

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
 Data File : BM020642.D
 Acq On : 31 May 2019 23:12
 Operator : JU/SJ
 Sample : K3072-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-1

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	3.069	9	14	28	rVB	2140344	2508588	20.09%	2.188%
2	5.416	406	413	426	rBV	6767114	9712525	77.77%	8.470%
3	6.998	673	682	694	rBV	6656345	11163099	89.38%	9.735%
4	7.369	736	745	753	rBV	6972672	11073669	88.66%	9.657%
5	7.834	817	824	831	rBV	1302501	2004065	16.05%	1.748%
6	8.151	869	878	886	rBV	5047086	8108387	64.92%	7.071%
7	8.998	1013	1022	1032	rBV	4090576	6856256	54.90%	5.979%
8	10.628	1288	1299	1307	rBV	1608454	2730738	21.86%	2.381%
9	13.092	1710	1718	1725	rBV	8637385	12489337	100.00%	10.891%
10	13.927	1854	1860	1869	rBV	1184429	1636084	13.10%	1.427%
11	14.469	1945	1952	1963	rVB2	2015189	3092317	24.76%	2.697%
12	15.957	2198	2205	2222	rBV	5441010	7739682	61.97%	6.749%
13	17.210	2409	2418	2425	rBV2	2044875	3016541	24.15%	2.631%
14	18.110	2565	2571	2581	rBV	765554	1130329	9.05%	0.986%
15	19.386	2784	2788	2797	rBV	345754	544938	4.36%	0.475%
16	19.839	2859	2865	2870	rBV	9887302	11532097	92.34%	10.057%
17	21.109	3076	3081	3091	rBV	1114902	1503053	12.03%	1.311%
18	21.392	3123	3129	3135	rBV	2278440	3000463	24.02%	2.617%
19	21.550	3152	3156	3161	rVB	630638	679642	5.44%	0.593%
20	21.997	3227	3232	3238	rBV	1080695	1376618	11.02%	1.200%
21	22.474	3308	3313	3318	rBV	1585592	1974625	15.81%	1.722%
22	23.003	3397	3403	3416	rVB	1452937	2347245	18.79%	2.047%
23	23.592	3497	3503	3510	rBV	1304709	2236500	17.91%	1.950%
24	23.686	3512	3519	3531	rVB	1574361	3149665	25.22%	2.747%
25	24.268	3612	3618	3626	rVV	937423	1730477	13.86%	1.509%
26	24.350	3628	3632	3645	rVB2	172101	365044	2.92%	0.318%
27	25.050	3745	3751	3760	rVB2	447420	969097	7.76%	0.845%

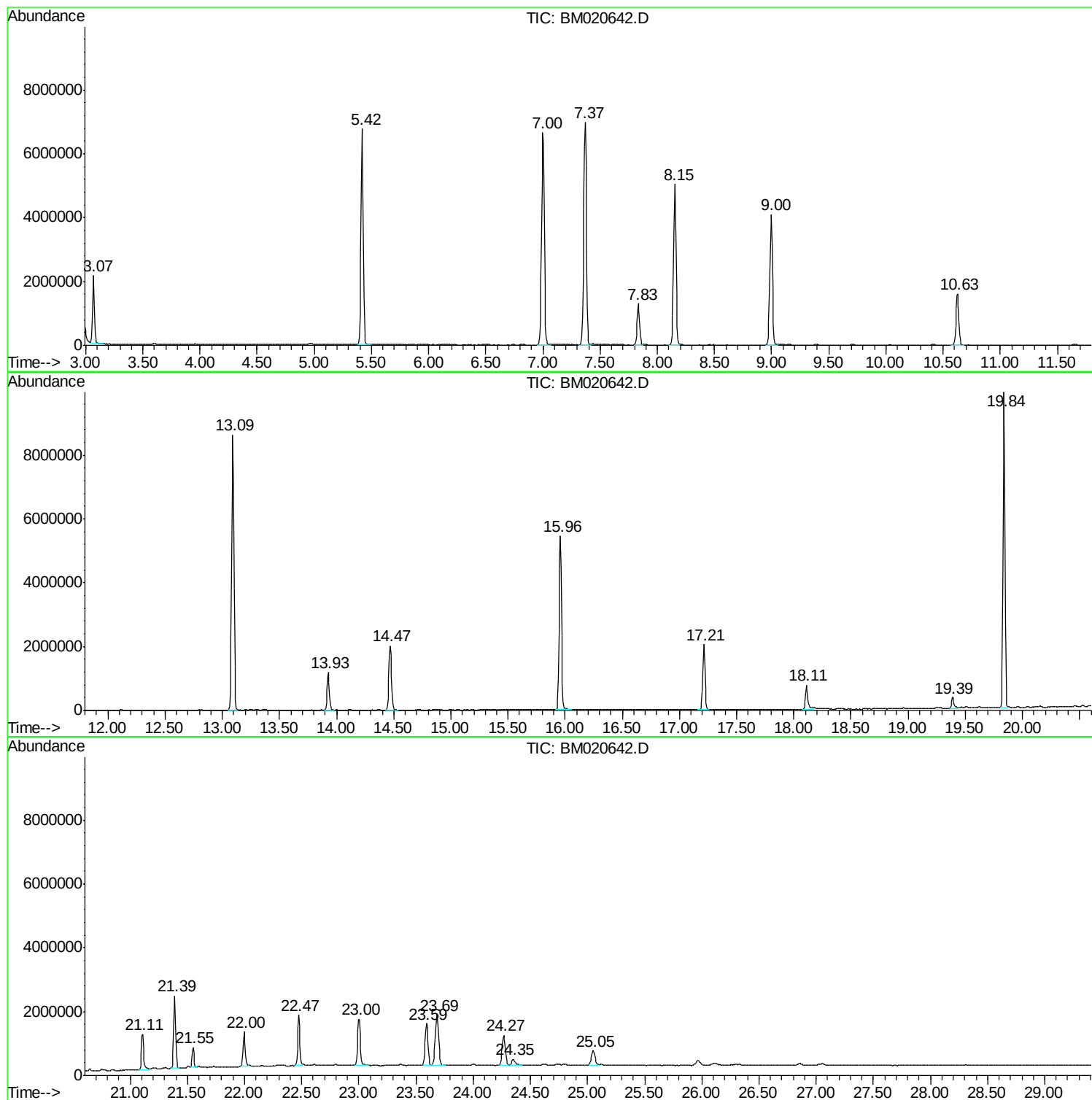
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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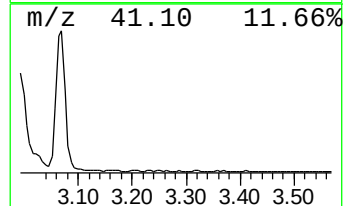
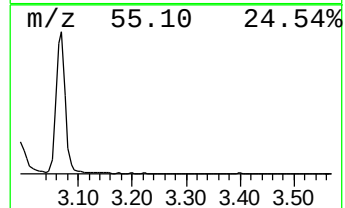
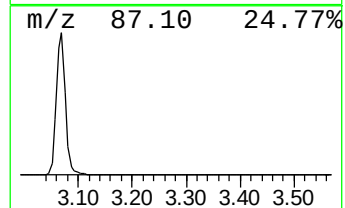
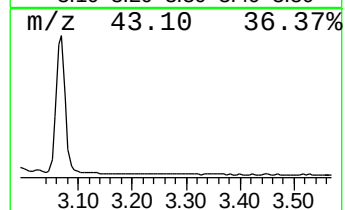
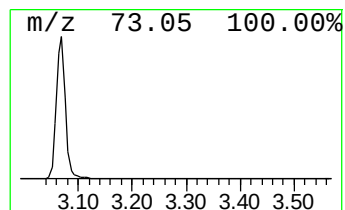
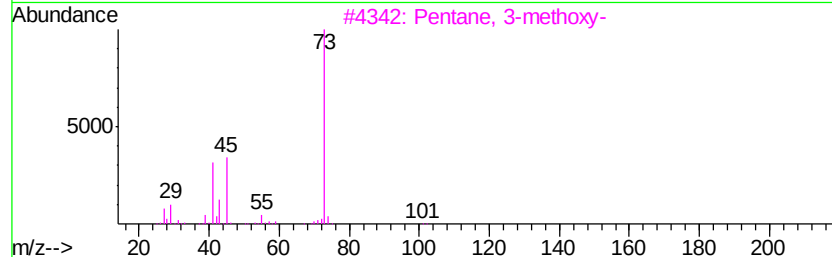
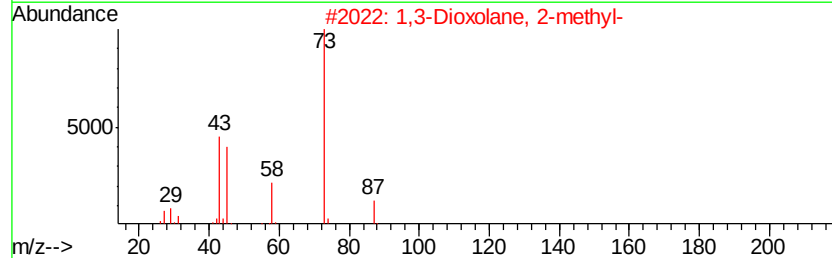
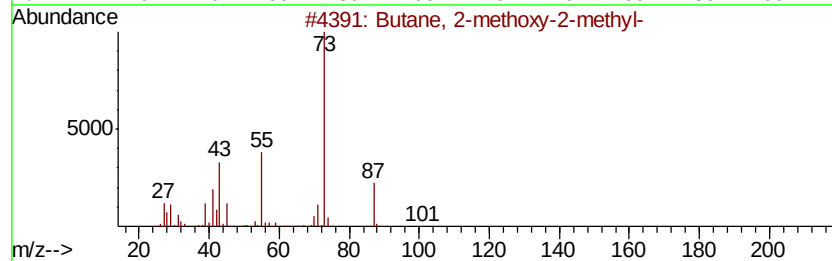
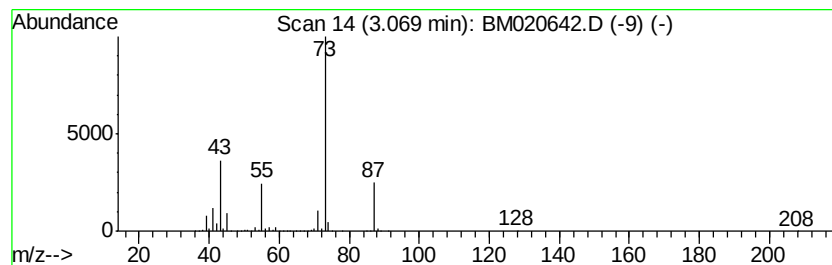
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
3.07	25.04 ng	2508590	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		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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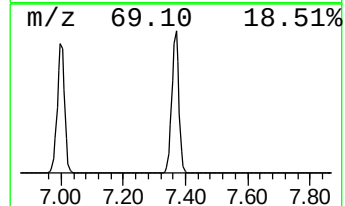
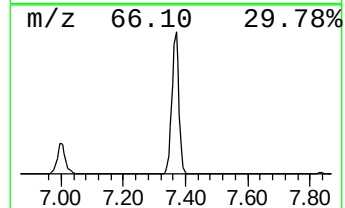
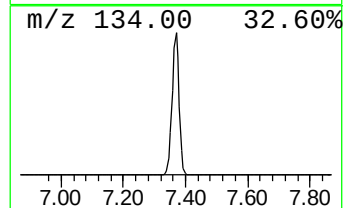
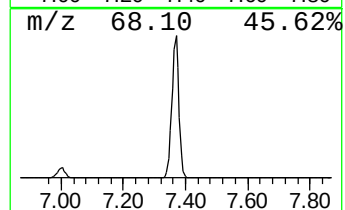
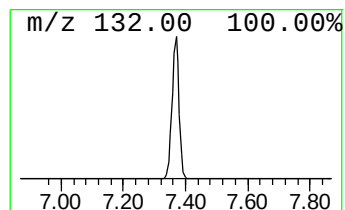
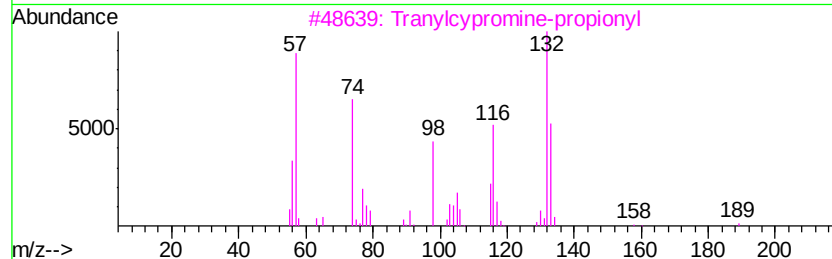
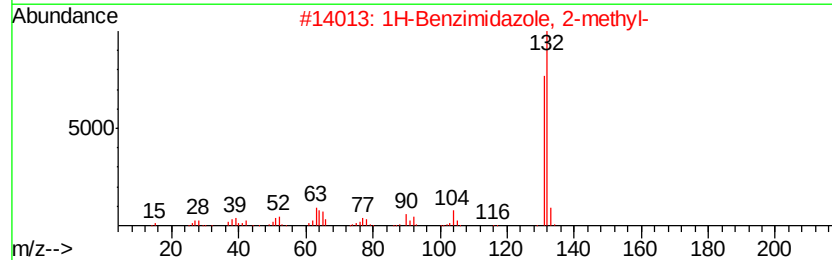
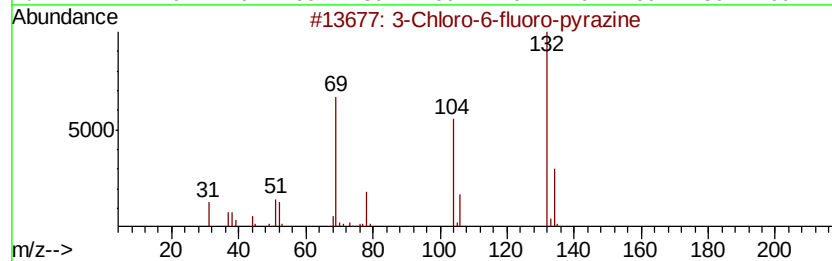
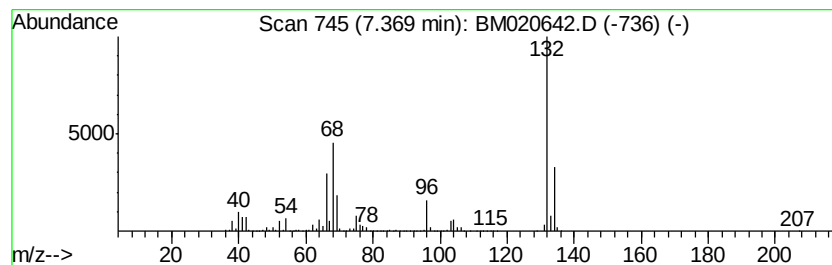
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TIC Library : C:\DATABASE\NIST02.L
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 Peak Number 2 unknown7.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	110.51 ng	11073700	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	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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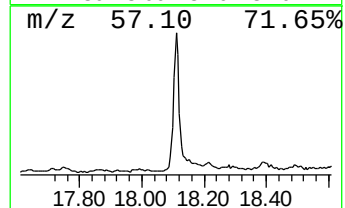
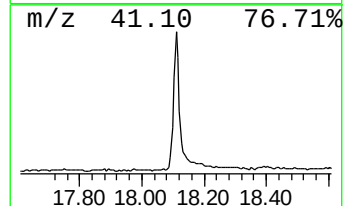
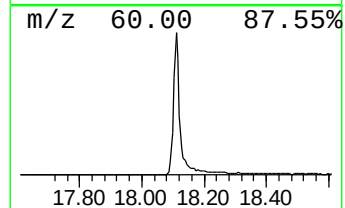
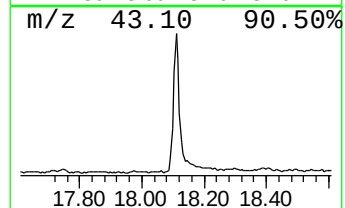
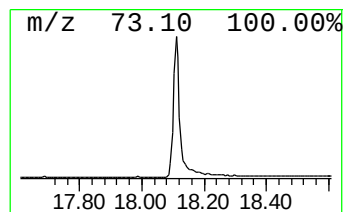
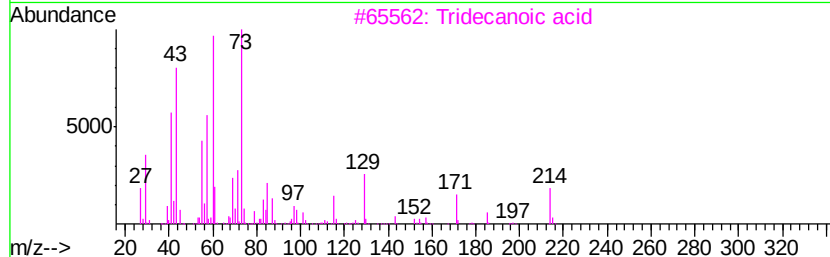
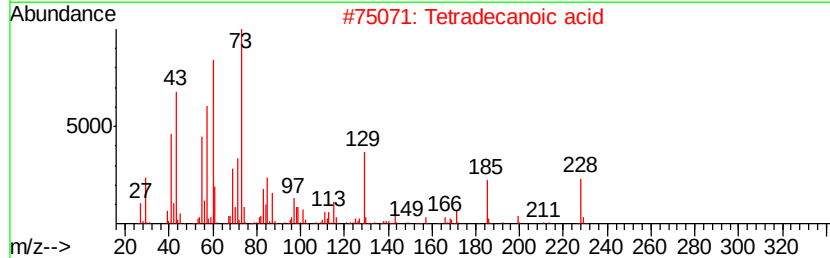
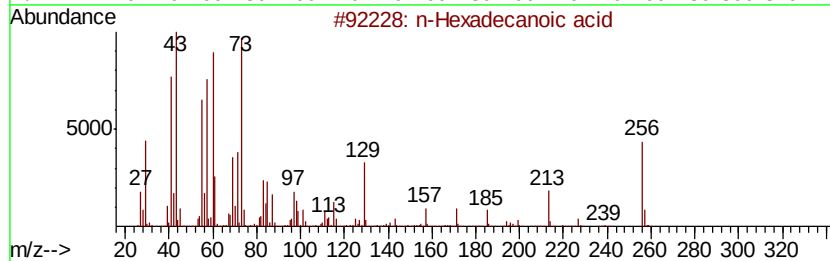
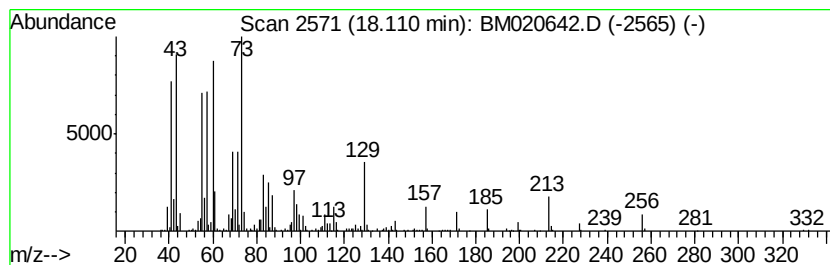
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	7.49 ng	1130330	Phenanthrene-d10	17.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	87
4	Estra-1,3,5(10)-trien-17.β.-ol	256	C18H24O	002529-64-8	30
5	Hexadecane	226	C16H34	000544-76-3	11



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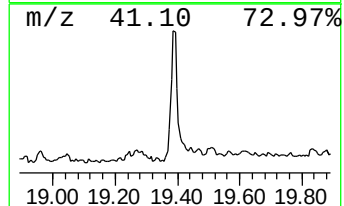
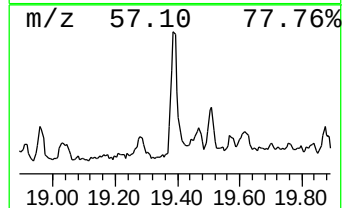
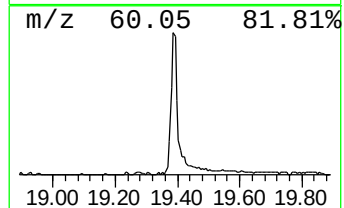
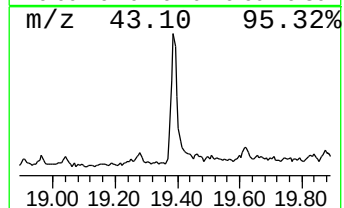
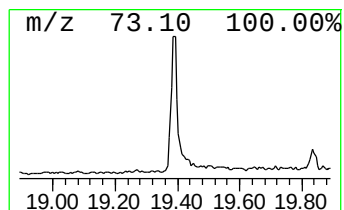
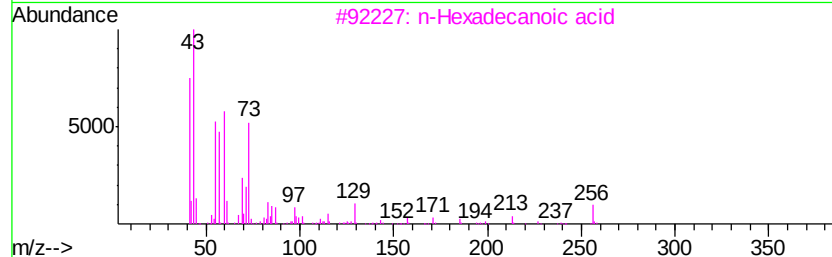
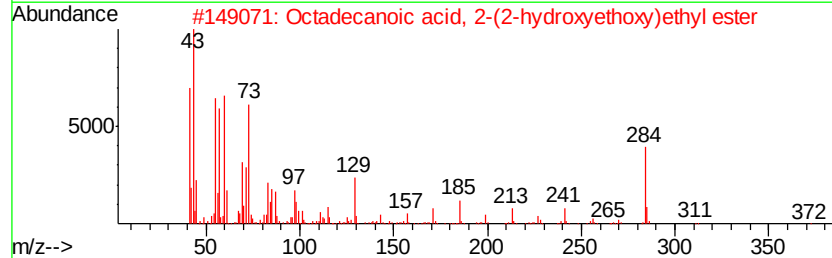
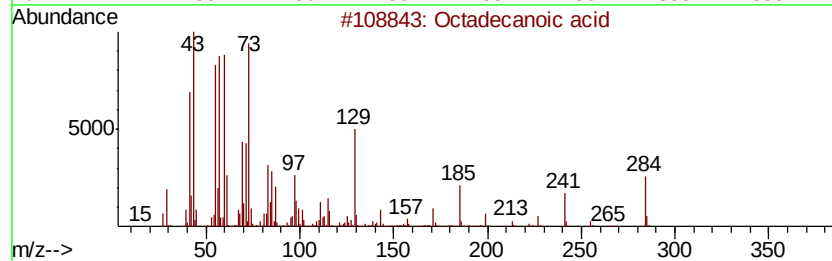
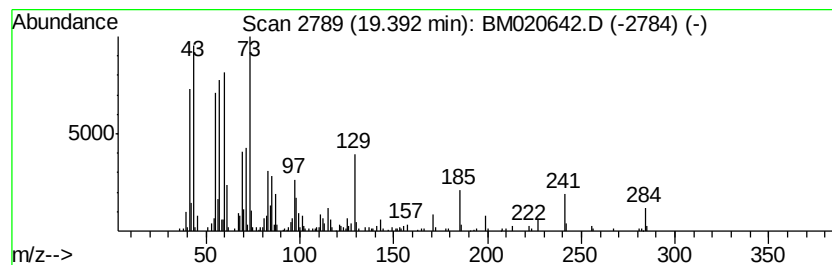
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	3.63 ng	544938	Chrysene-d12	21.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	91
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



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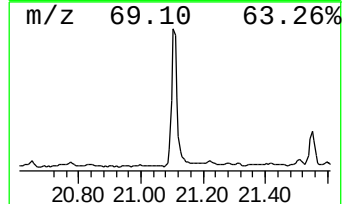
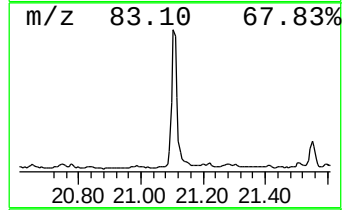
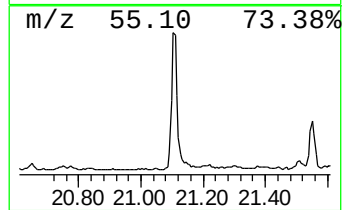
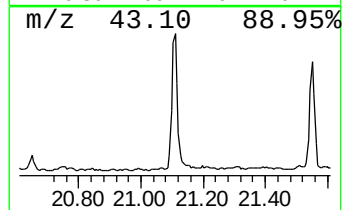
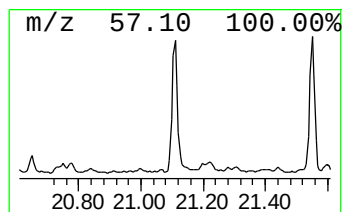
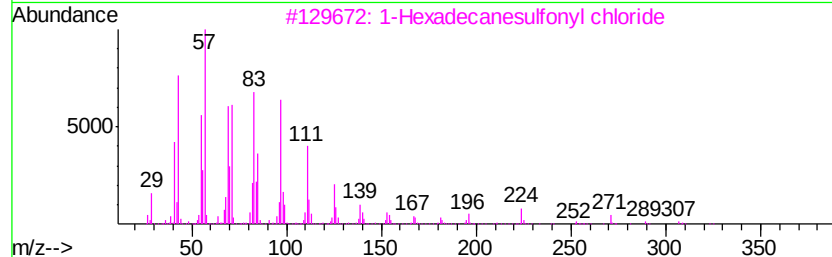
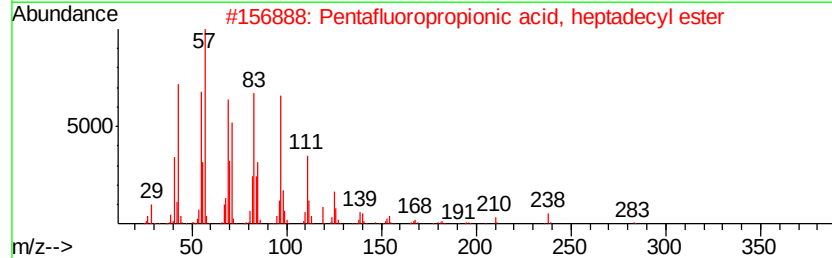
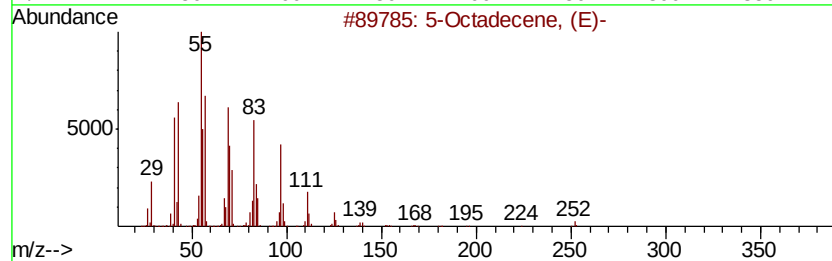
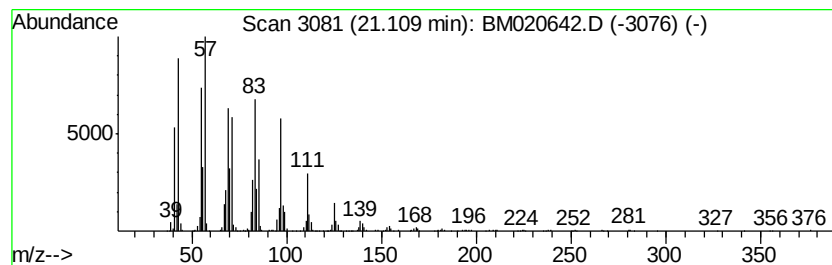
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Peak Number 6 5-Octadecene, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.11	10.02 ng	1503050	Chrysene-d12	21.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	94
3		1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	91
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
5		17-Pentatriacontene	491	C35H70	006971-40-0	90



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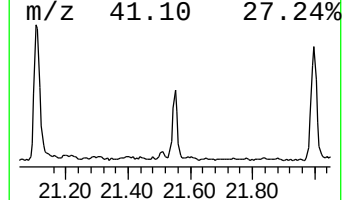
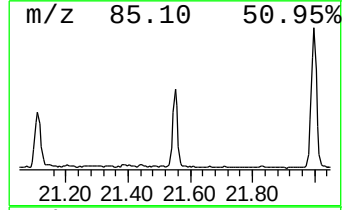
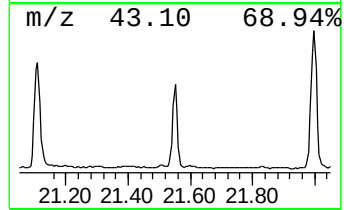
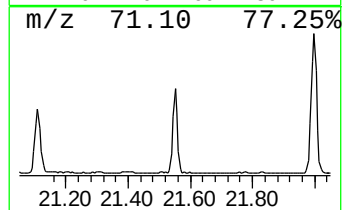
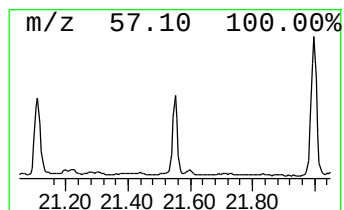
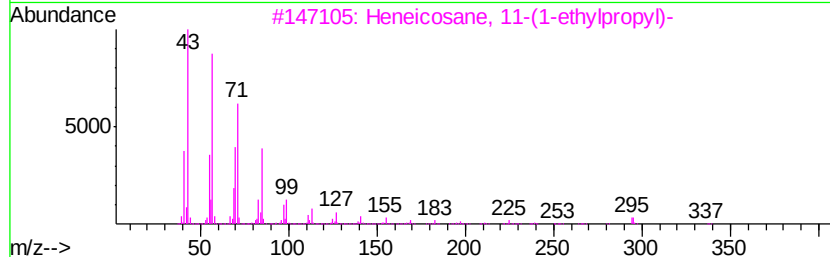
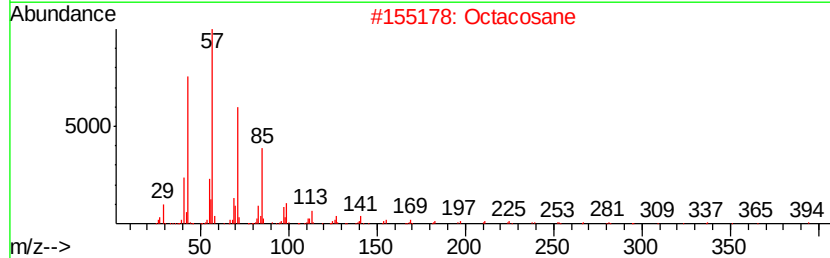
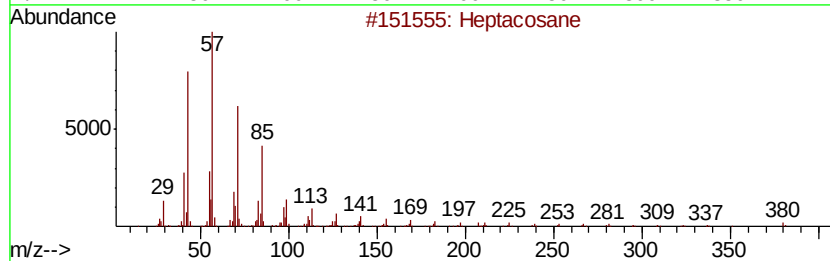
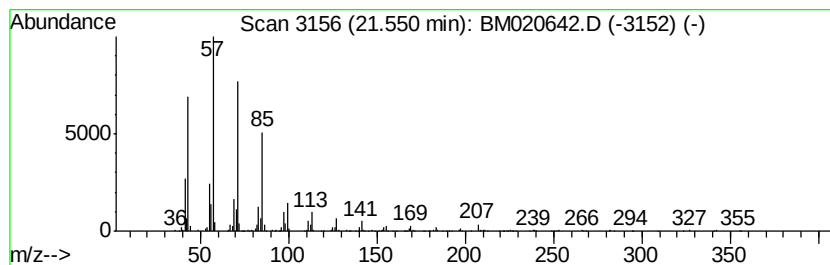
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM053119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.55	4.53 ng	679642	Chrysene-d12		21.39
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	91
2	Octacosane	394	C28H58	000630-02-4	91
3	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
Operator : JU/SJ
Sample : K3072-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
TP-1

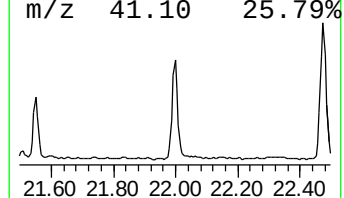
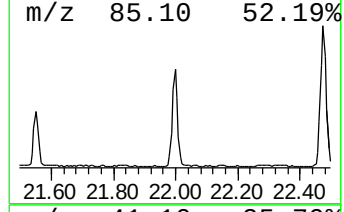
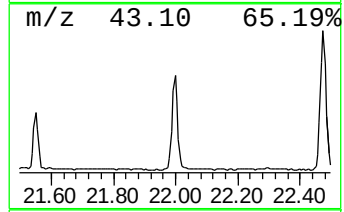
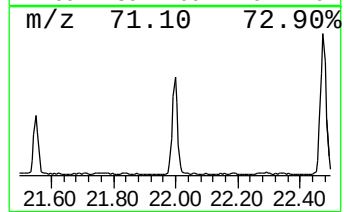
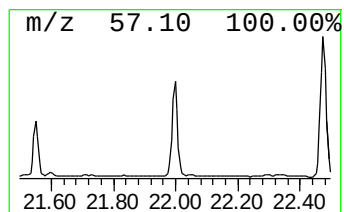
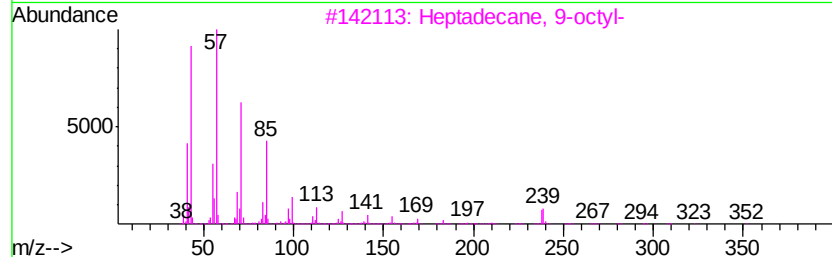
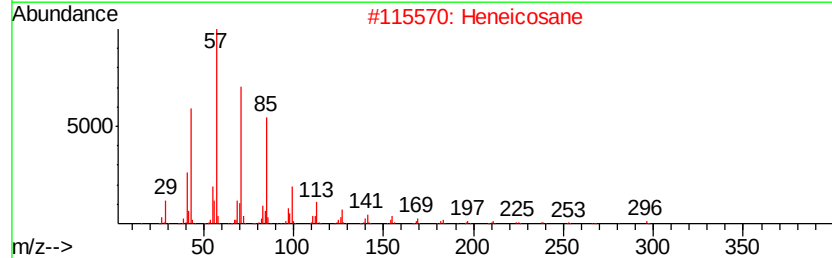
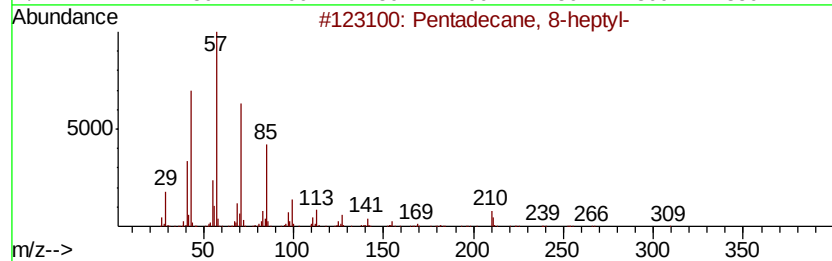
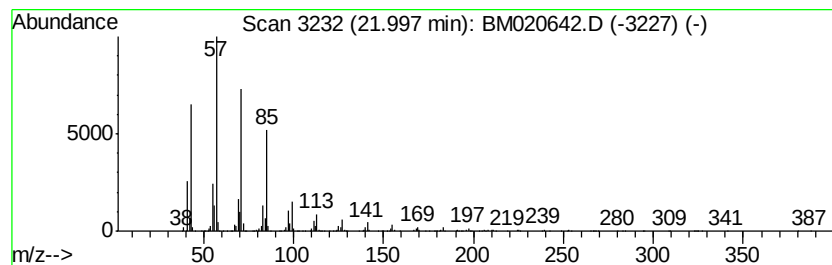
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.00	9.18 ng	1376620	Chrysene-d12	21.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Hexatriacontane	507	C36H74	000630-06-8	91
5		triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
Operator : JU/SJ
Sample : K3072-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

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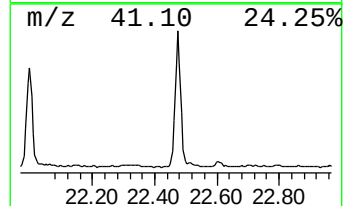
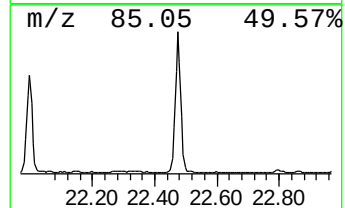
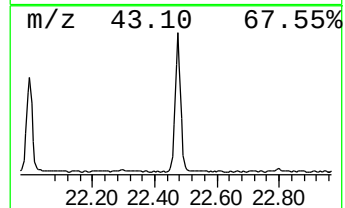
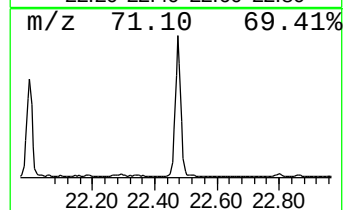
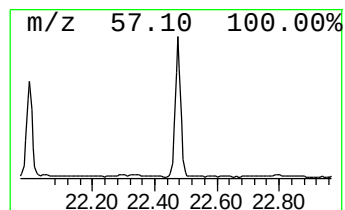
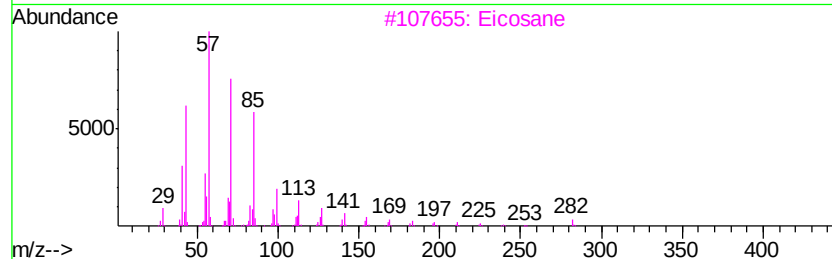
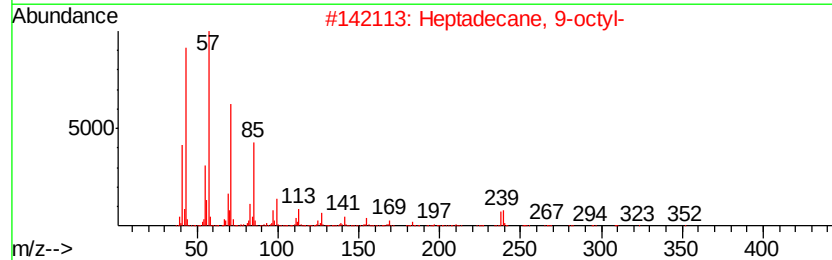
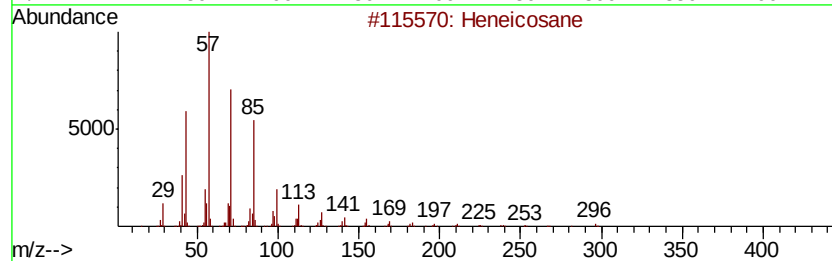
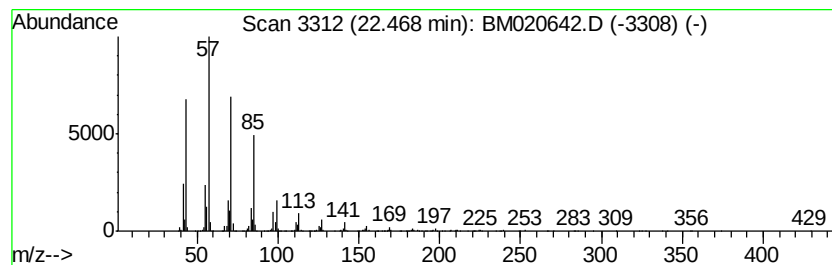
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.47	13.16 ng	1974630	Chrysene-d12	21.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
3	Eicosane		282	C20H42	000112-95-8	91
4	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91
5	Octacosane		394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
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Sample : K3072-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

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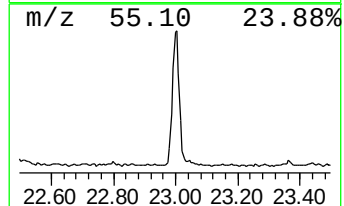
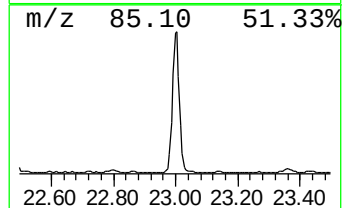
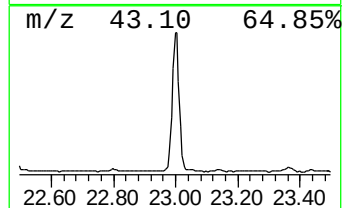
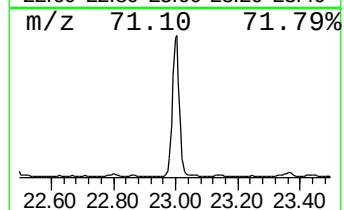
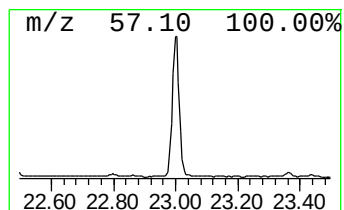
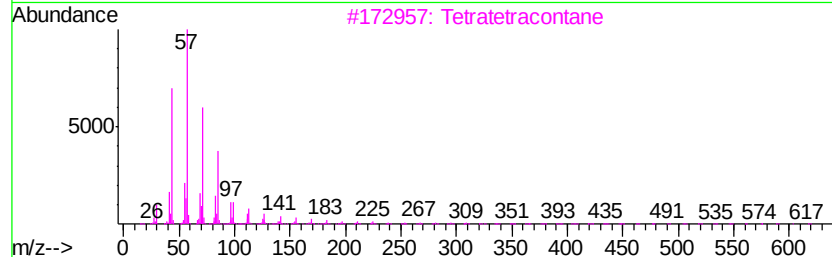
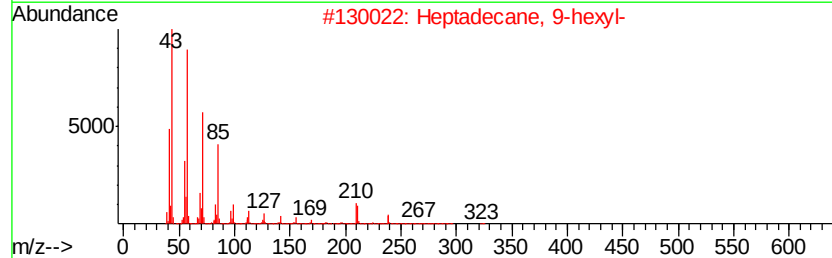
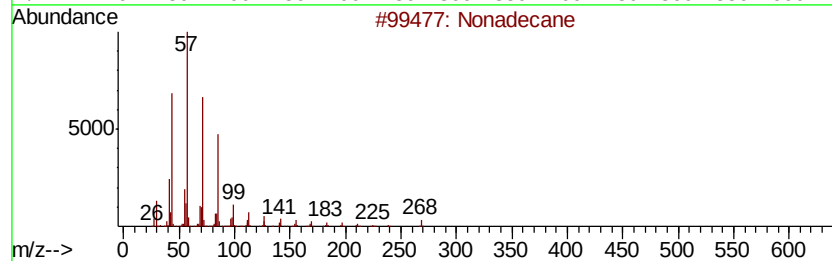
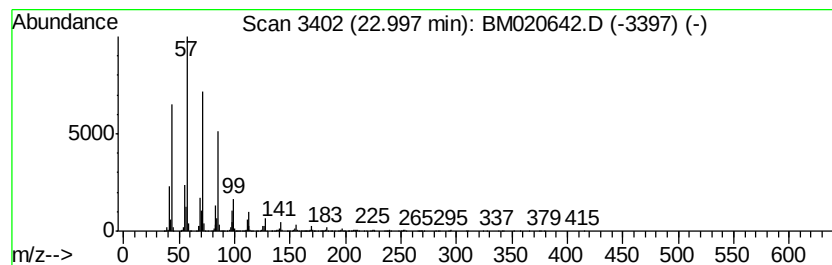
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Nonadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.00	14.90 ng	2347250	Perylene-d12	23.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
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Sample : K3072-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

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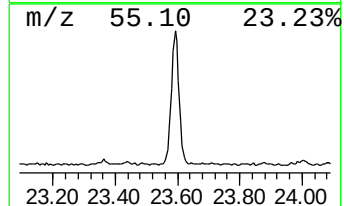
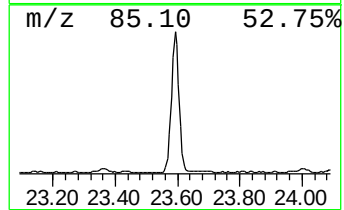
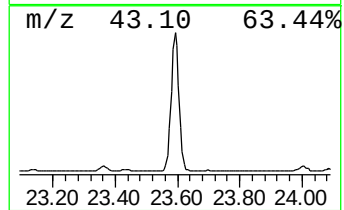
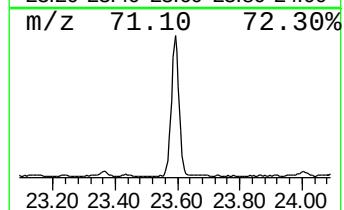
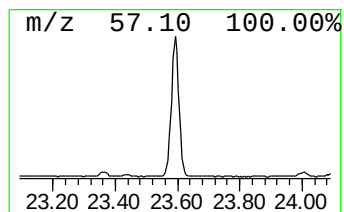
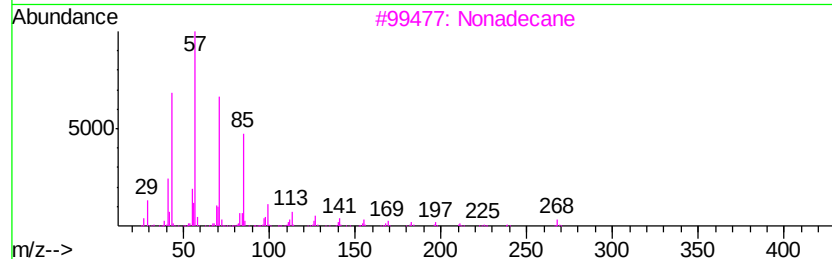
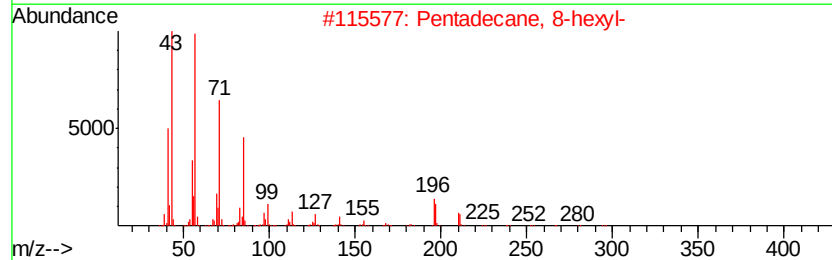
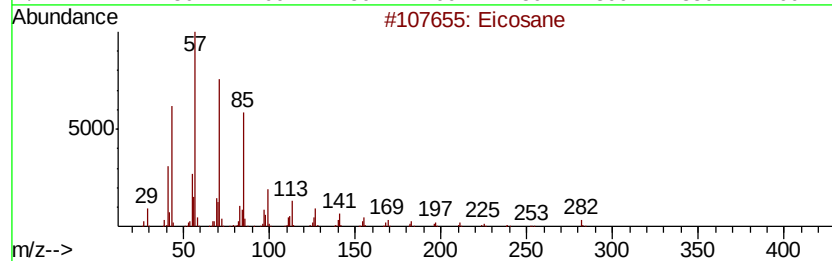
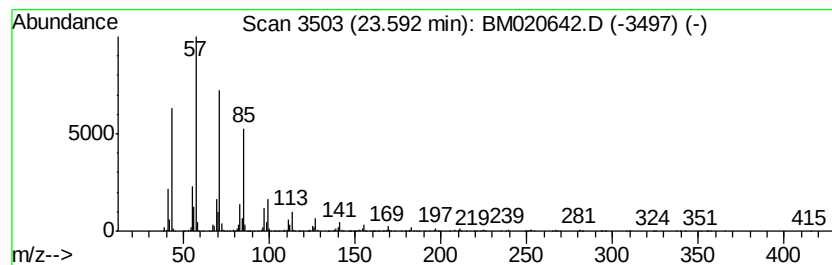
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.59	14.20 ng	2236500	Perylene-d12	23.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Nonadecane	268	C19H40	000629-92-5	93
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
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Sample : K3072-01
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Instrument :
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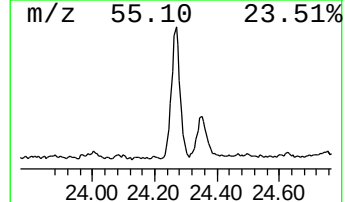
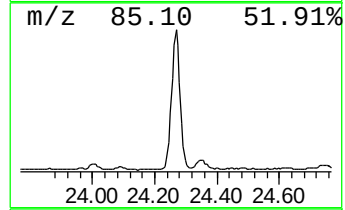
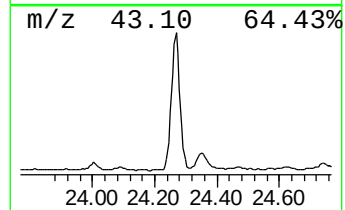
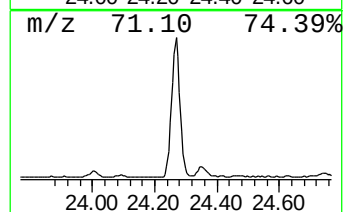
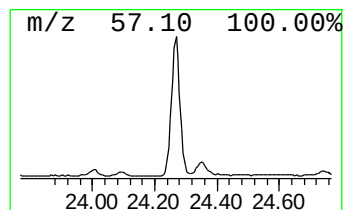
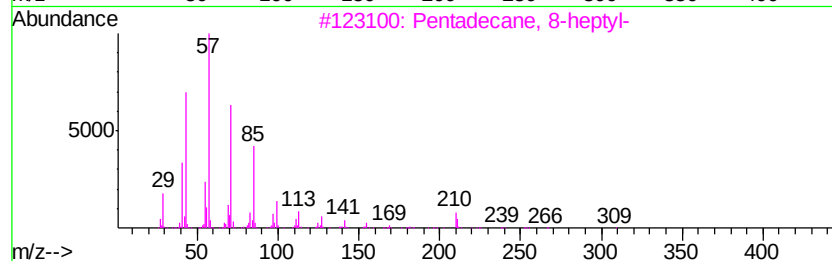
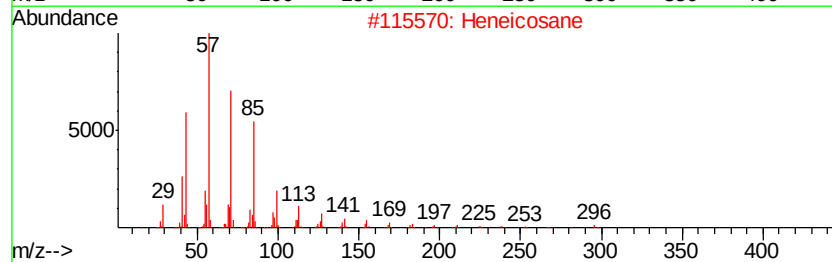
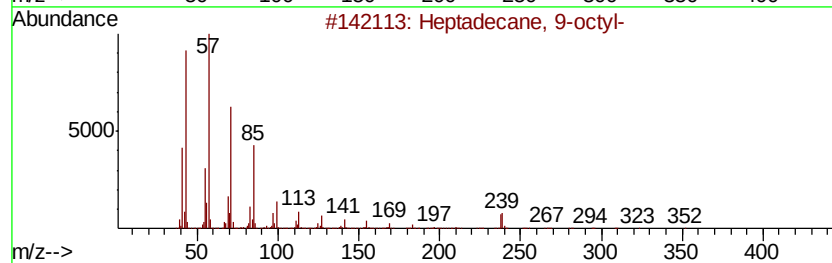
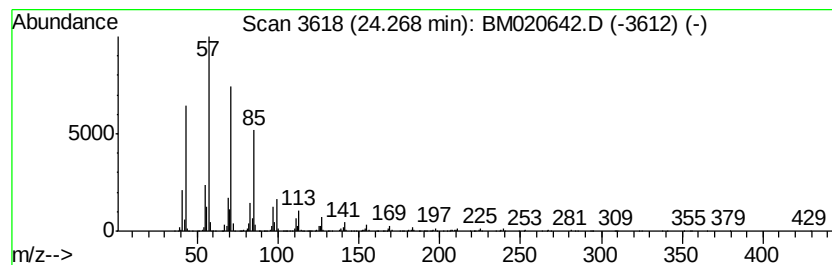
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptadecane, 9-octyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.27	10.99 ng	1730480	Perylene-d12	23.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Tetratriacontane	479	C34H70	014167-59-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
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Sample : K3072-01
Misc :
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ClientSampleId :
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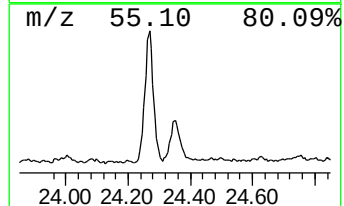
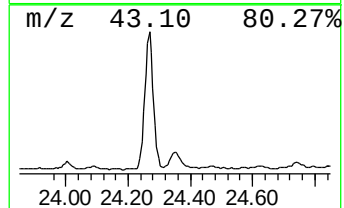
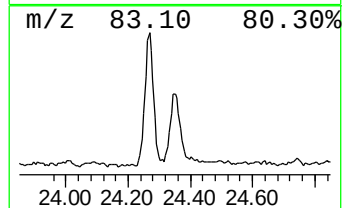
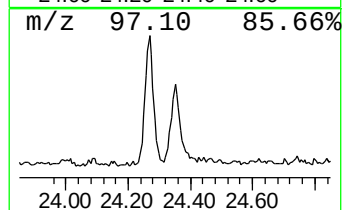
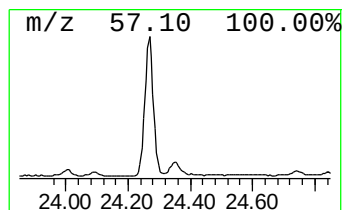
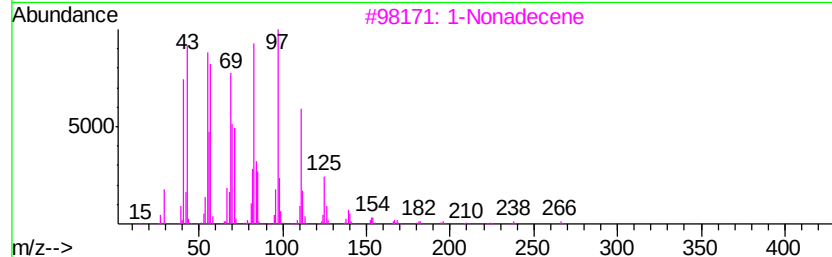
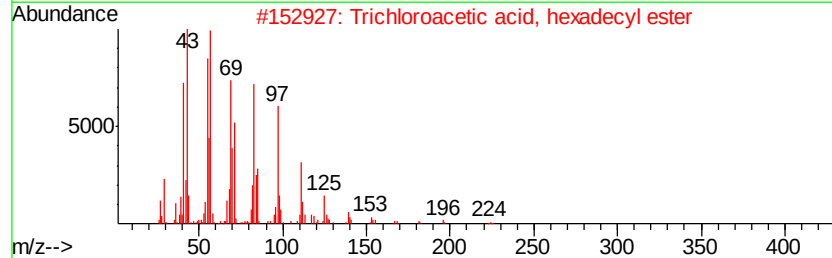
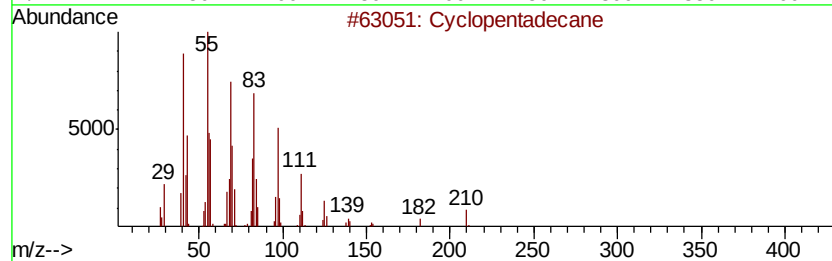
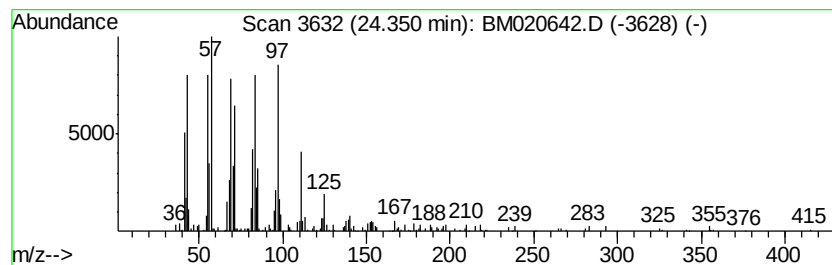
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopentadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.35	2.32 ng	365044	Perylene-d12	23.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	95
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
3		1-Nonadecene	266	C19H38	018435-45-5	83
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	81
5		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM053119\
Data File : BM020642.D
Acq On : 31 May 2019 23:12
Operator : JU/SJ
Sample : K3072-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown7.37	7.37	110.5	ng	11073700	1	7.83	2004070	20.0
n-Hexadecanoic acid	18.11	7.5	ng	1130330	4	17.22	3016540	20.0
Octadecanoic acid	19.39	3.6	ng	544938	5	21.39	3000460	20.0
5-Octadecene, (E)-	21.11	10.0	ng	1503050	5	21.39	3000460	20.0
Heptacosane	21.55	4.5	ng	679642	5	21.39	3000460	20.0
Pentadecane, 8-he...	22.00	9.2	ng	1376620	5	21.39	3000460	20.0
Heneicosane	22.47	13.2	ng	1974630	5	21.39	3000460	20.0
Nonadecane	23.00	14.9	ng	2347250	6	23.69	3149670	20.0
Eicosane	23.59	14.2	ng	2236500	6	23.69	3149670	20.0
Heptadecane, 9-oc...	24.27	11.0	ng	1730480	6	23.69	3149670	20.0
Cyclopentadecane	24.35	2.3	ng	365044	6	23.69	3149670	20.0