

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
 Operator : HP/JU
 Sample : K3072-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 TP-4

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-BM053119.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	3.063	8	13	33	rVB	2229601	2596401	18.88%	2.414%
2	5.416	406	413	424	rBV	5764743	8400050	61.08%	7.810%
3	6.992	673	681	697	rBV	5666422	9474862	68.89%	8.809%
4	7.363	736	744	753	rBV	5889313	9355932	68.03%	8.698%
5	7.828	817	823	830	rBV	1077896	1699977	12.36%	1.580%
6	8.151	870	878	885	rBV	4104049	6670395	48.50%	6.202%
7	8.992	1013	1021	1033	rBV	3374211	5622366	40.88%	5.227%
8	10.622	1291	1298	1305	rVB	1376130	2298429	16.71%	2.137%
9	13.086	1710	1717	1727	rVB	7571078	11151198	81.08%	10.367%
10	13.921	1853	1859	1867	rBV	921629	1268814	9.23%	1.180%
11	14.463	1940	1951	1960	rBV2	1842827	2889149	21.01%	2.686%
12	15.957	2197	2205	2221	rBV	6188274	8714914	63.37%	8.102%
13	17.209	2412	2418	2426	rVB2	2286487	3262798	23.72%	3.033%
14	18.109	2565	2571	2580	rBV	1324496	1823161	13.26%	1.695%
15	18.751	2676	2680	2690	rVB2	174453	245592	1.79%	0.228%
16	19.268	2765	2768	2777	rVB8	87608	176764	1.29%	0.164%
17	19.386	2783	2788	2797	rBV	574484	853365	6.20%	0.793%
18	19.656	2830	2834	2839	rVB	127889	154333	1.12%	0.143%
19	19.839	2859	2865	2870	rBV	11120153	13752970	100.00%	12.786%
20	20.150	2915	2918	2922	rVB	177747	177330	1.29%	0.165%
21	20.521	2976	2981	2988	rVB3	85763	177976	1.29%	0.165%
22	20.645	2998	3002	3006	rBV2	306182	351135	2.55%	0.326%
23	21.103	3076	3080	3091	rBV2	1492435	1917179	13.94%	1.782%
24	21.386	3123	3128	3134	rBV	2756242	3464329	25.19%	3.221%
25	21.544	3152	3155	3160	rVB	625906	677349	4.93%	0.630%
26	21.991	3227	3231	3237	rBV	671806	878528	6.39%	0.817%
27	22.468	3308	3312	3318	rBV	647075	869119	6.32%	0.808%
28	22.603	3330	3335	3343	rVB	1120727	1564237	11.37%	1.454%
29	22.997	3397	3402	3407	rBV	600685	873450	6.35%	0.812%
30	23.586	3497	3502	3509	rVB	503549	838500	6.10%	0.780%
31	23.680	3511	3518	3527	rVB	1769918	3460780	25.16%	3.218%
32	24.262	3611	3617	3623	rVB	319443	622085	4.52%	0.578%
33	24.344	3627	3631	3644	rVB3	189490	387000	2.81%	0.360%
34	26.850	4051	4057	4068	rVB2	316173	889989	6.47%	0.827%

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TP-4

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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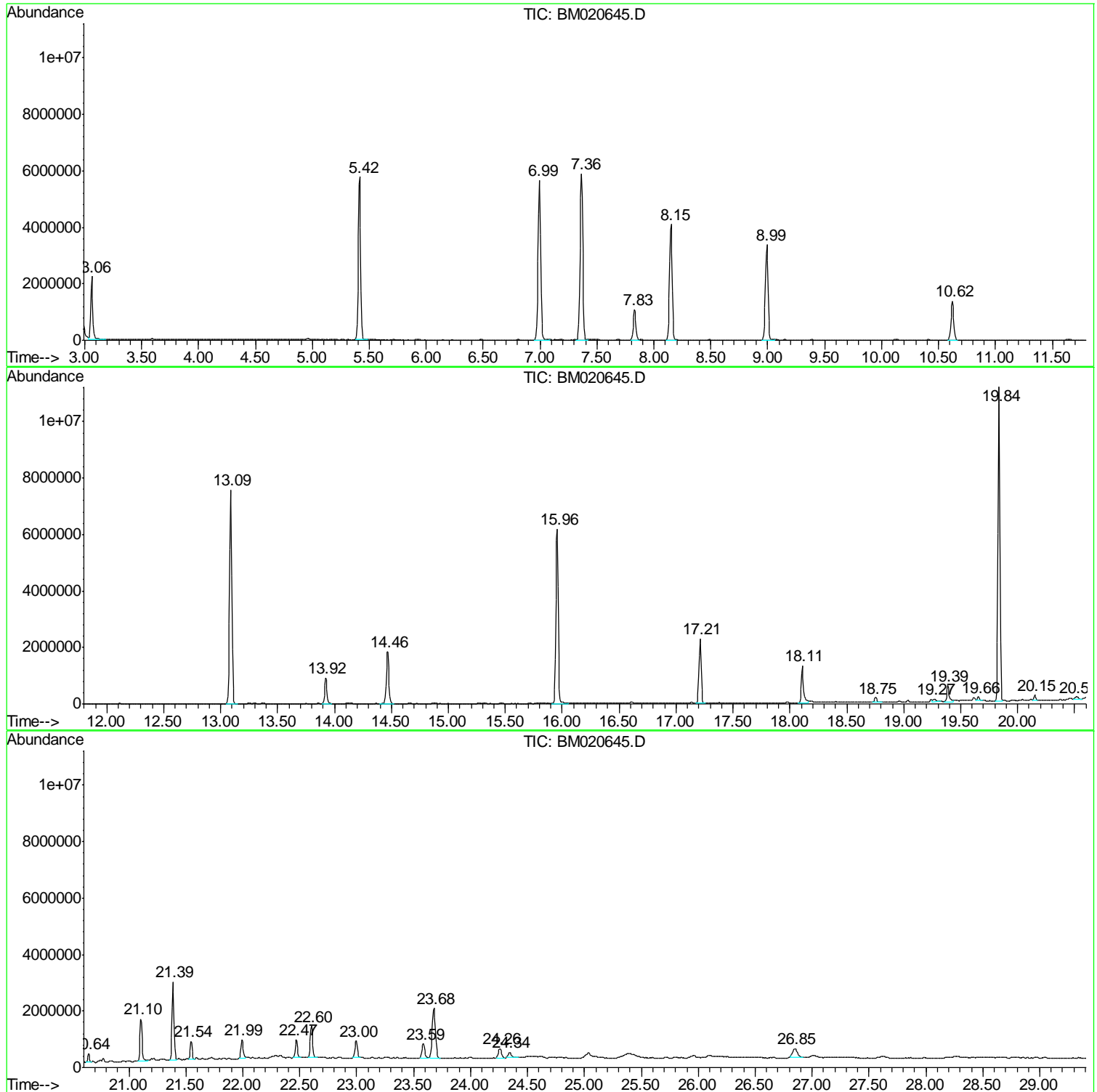
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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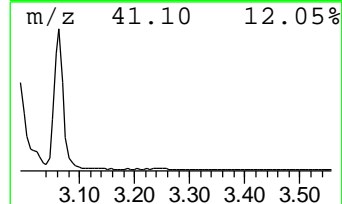
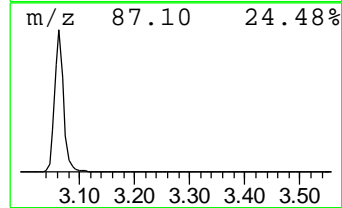
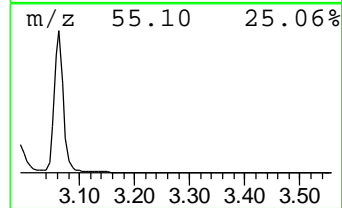
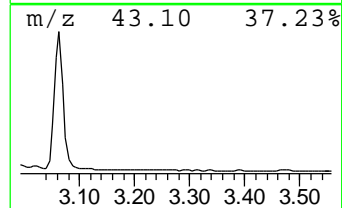
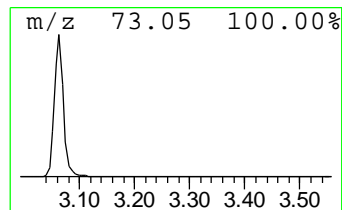
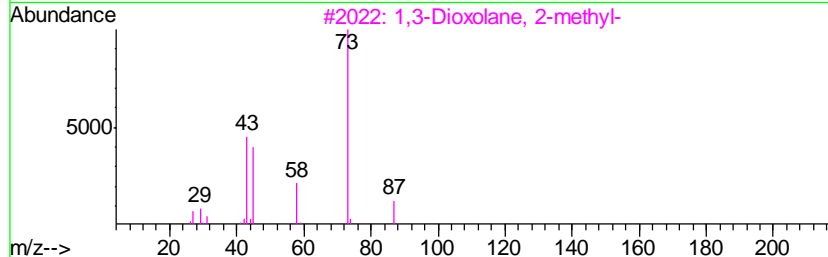
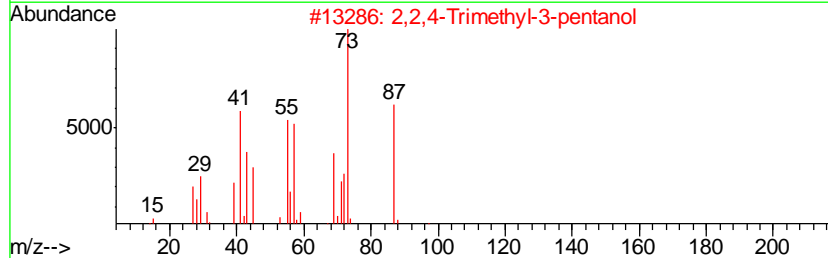
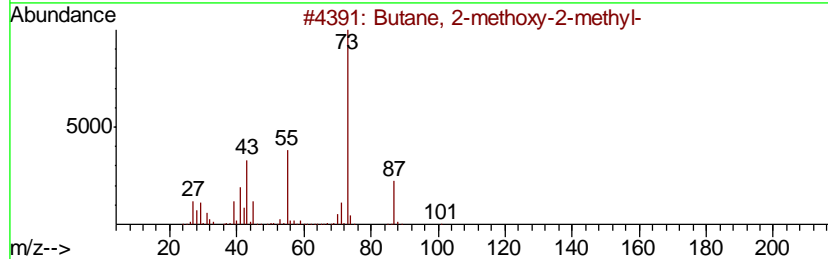
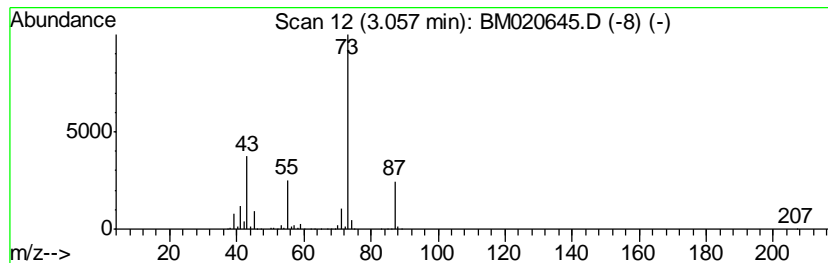
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	30.55 ng	2596400	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	56
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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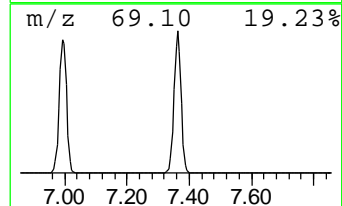
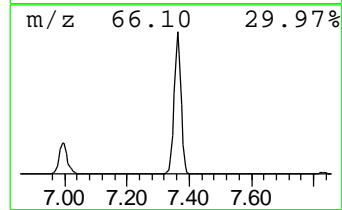
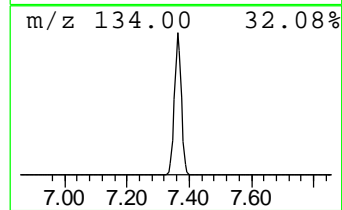
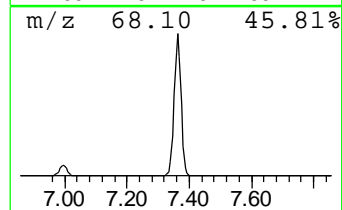
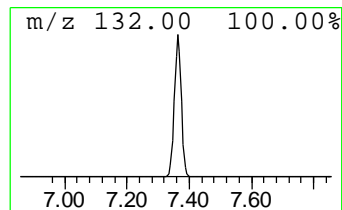
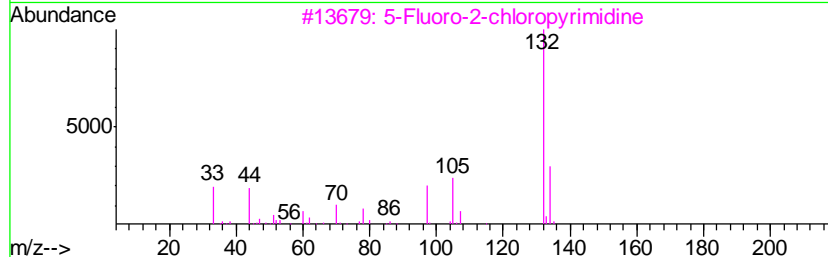
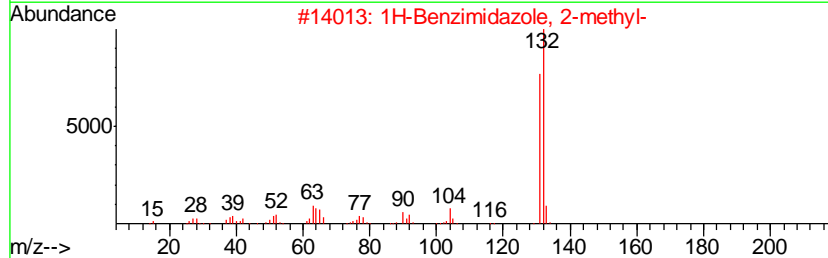
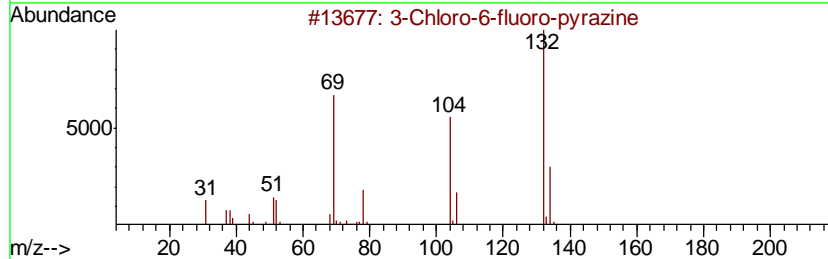
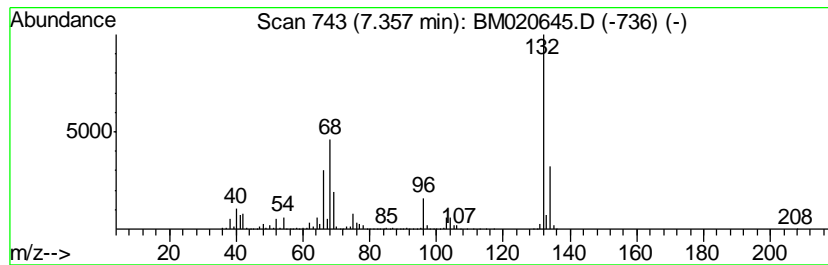
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 2 unknown7.36 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	110.07 ng	9355930	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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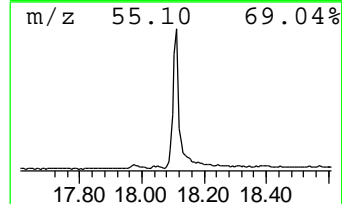
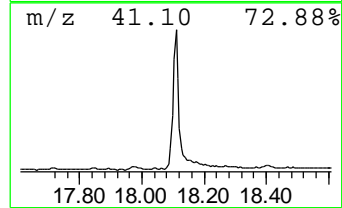
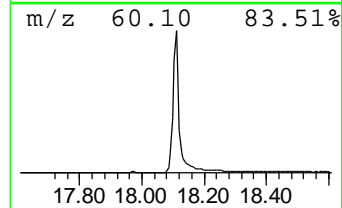
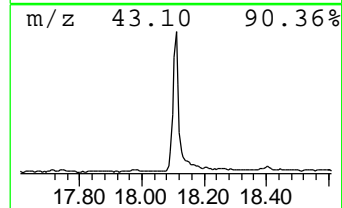
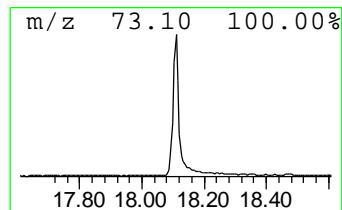
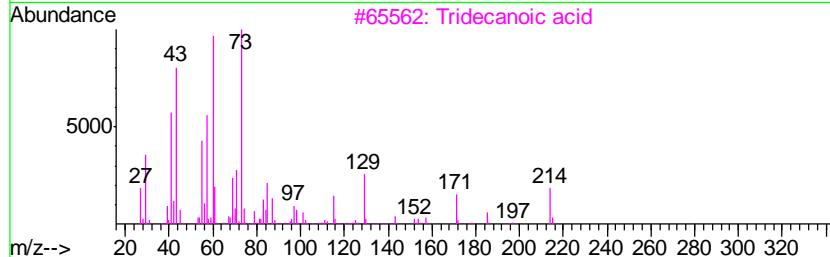
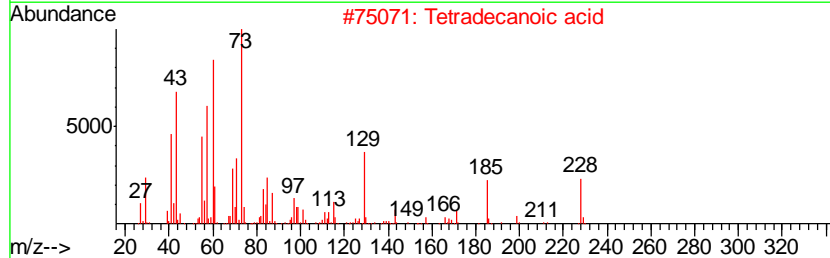
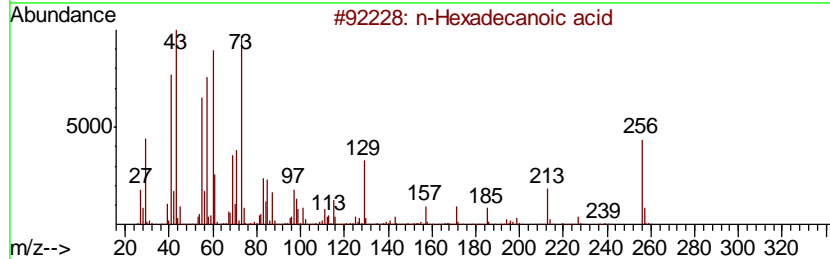
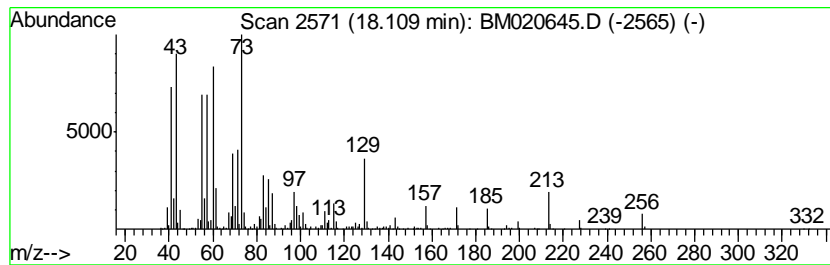
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	11.18 ng	1823160	Phenanthrene-d10	17.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3		Tridecanoic acid	214	C13H26O2	000638-53-9	76
4		Nonanoic acid	158	C9H18O2	000112-05-0	38
5		Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	30



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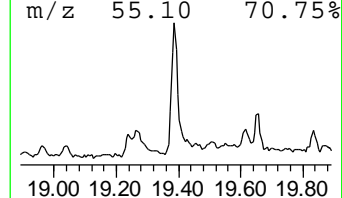
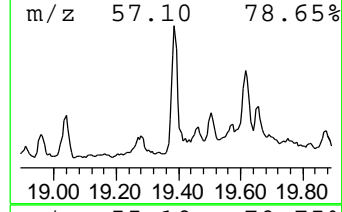
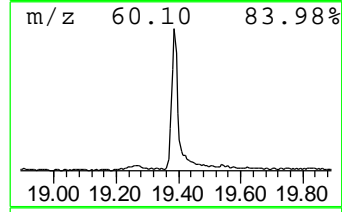
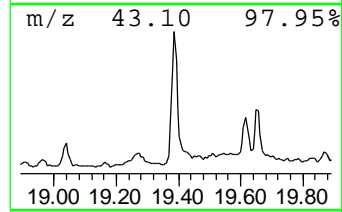
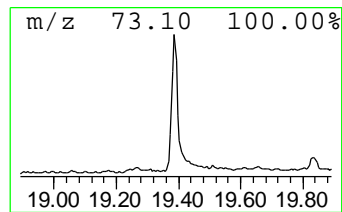
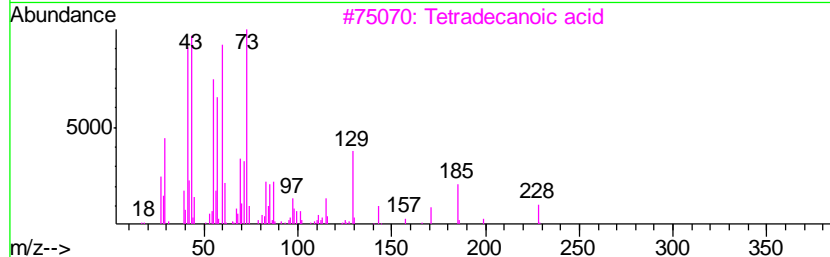
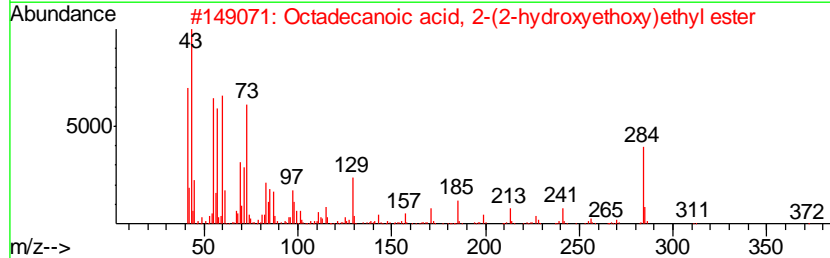
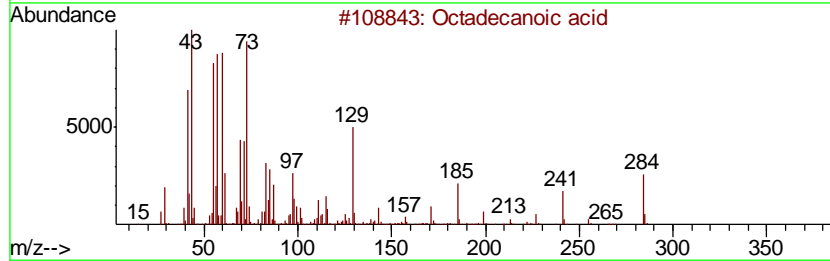
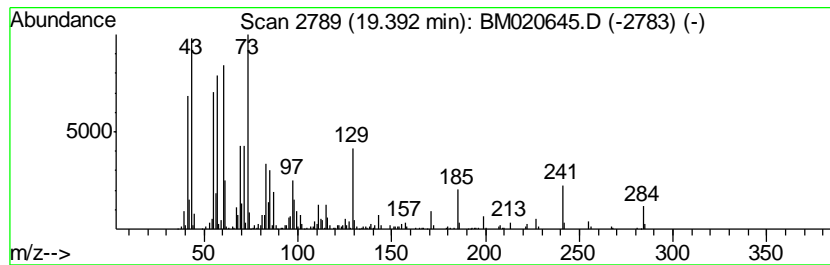
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TIC Library : C:\DATABASE\NIST02.L
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 Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	4.93 ng	853365	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		n-Decanoic acid	172	C10H20O2	000334-48-5	68
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	68



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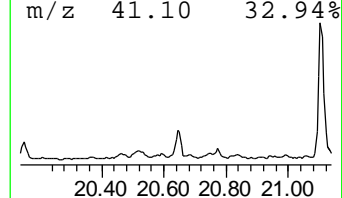
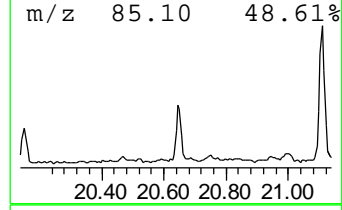
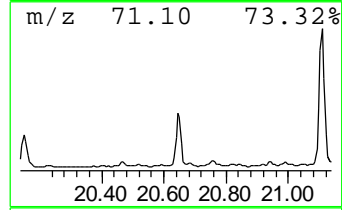
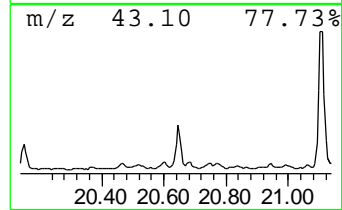
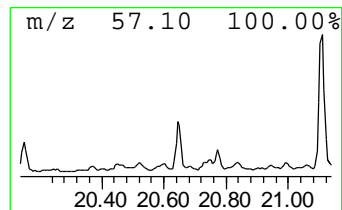
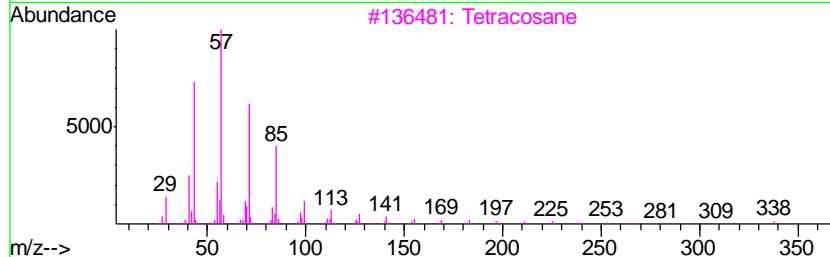
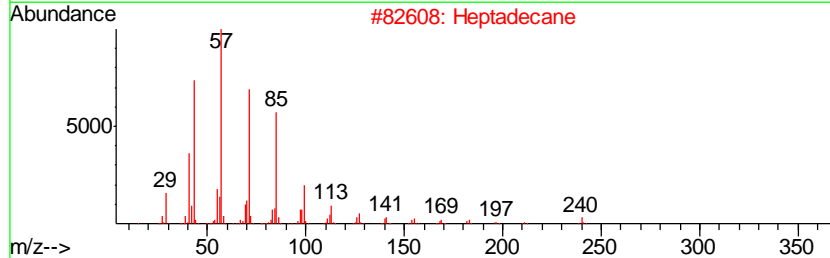
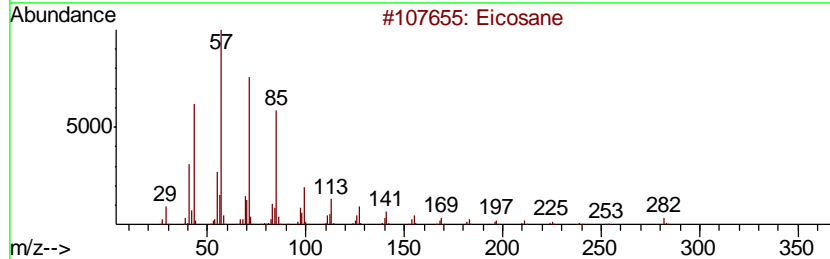
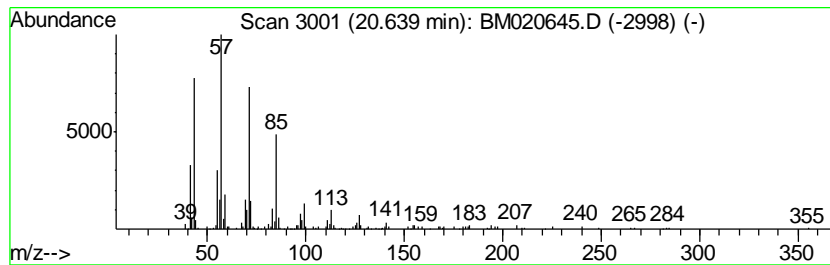
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 Peak Number 6 Eicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.64	2.03 ng	351135	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Heptadecane	240	C17H36	000629-78-7	93
3		Tetracosane	338	C24H50	000646-31-1	90
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	90
5		Octadecane	254	C18H38	000593-45-3	87



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 Misc :
 ALS Vial : 3 Sample Multiplier: 1

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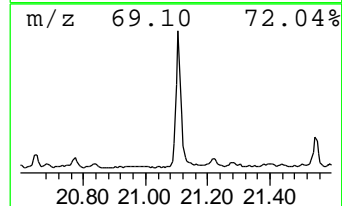
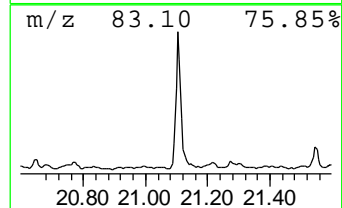
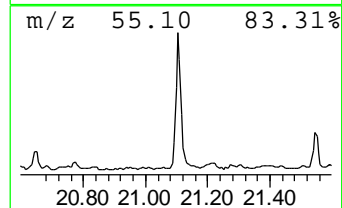
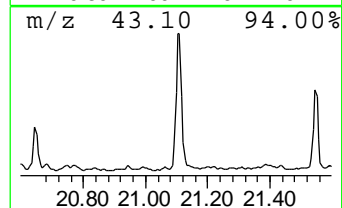
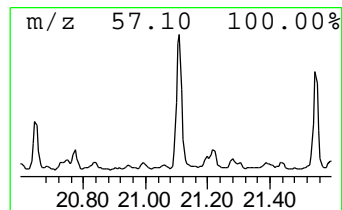
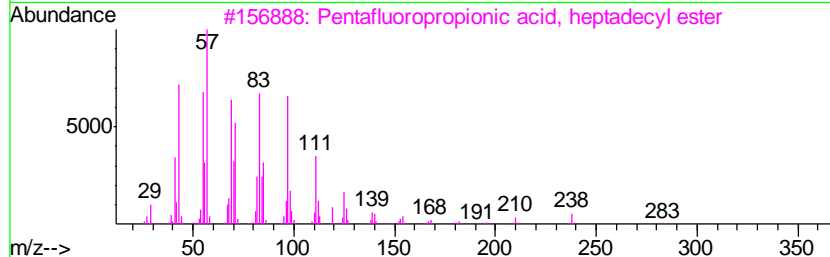
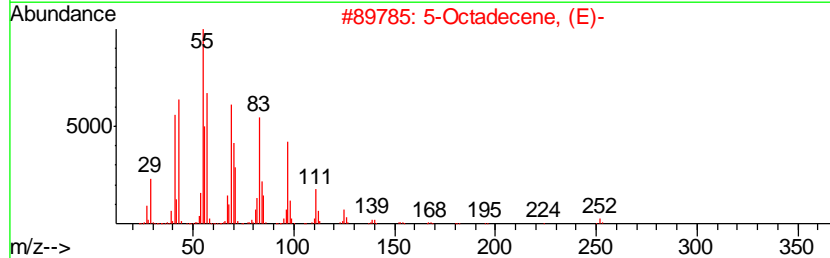
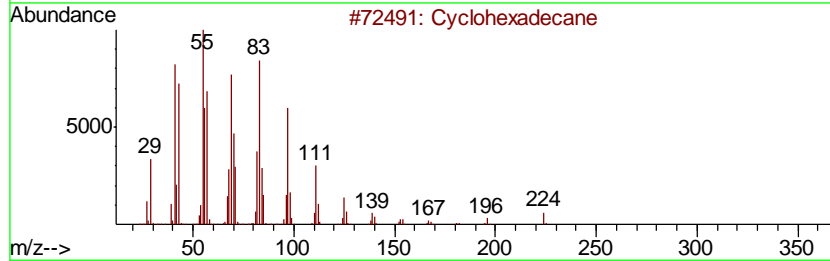
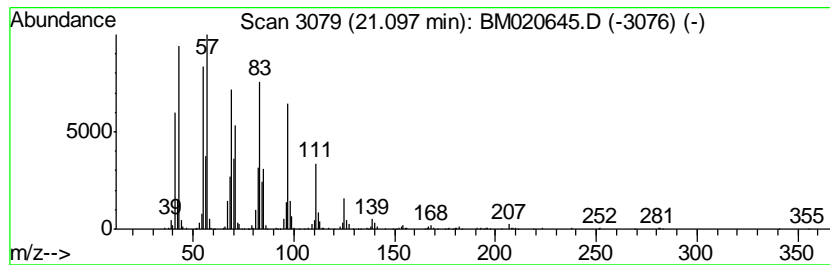
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 7 Cyclohexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.10	11.07 ng	1917180	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	98
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	95
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	91
4		2-Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	87
5		17-Pentatriacontene	491	C35H70	006971-40-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
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 Sample : K3072-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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 ClientSampled :
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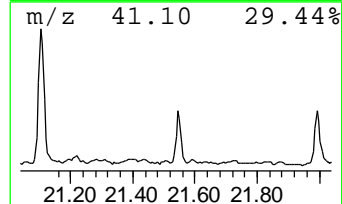
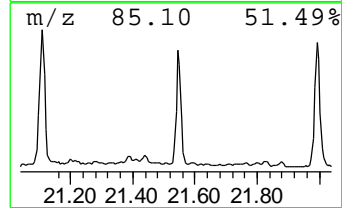
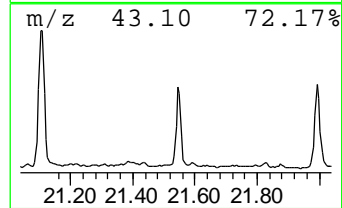
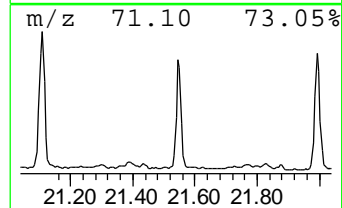
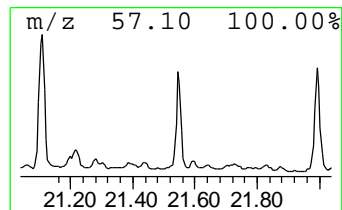
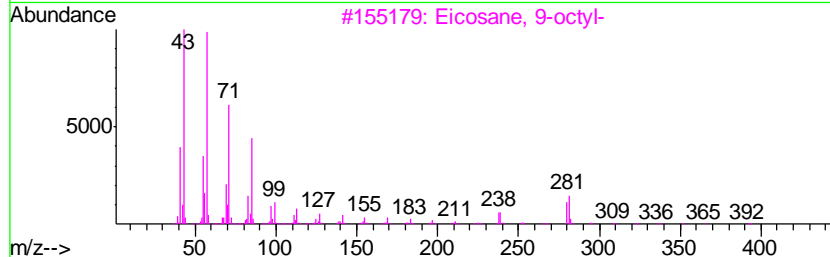
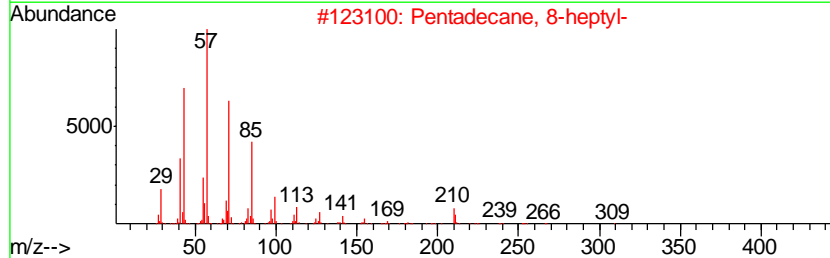
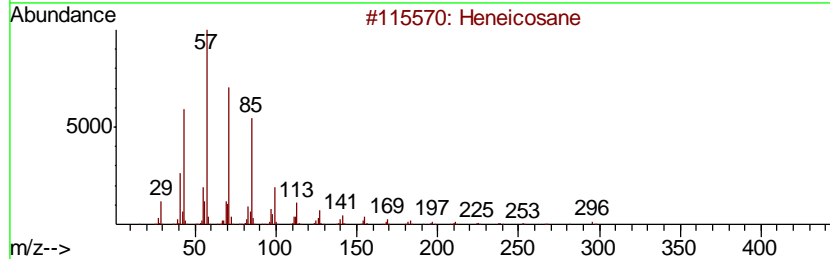
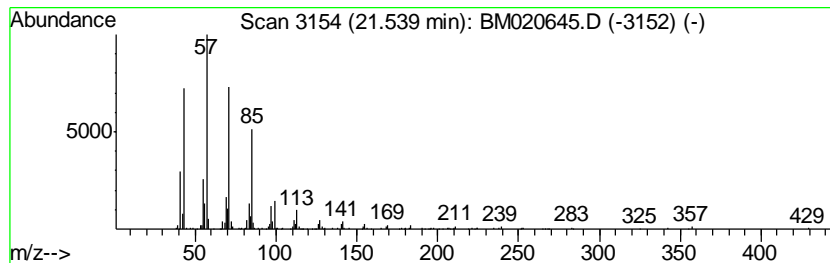
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 8 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.54	3.91 ng	677349	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Eicosane, 9-octyl-	394	C28H58	013475-77-9	91
4		Octacosane	394	C28H58	000630-02-4	90
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
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Instrument :
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 ClientSampled :
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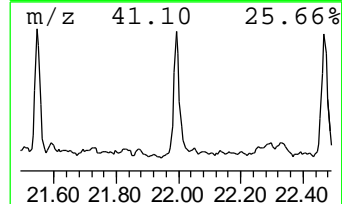
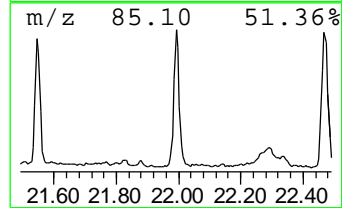
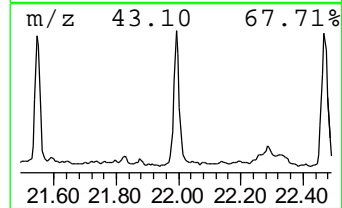
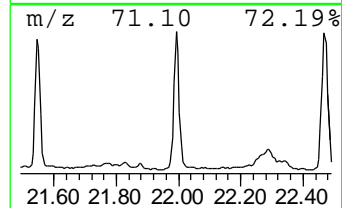
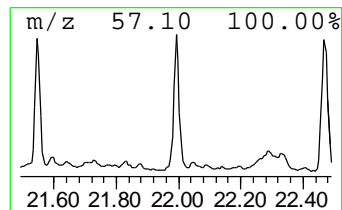
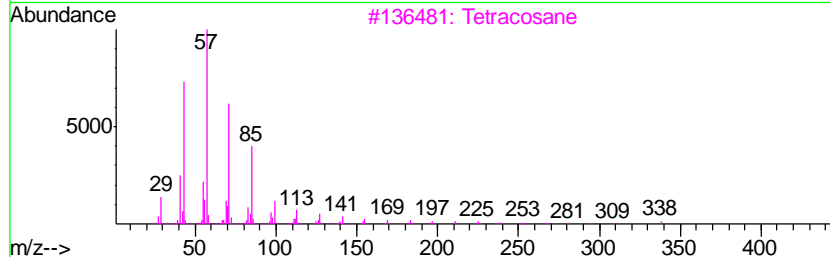
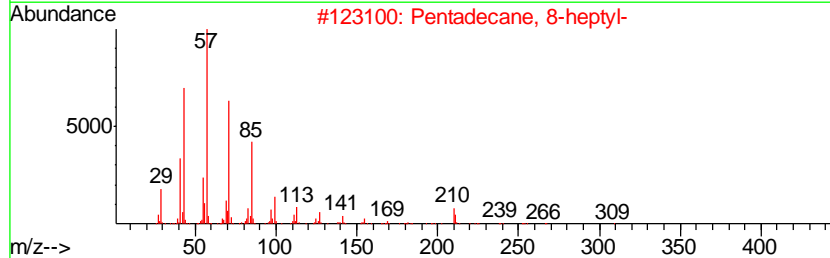
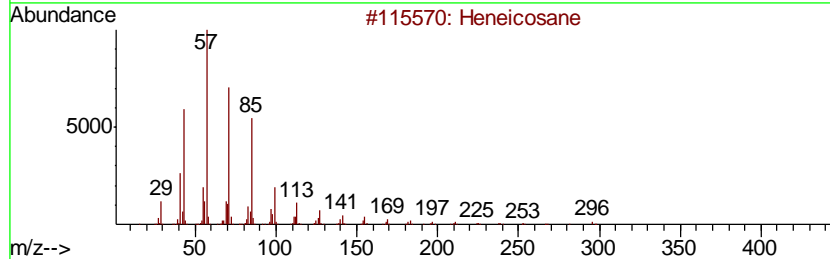
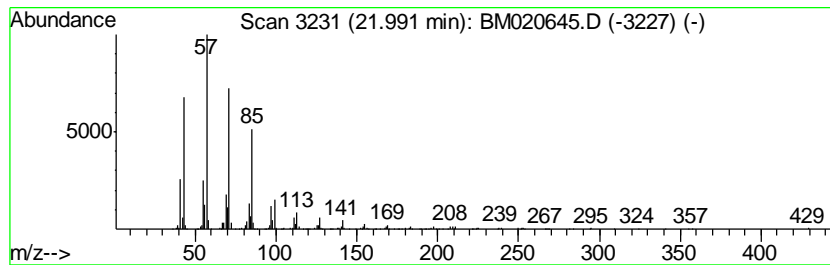
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 9 Pentadecane, 8-heptyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.99	5.07 ng	878528	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
5		Tricosane	324	C23H48	000638-67-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
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 Acq On : 01 Jun 2019 11:56
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Instrument :
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 ClientSampled :
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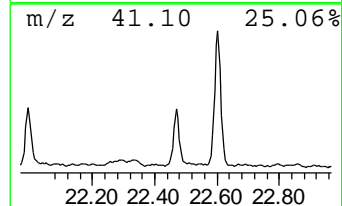
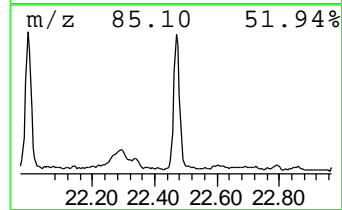
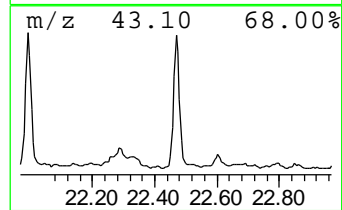
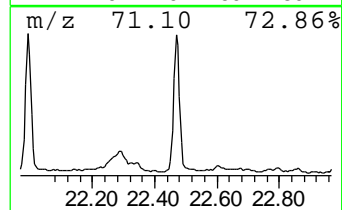
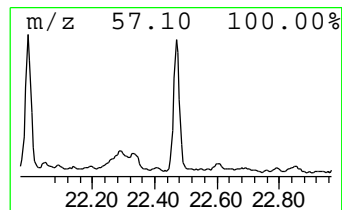
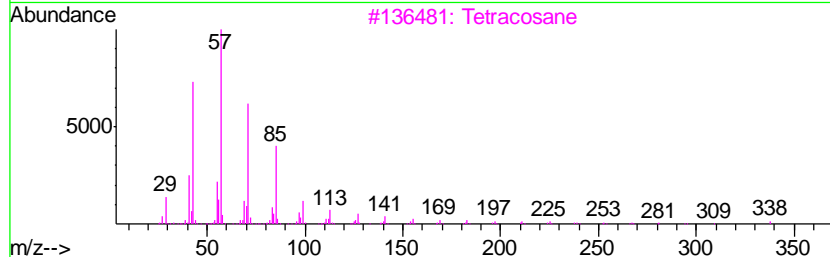
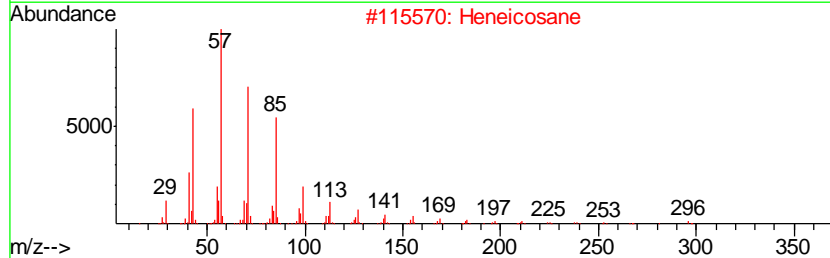
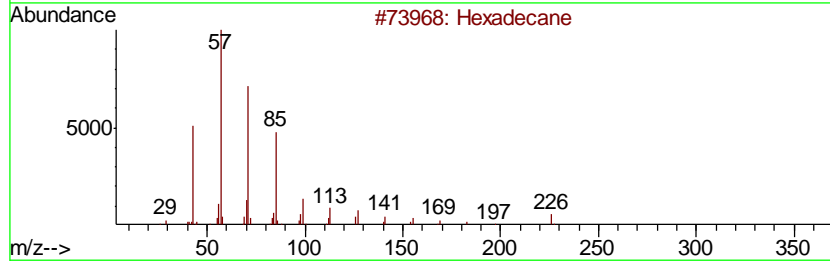
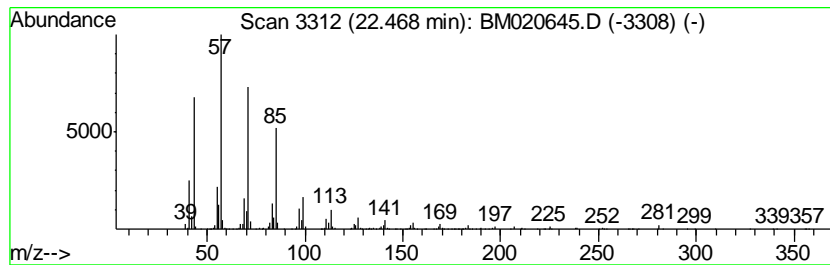
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 10 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.47	5.02 ng	869119	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
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Instrument :
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 ClientSampled :
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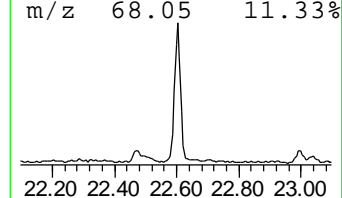
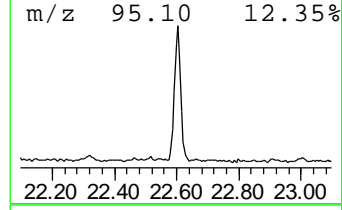
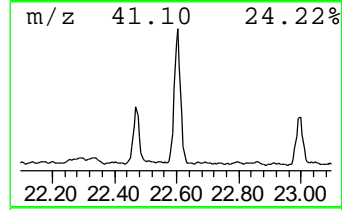
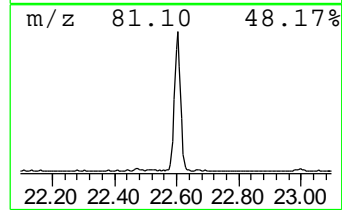
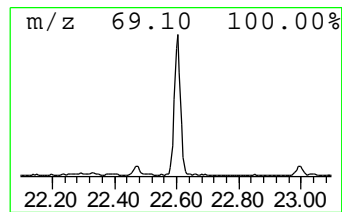
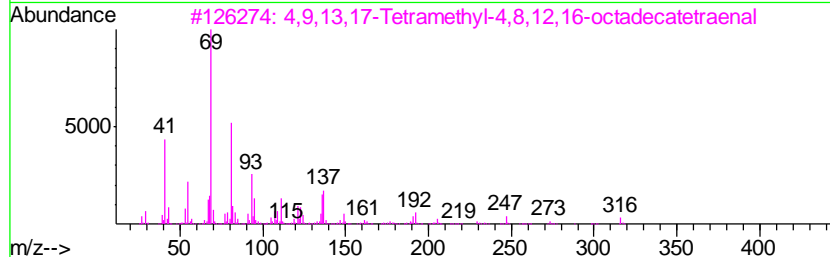
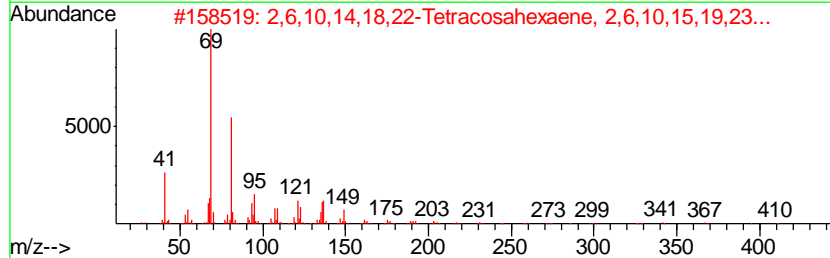
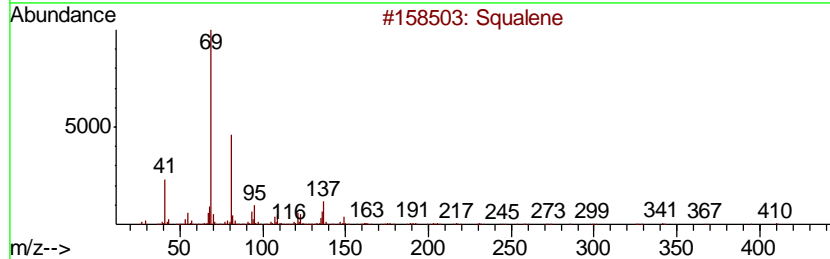
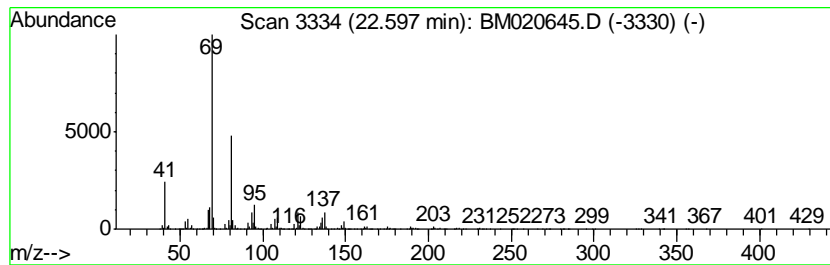
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TIC Library : C:\DATABASE\NIST02.L
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 Peak Number 11 Squalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.60	9.04 ng	1564240	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	91
2		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	91
3		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	74
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5		Farnesol isomer a	222	C15H26O	1000108-92-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
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Instrument :
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 ClientSampled :
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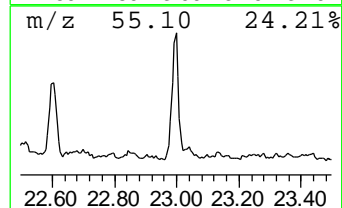
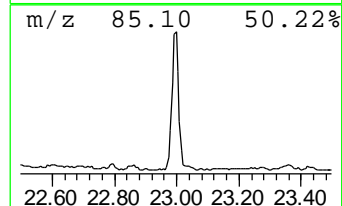
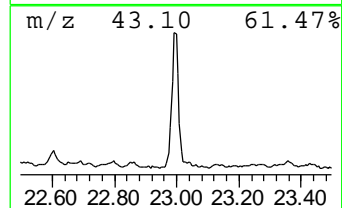
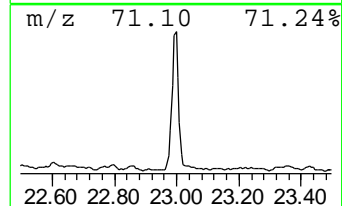
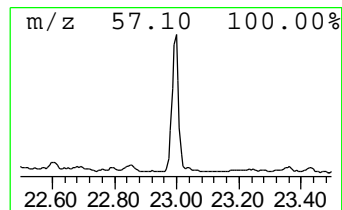
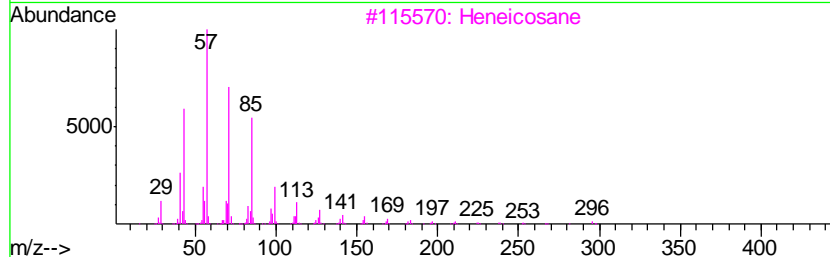
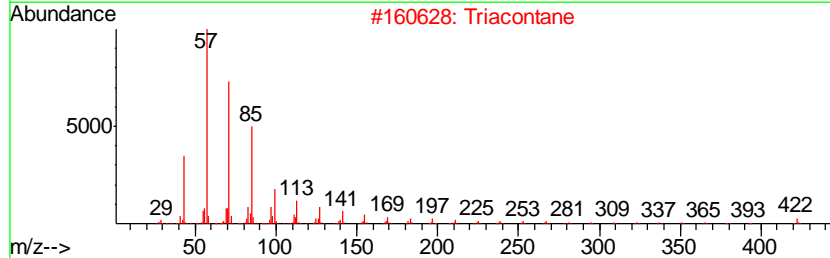
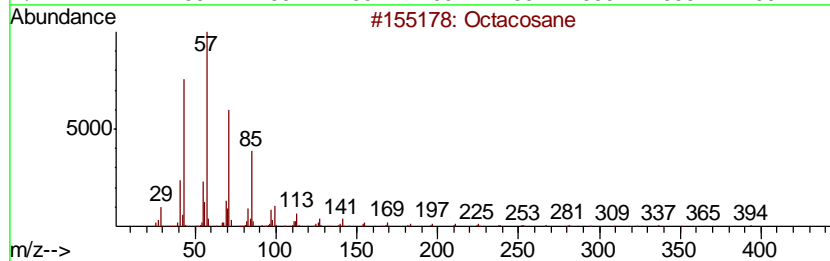
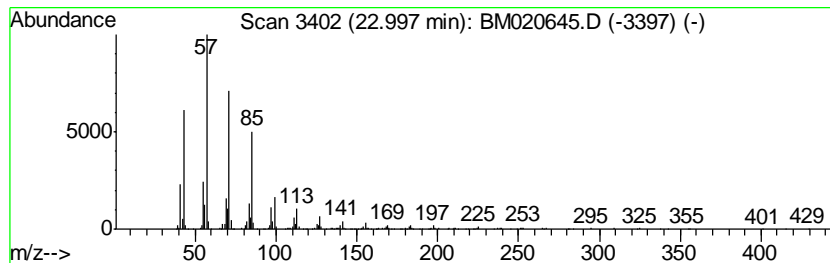
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 12 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.00	5.05 ng	873450	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		triacontane	422	C30H62	000638-68-6	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



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Instrument :
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 ClientSampled :
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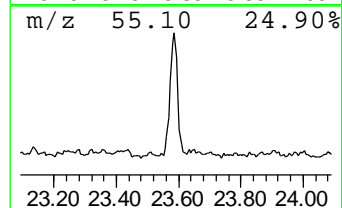
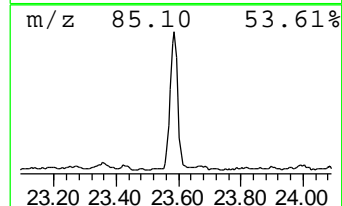
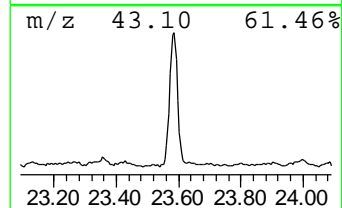
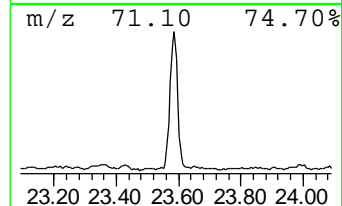
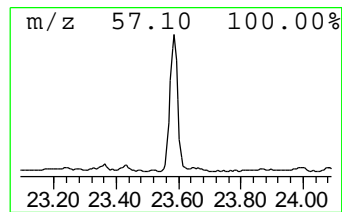
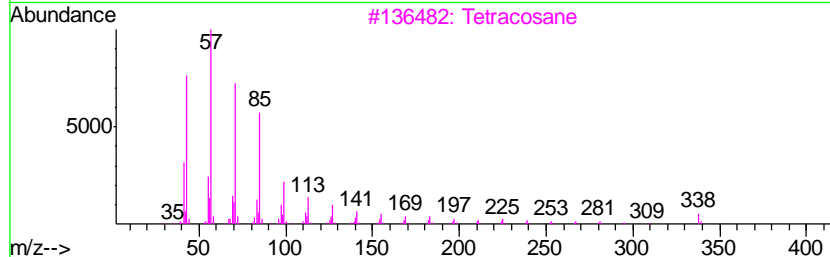
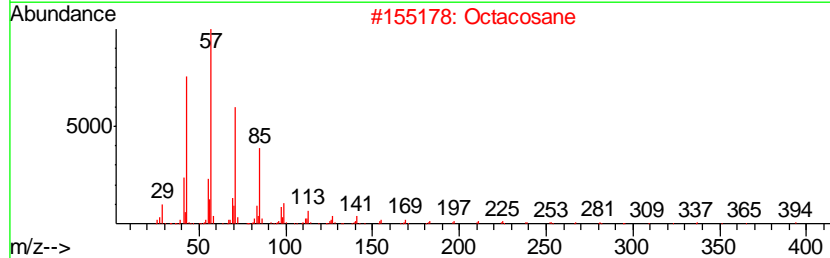
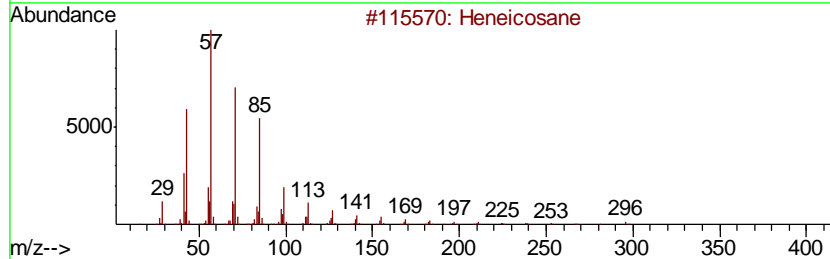
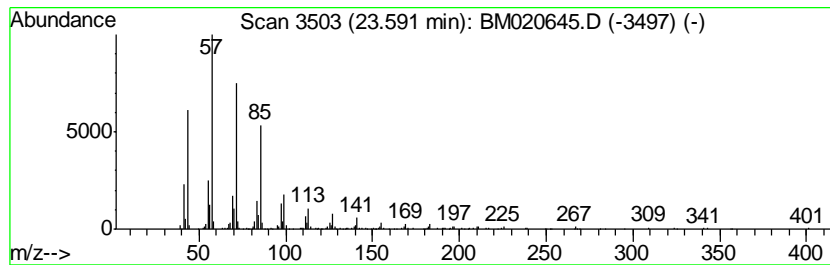
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TIC Library : C:\DATABASE\NIST02.L
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 Peak Number 13 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.59	4.85 ng	838500	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
 Operator : HP/JU
 Sample : K3072-04
 Misc :
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Instrument :
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 ClientSampled :
 TP-4

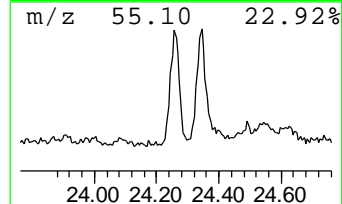
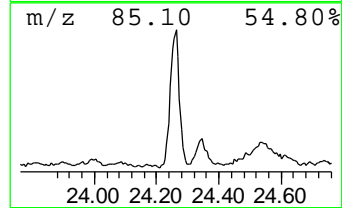
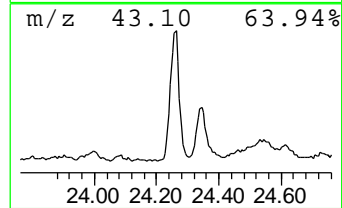
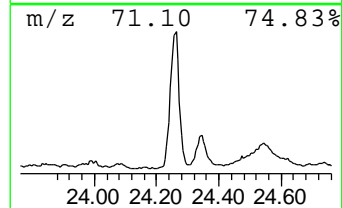
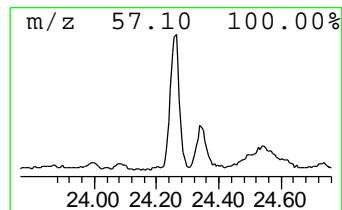
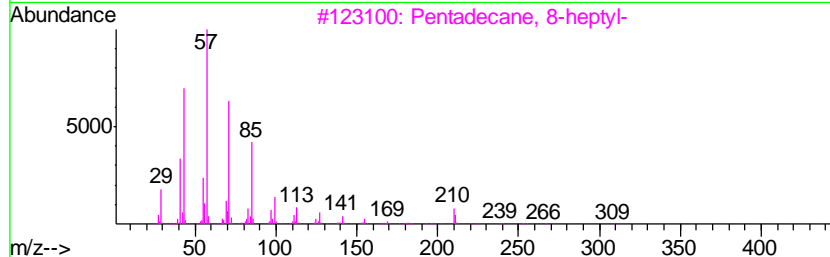
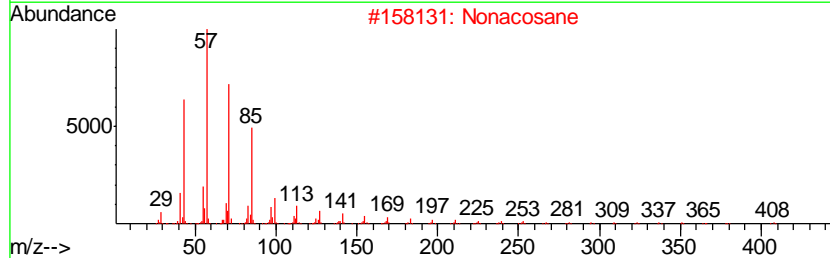
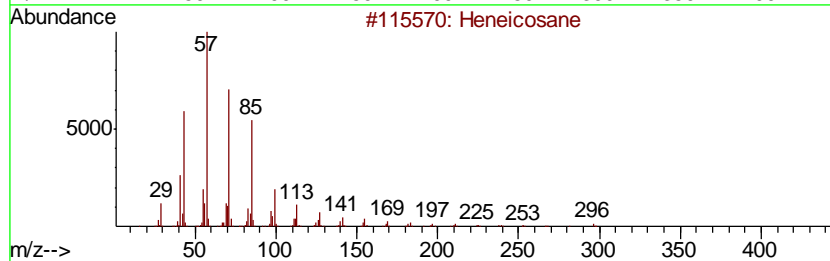
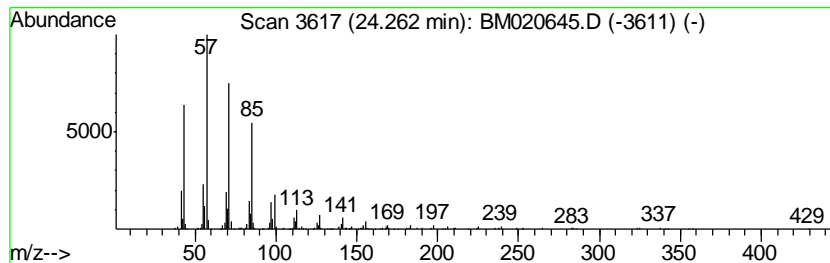
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 14 Nonacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.26	3.60 ng	622085	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Nonacosane	408	C29H60	000630-03-5	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Octadecane	254	C18H38	000593-45-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
 Operator : HP/JU
 Sample : K3072-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 TP-4

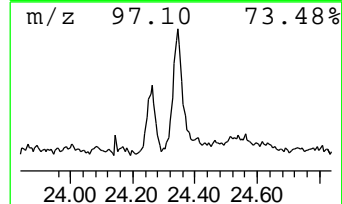
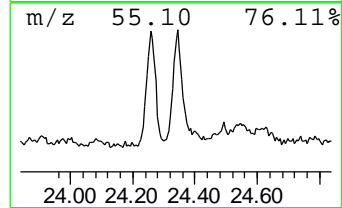
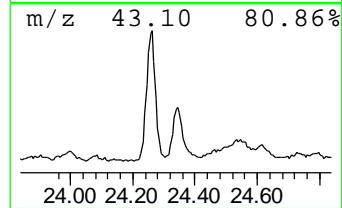
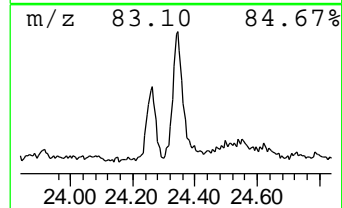
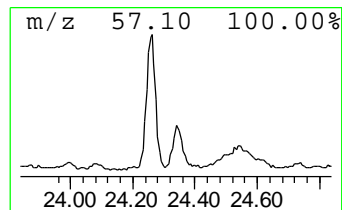
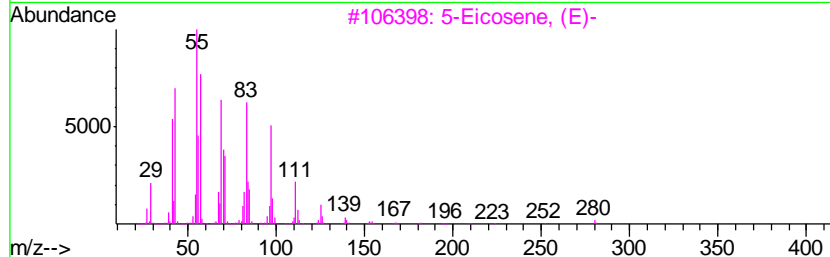
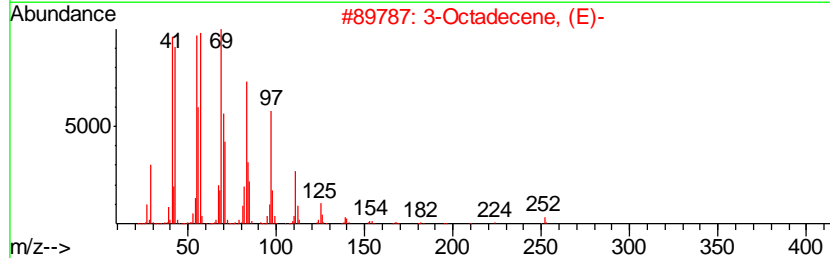
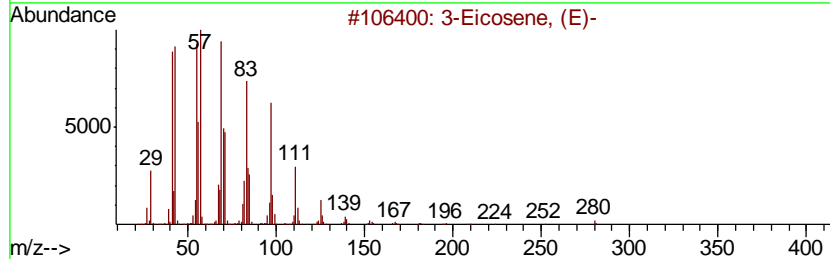
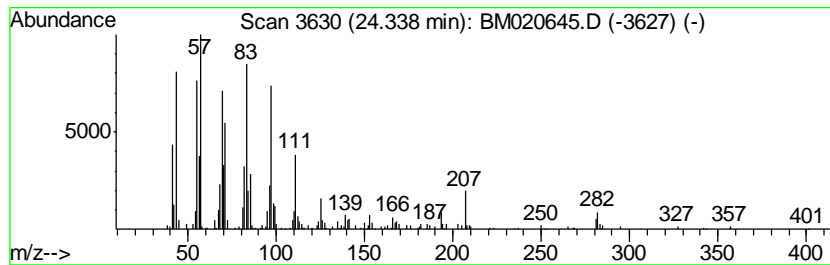
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 15 3-Eicosene, (E)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.34	2.24 ng	387000	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
2		3-Octadecene, (E)-	252	C18H36	007206-19-1	86
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	78
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	76
5		17-Pentatriacontene	491	C35H70	006971-40-0	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
 Operator : HP/JU
 Sample : K3072-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 TP-4

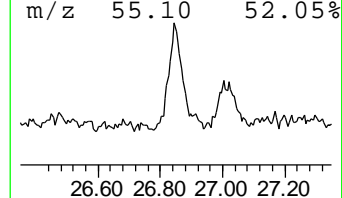
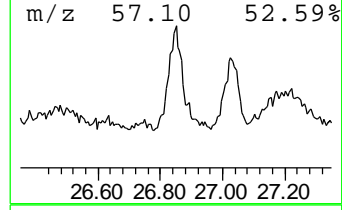
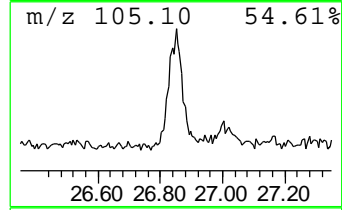
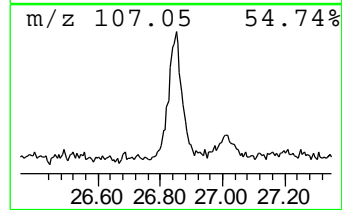
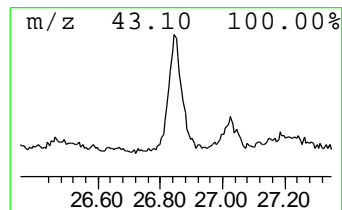
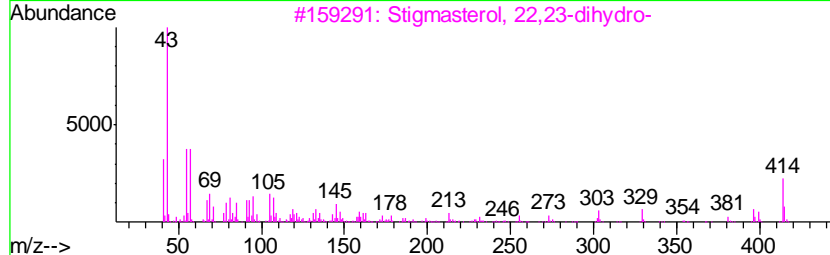
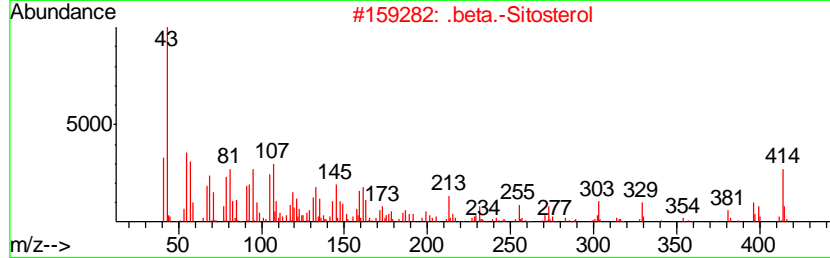
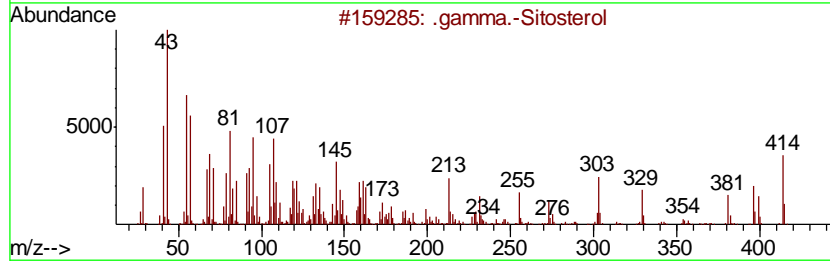
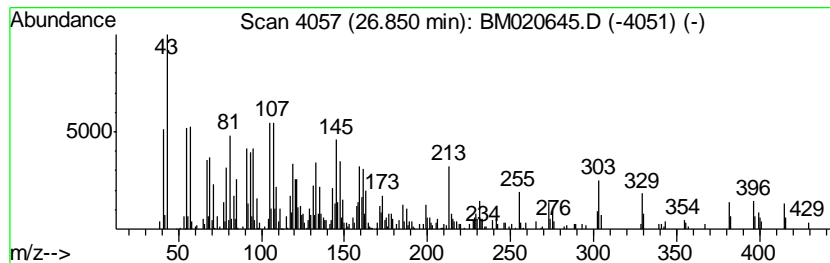
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 16 .gamma.-Sitosterol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.85	5.14 ng	889989	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	98
3		Stigmasterol, 22,23-dihydro-	414	C29H50O	1000214-20-7	97
4		Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	95
5		.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060119\
 Data File : BM020645.D
 Acq On : 01 Jun 2019 11:56
 Operator : HP/JU
 Sample : K3072-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM053119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.06	30.6	ng	2596400	1	7.83	1699980	20.0
unknown7.36	7.36	110.1	ng	9355930	1	7.83	1699980	20.0
n-Hexadecanoic acid	18.11	11.2	ng	1823160	4	17.21	3262800	20.0
Octadecanoic acid	19.39	4.9	ng	853365	5	21.39	3464330	20.0
Eicosane	20.64	2.0	ng	351135	5	21.39	3464330	20.0
Cyclohexadecane	21.10	11.1	ng	1917180	5	21.39	3464330	20.0
Heneicosane	21.54	3.9	ng	677349	5	21.39	3464330	20.0
Pentadecane, 8-he...	21.99	5.1	ng	878528	5	21.39	3464330	20.0
Hexadecane	22.47	5.0	ng	869119	5	21.39	3464330	20.0
Squalene	22.60	9.0	ng	1564240	6	23.68	3460780	20.0
Octacosane	23.00	5.0	ng	873450	6	23.68	3460780	20.0
Tetracosane	23.59	4.8	ng	838500	6	23.68	3460780	20.0
Nonacosane	24.26	3.6	ng	622085	6	23.68	3460780	20.0
3-Eicosene, (E)-	24.34	2.2	ng	387000	6	23.68	3460780	20.0
.gamma.-Sitosterol	26.85	5.1	ng	889989	6	23.68	3460780	20.0