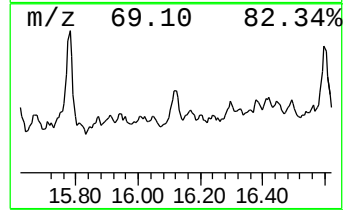
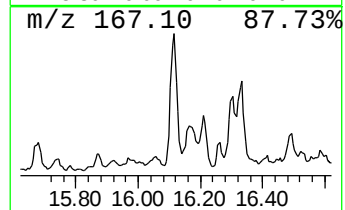
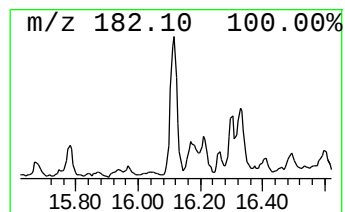
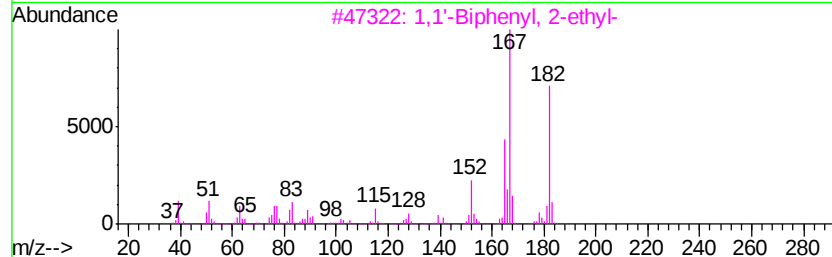
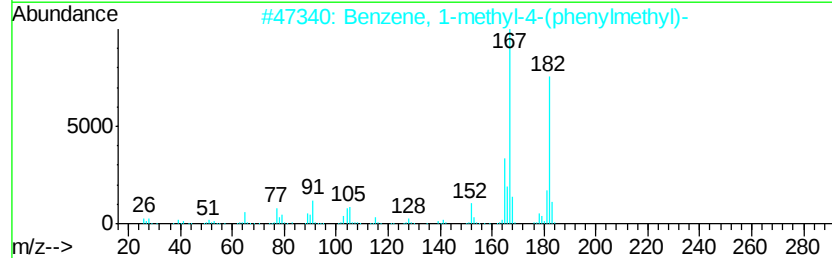
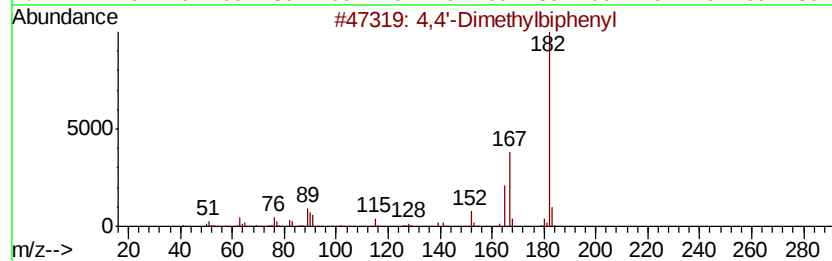
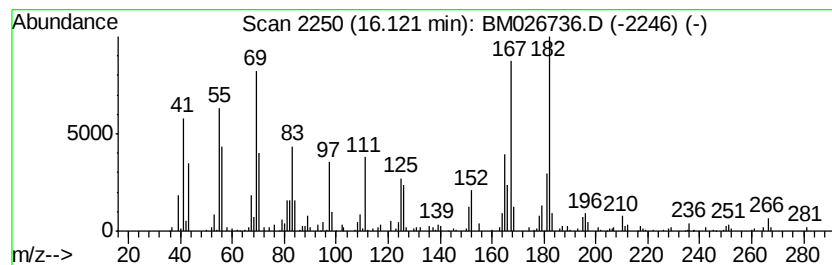
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E3-41-BOILER-ROOM

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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 4,4'-Dimethylbiphenyl Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.12	4.61 ng	247887	Phenanthrene-d10	16.74	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	76
2	Benzene, 1-methyl-4-(phenylmethyl)-	182	C14H14	000620-83-7	74
3	1,1'-Biphenyl, 2-ethyl-	182	C14H14	001812-51-7	72
4	Benzene, 1-methyl-2-(phenylmethyl)-	182	C14H14	000713-36-0	64
5	7-Phenylbicyclo[3.2.1]octa-2,6-d...	182	C14H14	1000201-01-7	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
Data File : BM026736.D
Acq On : 10 Jul 2020 16:24
Operator : JU/CG
Sample : L3217-14
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ClientSampleId :
E3-41-BOILER-ROOM

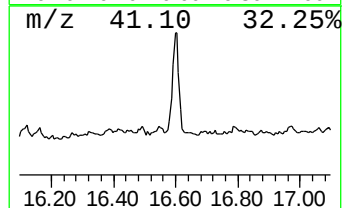
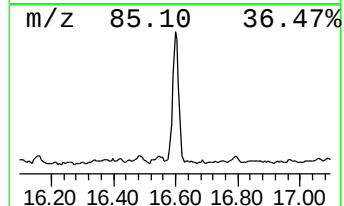
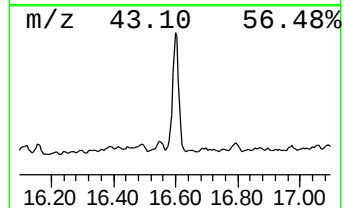
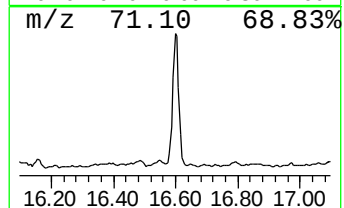
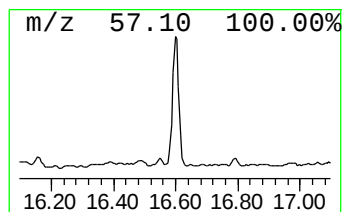
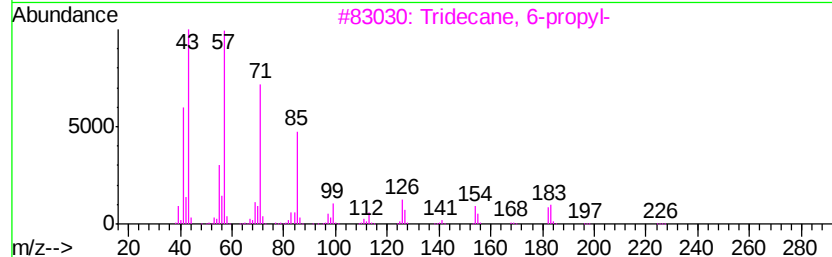
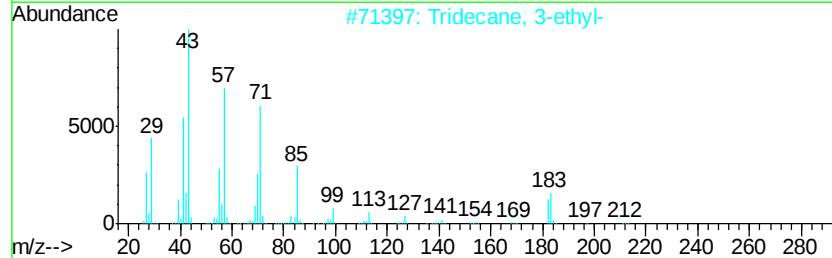
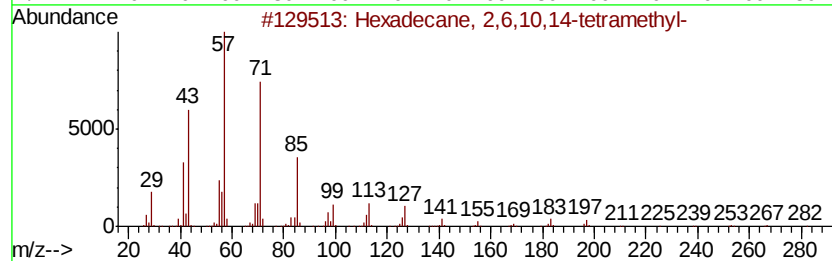
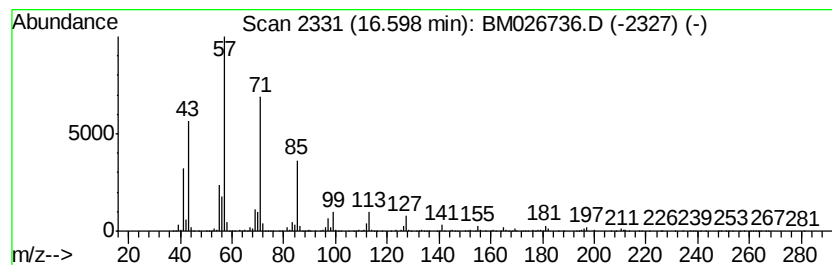
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane, 2,6,10,14-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.60	28.16 ng	1513150	Phenanthrene-d10	16.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Tridecane, 3-ethyl-	212	C15H32	013286-73-2	87
3		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5		Heptacosane	380	C27H56	000593-49-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
Data File : BM026736.D
Acq On : 10 Jul 2020 16:24
Operator : JU/CG
Sample : L3217-14
Misc :
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Instrument :
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ClientSampleId :
E3-41-BOILER-ROOM

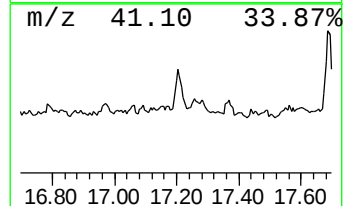
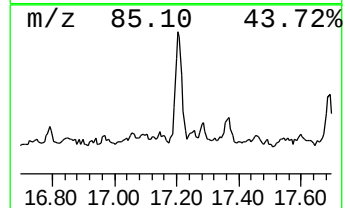
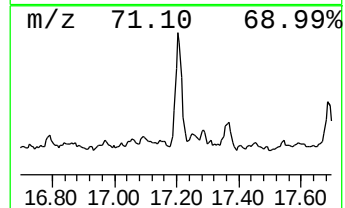
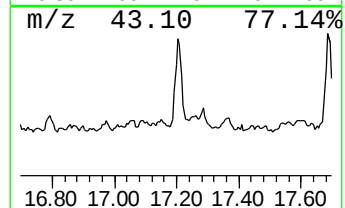
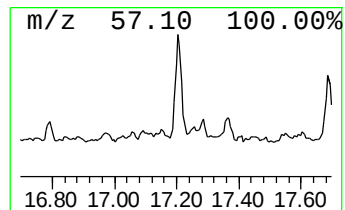
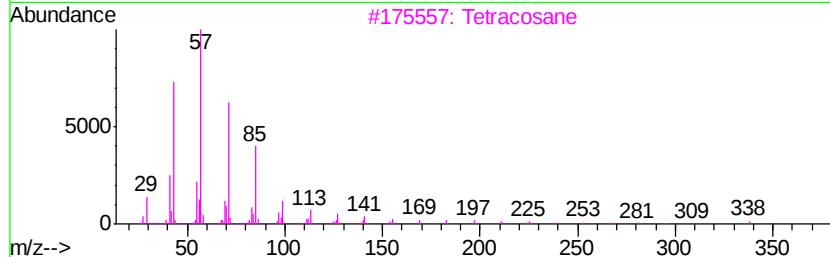
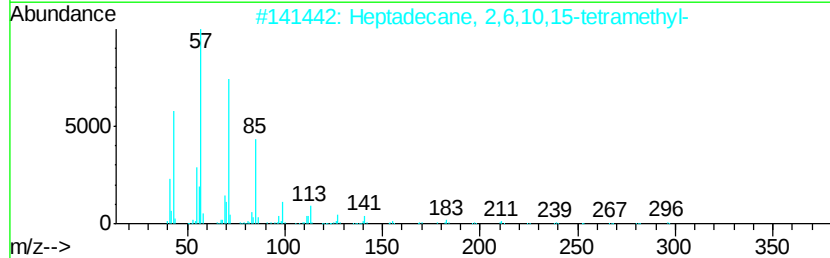
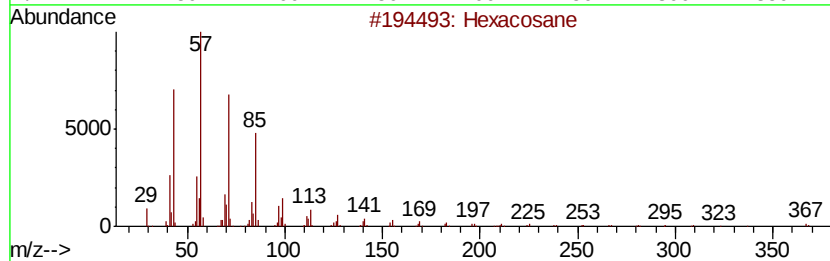
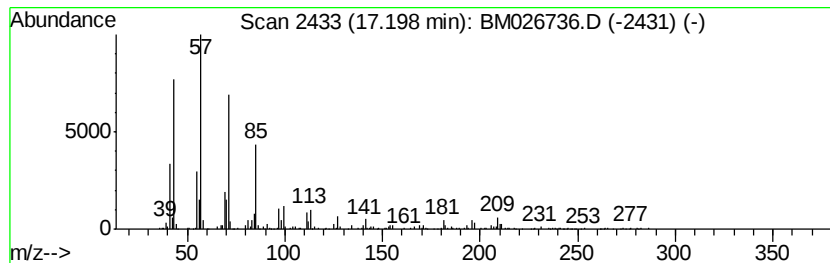
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.20	8.54 ng	459103	Phenanthrene-d10	16.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3	Tetracosane	338	C24H50	000646-31-1	90
4	Nonadecane	268	C19H40	000629-92-5	90
5	Tetradecane	198	C14H30	000629-59-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
Data File : BM026736.D
Acq On : 10 Jul 2020 16:24
Operator : JU/CG
Sample : L3217-14
Misc :
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Instrument :
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ClientSampleId :
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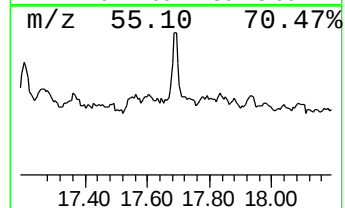
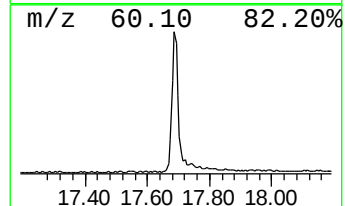
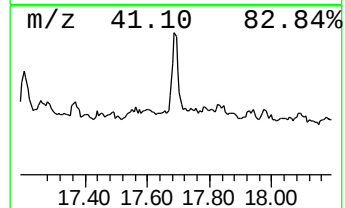
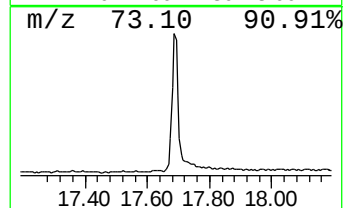
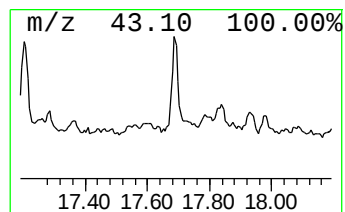
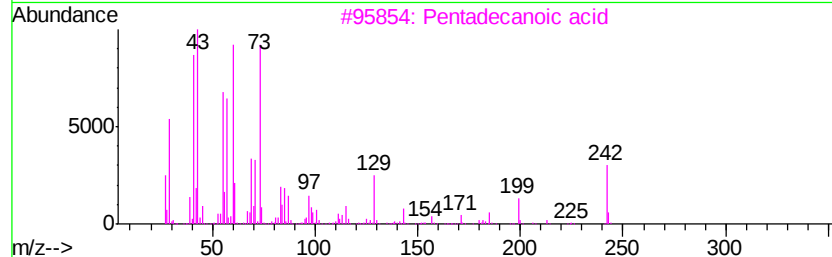
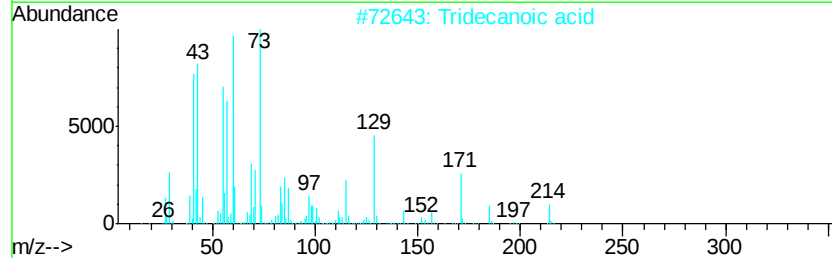
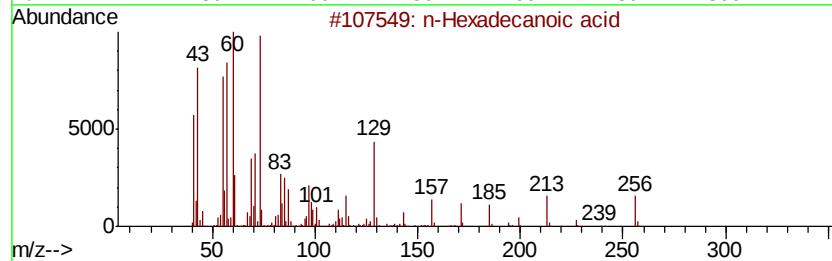
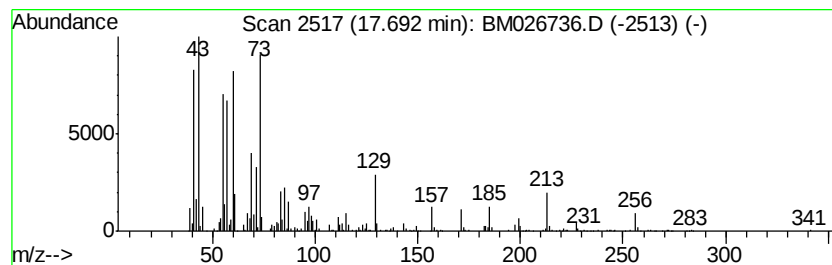
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.69	10.97 ng	589260	Phenanthrene-d10	16.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	91
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	78
5			Octadecanoic acid	284	C18H36O2	000057-11-4	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
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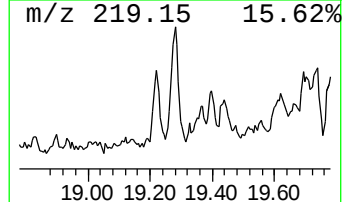
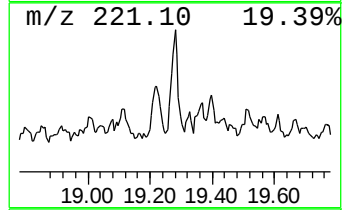
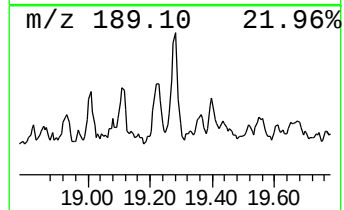
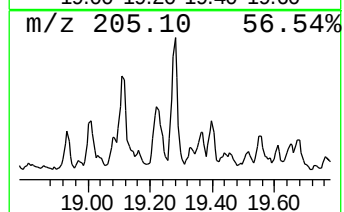
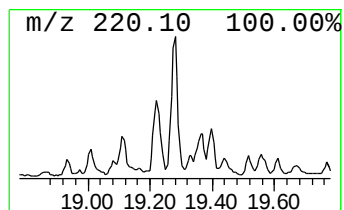
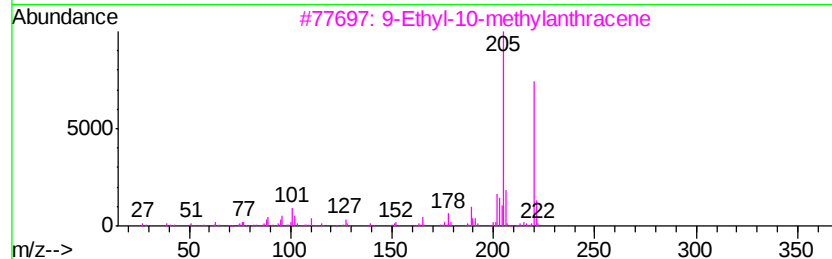
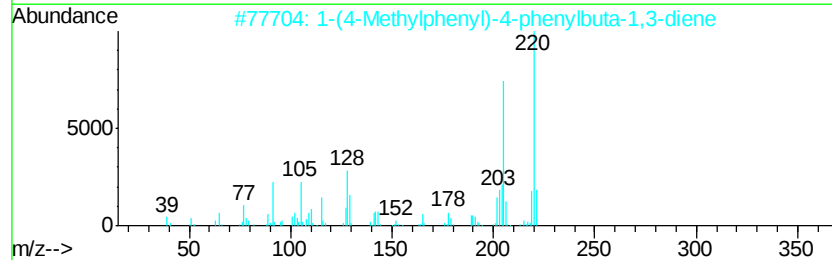
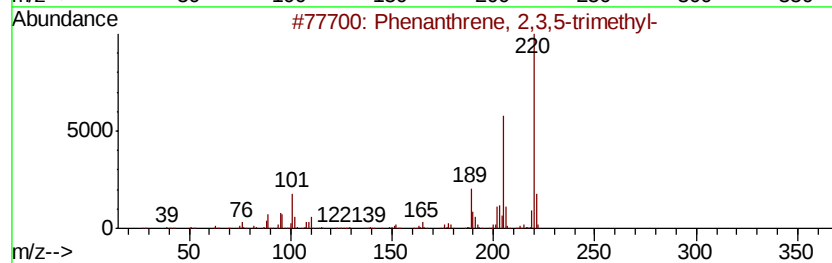
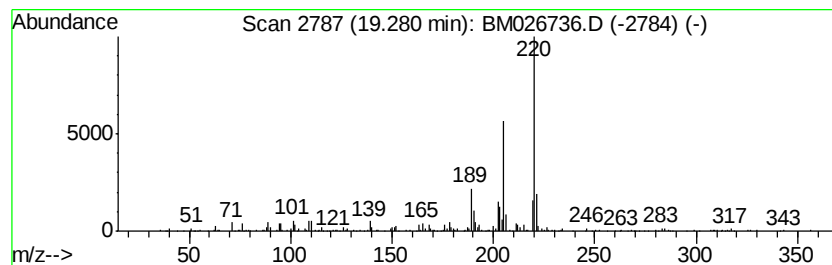
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.28	5.01 ng	325749	Chrysene-d12	20.97		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	95
2		1-(4-Methylphenyl)-4-phenylbuta-...	220	C17H16	037985-11-8	87
3		9-Ethyl-10-methylanthracene	220	C17H16	019713-49-6	64
4		Benzo[b]dihdropyran, 6-hydroxy-...	220	C14H20O2	050442-70-1	53
5		3,5-Cyclohexadiene-1,2-dione, 3,...	220	C14H20O2	003383-21-9	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
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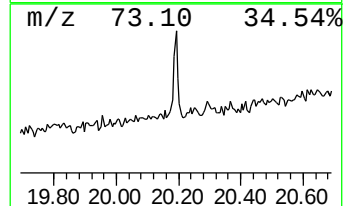
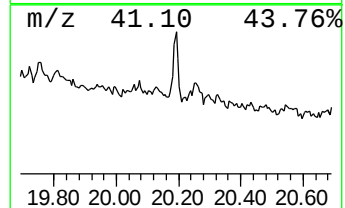
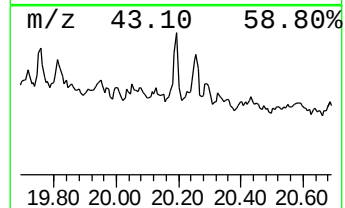
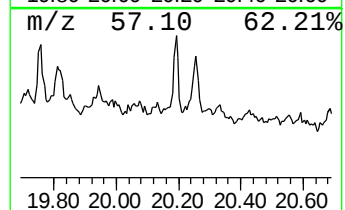
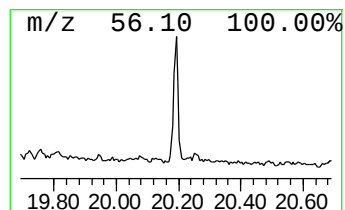
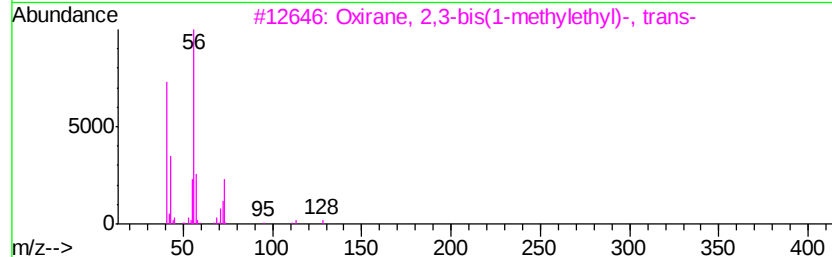
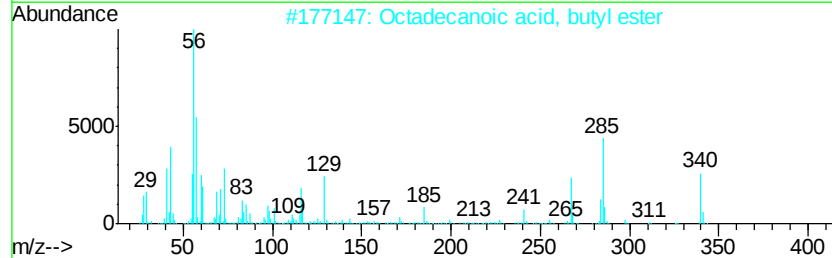
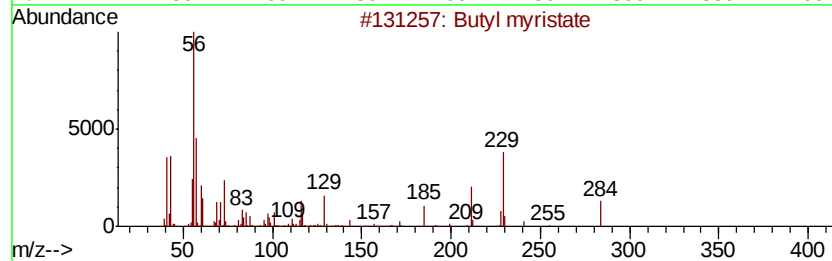
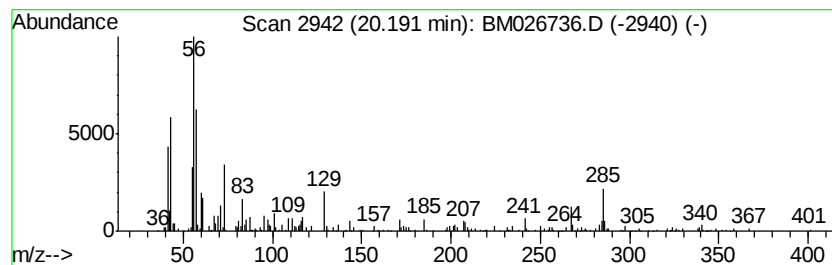
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Butyl myristate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.19	2.17 ng	141419	Chrysene-d12	20.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butyl myristate	284	C18H36O2	000110-36-1	50
2		Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	43
3		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	38
4		2-(Cyanomethoxy)tetrahydropyran	141	C7H11NO2	017521-49-2	35
5		Octadecanoic acid	284	C18H36O2	000057-11-4	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071020\
Data File : BM026736.D
Acq On : 10 Jul 2020 16:24
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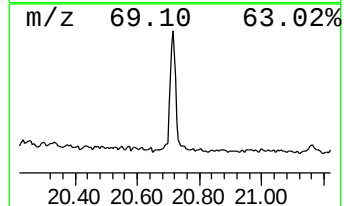
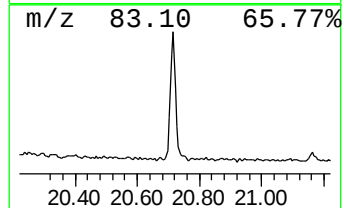
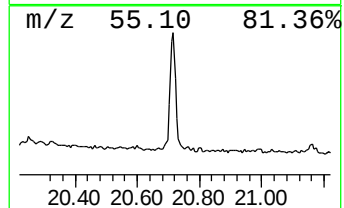
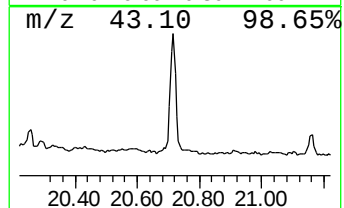
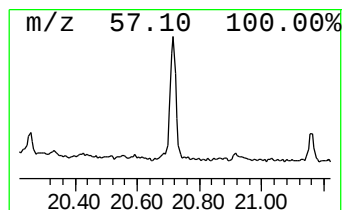
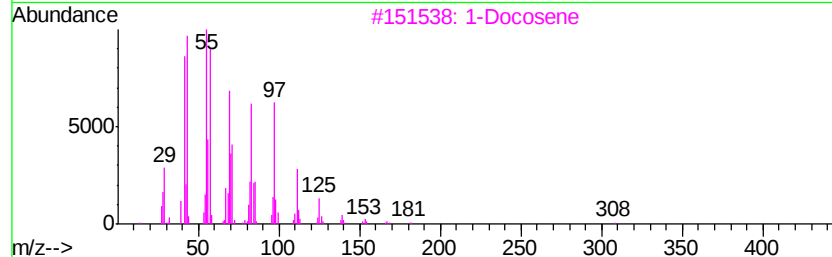
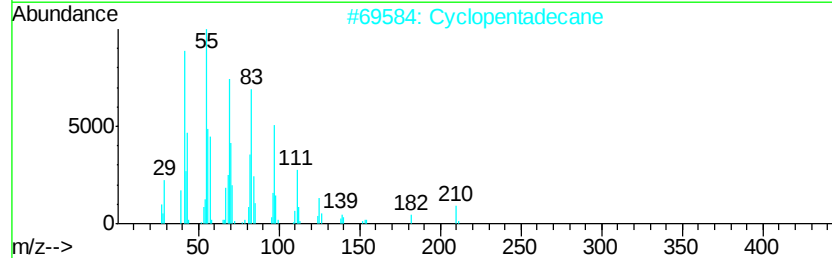
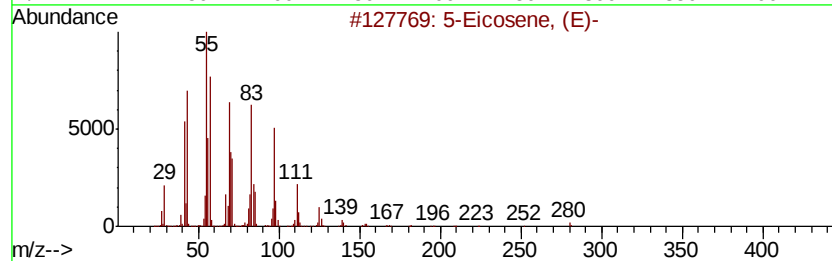
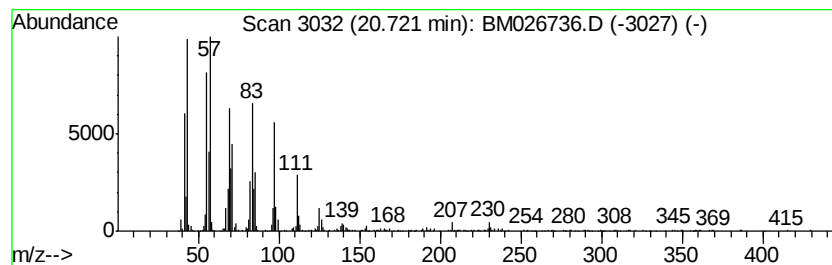
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.72	10.71 ng	696518	Chrysene-d12	20.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2		Cyclopentadecane	210	C15H30	000295-48-7	96
3		1-Docosene	308	C22H44	001599-67-3	94
4		1-Octadecene	252	C18H36	000112-88-9	92
5		Octacosanol	410	C28H58O	000557-61-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM071020\
 Data File : BM026736.D
 Acq On : 10 Jul 2020 16:24
 Operator : JU/CG
 Sample : L3217-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E3-41-BOILER-ROOM

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM070820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-eth...	6.96	4.7	ng	174750	1	7.32	742736	20.0
Sulfurous acid, p...	14.54	2.3	ng	120055	3	13.98	1047010	20.0
Undecane, 3,6-dim...	15.30	7.7	ng	403418	3	13.98	1047010	20.0
Tridecane, 5-propyl-	15.78	38.0	ng	2041300	4	16.74	1074660	20.0
Naphthalene, 1,2,...	15.87	3.4	ng	183172	4	16.74	1074660	20.0
4,4'-Dimethylbiph...	16.12	4.6	ng	247887	4	16.74	1074660	20.0
Hexadecane, 2,6,1...	16.60	28.2	ng	1513150	4	16.74	1074660	20.0
Hexacosane	17.20	8.5	ng	459103	4	16.74	1074660	20.0
n-Hexadecanoic acid	17.69	11.0	ng	589260	4	16.74	1074660	20.0
Phenanthrene, 2,3...	19.28	5.0	ng	325749	5	20.97	1301060	20.0
Butyl myristate	20.19	2.2	ng	141419	5	20.97	1301060	20.0
5-Eicosene, (E)-	20.72	10.7	ng	696518	5	20.97	1301060	20.0