

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114129.D
 Acq On : 10 May 2019 19:15
 Operator : JU/SJ
 Sample : K2762-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P021-SS001-1824-01

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_F\METHODS\8270-BF051019.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	2.134	3	8	28	rVB	1354821	1891437	24.98%	2.638%
2	5.481	572	577	580	rBV	7285080	6997554	92.40%	9.759%
3	6.140	686	689	692	rBV	121424	104009	1.37%	0.145%
4	6.487	743	748	751	rBV	7334605	7572899	100.00%	10.561%
5	6.557	758	760	767	rVB	84102	77362	1.02%	0.108%
6	6.622	767	771	775	rBV	7361884	7209303	95.20%	10.054%
7	6.851	806	810	814	rBV	2609020	2094129	27.65%	2.920%
8	7.010	833	837	840	rBV	6398047	5429737	71.70%	7.572%
9	7.410	900	905	908	rBV	5117124	4657391	61.50%	6.495%
10	8.133	1024	1028	1031	rBV	2877903	2598985	34.32%	3.624%
11	9.204	1206	1210	1214	rBV	7553967	6891193	91.00%	9.610%
12	9.592	1273	1276	1280	rBV	1229822	1079037	14.25%	1.505%
13	9.886	1322	1326	1329	rVB	2914266	2585858	34.15%	3.606%
14	10.669	1455	1459	1463	rBV	4337707	4109811	54.27%	5.731%
15	11.369	1574	1578	1581	rBV	2779821	2396068	31.64%	3.341%
16	11.898	1664	1668	1680	rBV	844605	1004338	13.26%	1.401%
17	12.580	1781	1784	1787	rBV	149385	146403	1.93%	0.204%
18	12.674	1797	1800	1804	rBV	256219	285574	3.77%	0.398%
19	12.804	1819	1822	1825	rBV	208007	182907	2.42%	0.255%
20	12.951	1843	1847	1850	rBV	5565367	4904060	64.76%	6.839%
21	13.527	1943	1945	1951	rVB2	86029	94642	1.25%	0.132%
22	13.798	1988	1991	1995	rBV5	151661	190057	2.51%	0.265%
23	13.851	1997	2000	2012	rVB	901837	1217086	16.07%	1.697%
24	14.015	2023	2028	2031	rBV	2190855	2155457	28.46%	3.006%
25	14.039	2031	2032	2040	rVB	105122	84893	1.12%	0.118%
26	14.186	2054	2057	2062	rVB	221326	207288	2.74%	0.289%
27	14.504	2108	2111	2116	rBV	330421	394743	5.21%	0.550%
28	14.821	2162	2165	2168	rVB	381766	341746	4.51%	0.477%
29	14.898	2175	2178	2181	rVB	155379	167719	2.21%	0.234%
30	15.157	2218	2222	2226	rBV	709776	864777	11.42%	1.206%
31	15.504	2277	2281	2285	rBV	1593202	2052833	27.11%	2.863%
32	15.539	2285	2287	2294	rVB	516404	629019	8.31%	0.877%
33	15.974	2357	2361	2367	rBV	504230	771337	10.19%	1.076%
34	16.480	2443	2447	2452	rVB	195922	317432	4.19%	0.443%

LSC Area Percent Report

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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_F\METHODS\8270-BF051019.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

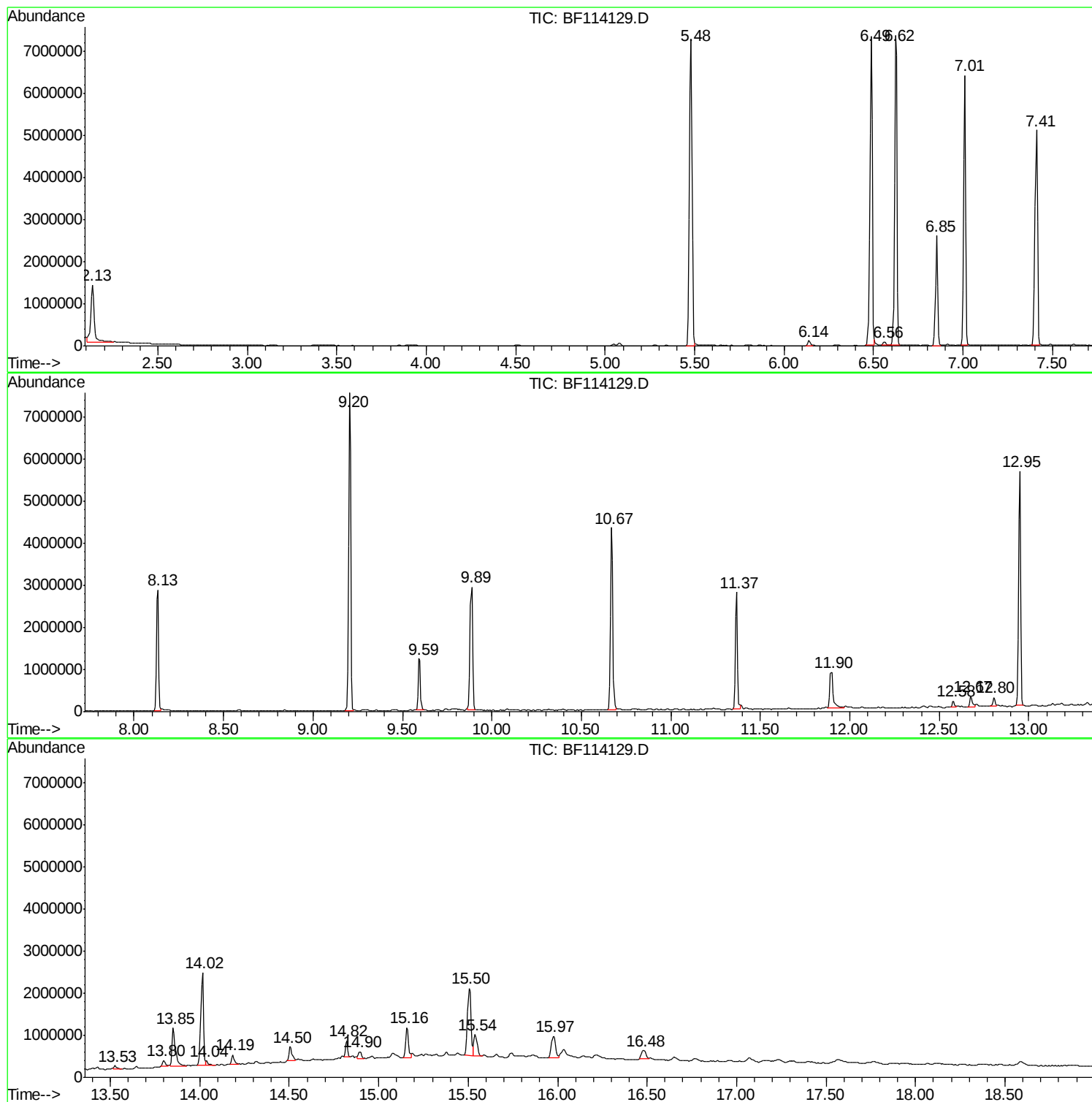
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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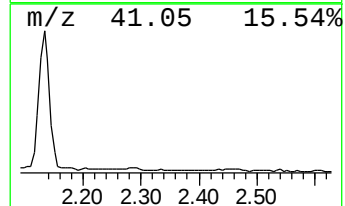
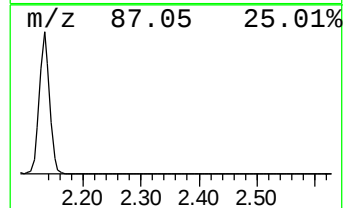
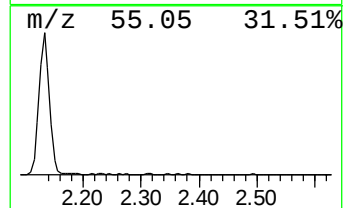
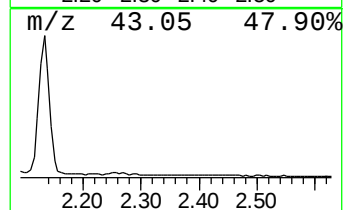
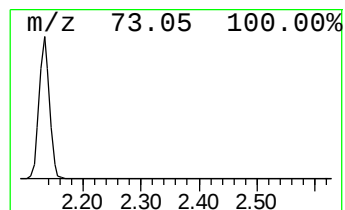
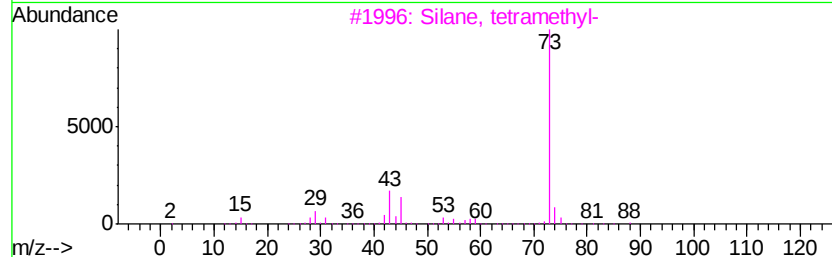
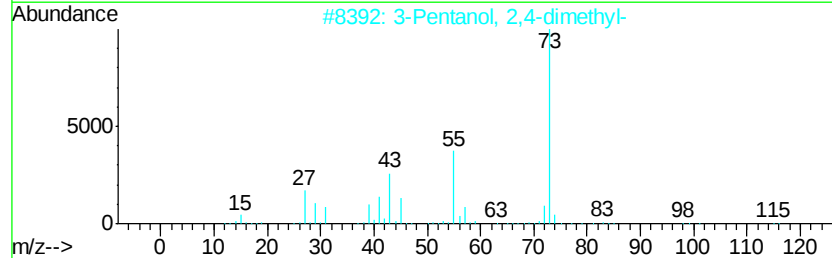
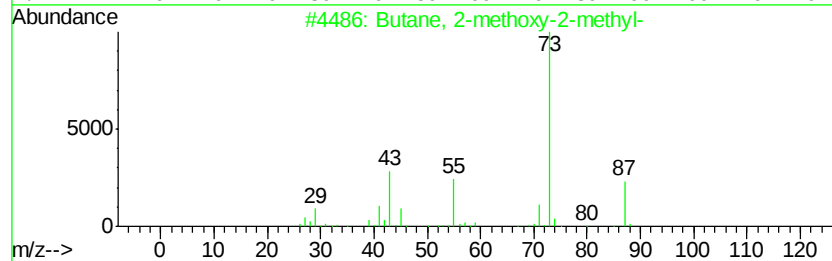
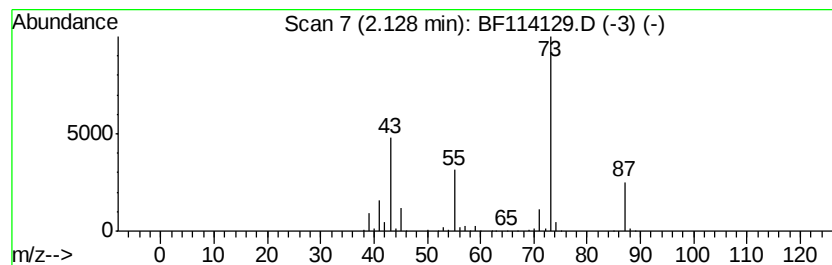
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	18.06 ng	1891440	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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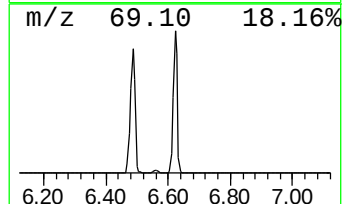
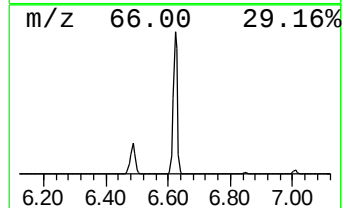
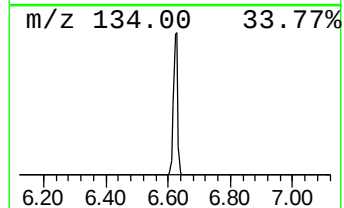
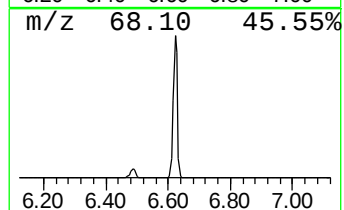
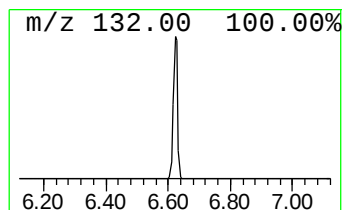
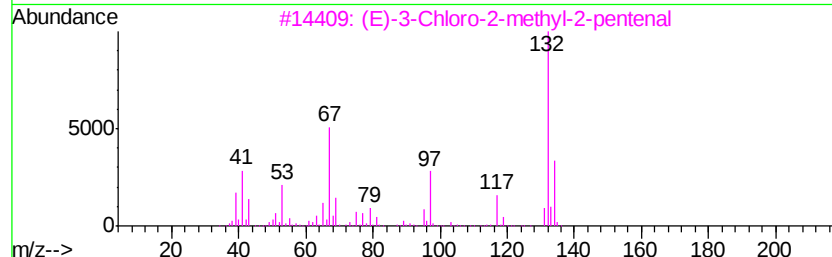
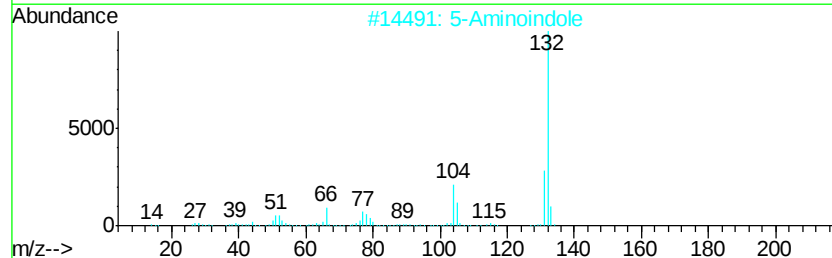
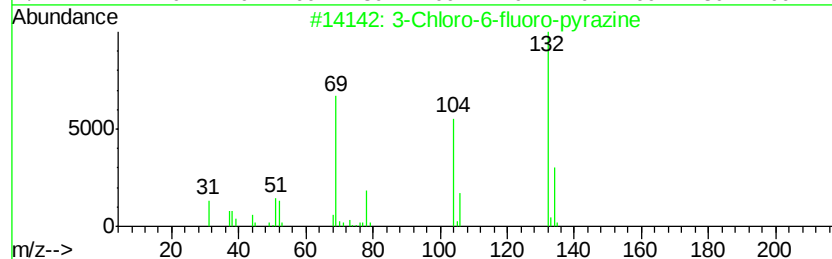
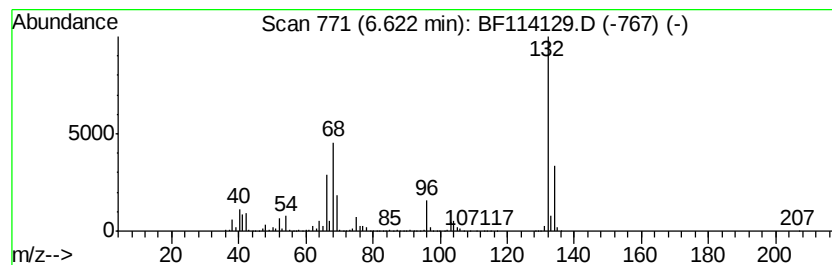
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	68.85 ng	7209300	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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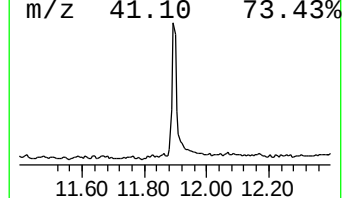
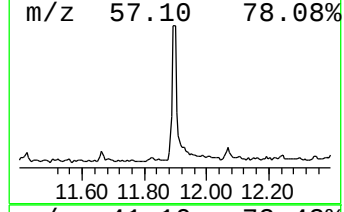
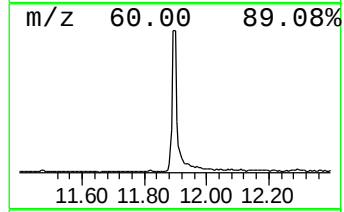
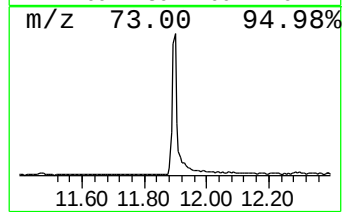
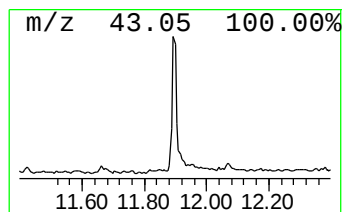
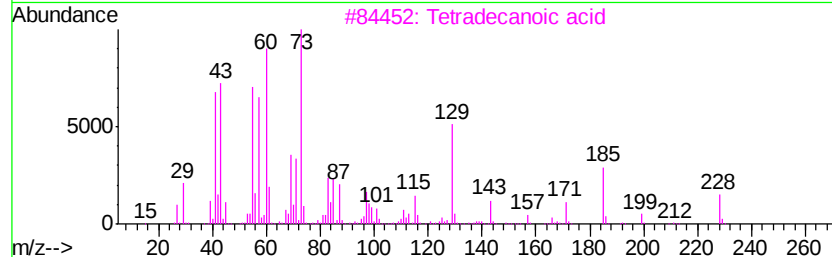
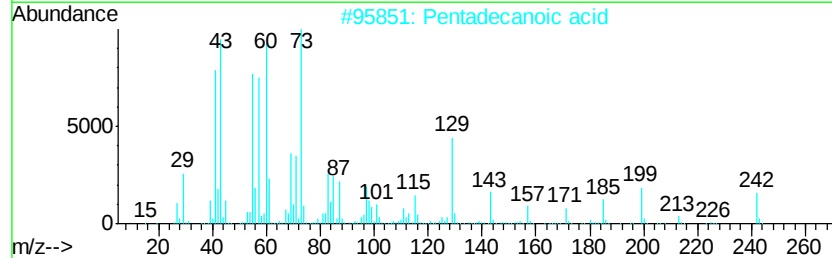
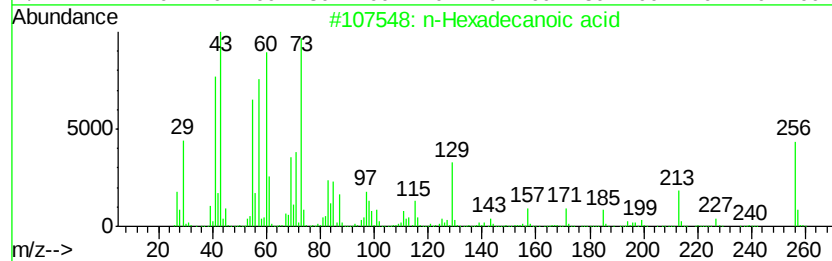
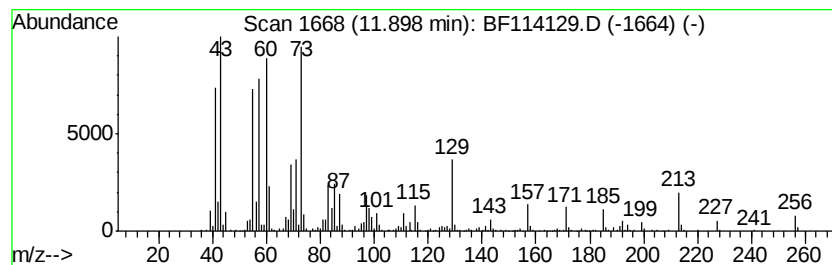
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.90	8.38 ng	1004340	Phenanthrene-d10	11.37	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	76
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	Undecanoic acid	186	C11H22O2	000112-37-8	68



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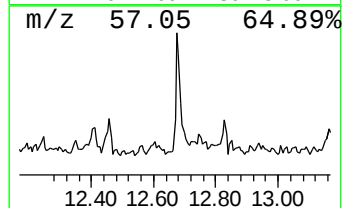
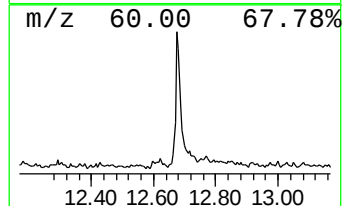
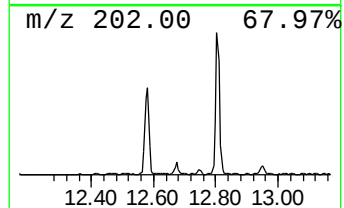
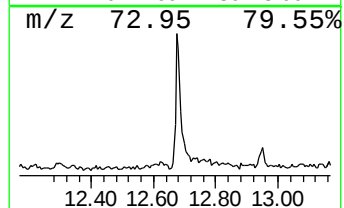
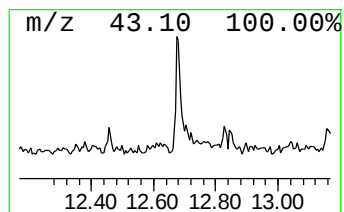
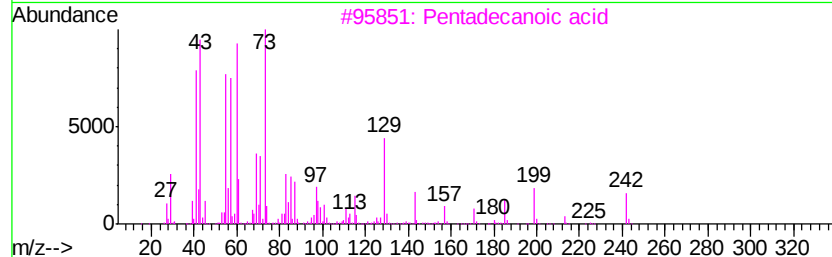
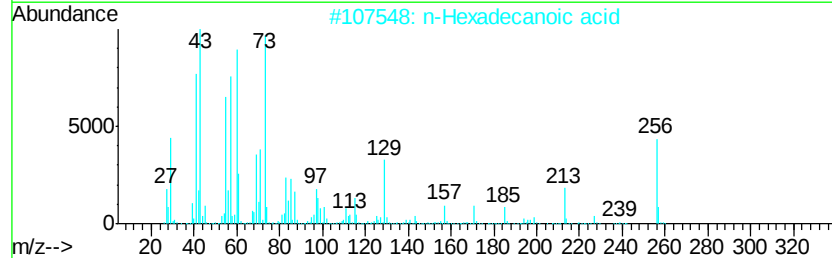
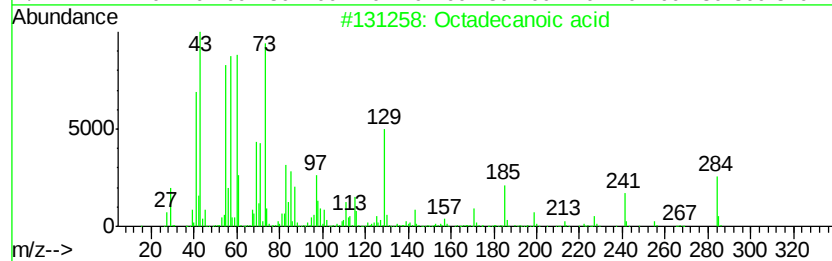
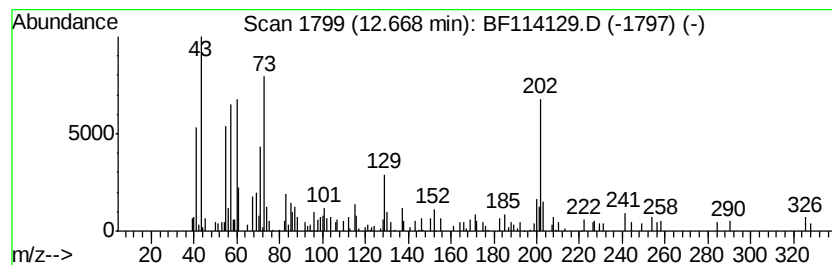
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	2.38 ng	285574	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4		Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	64
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



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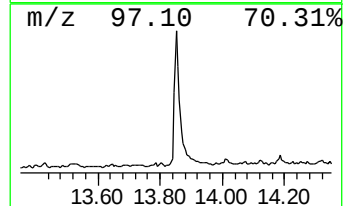
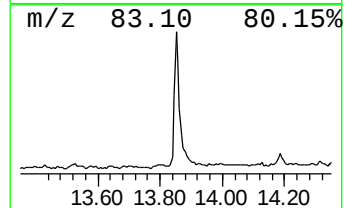
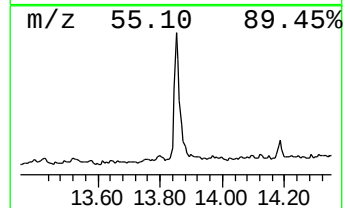
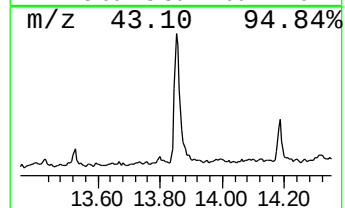
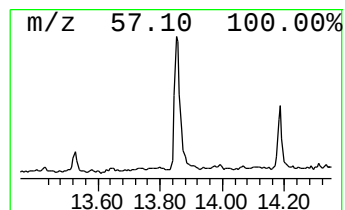
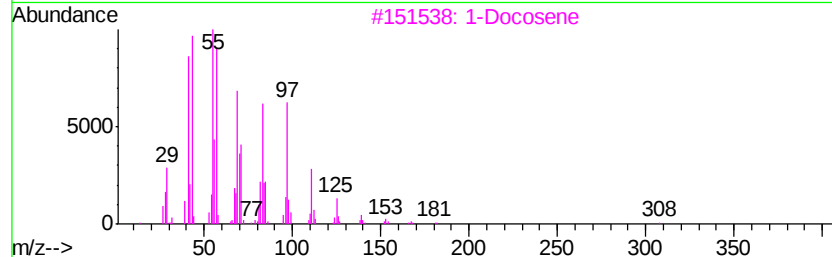
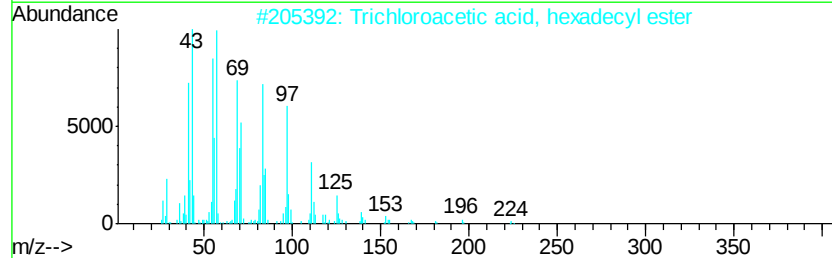
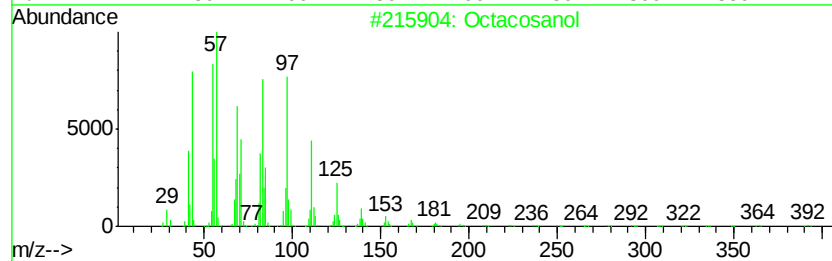
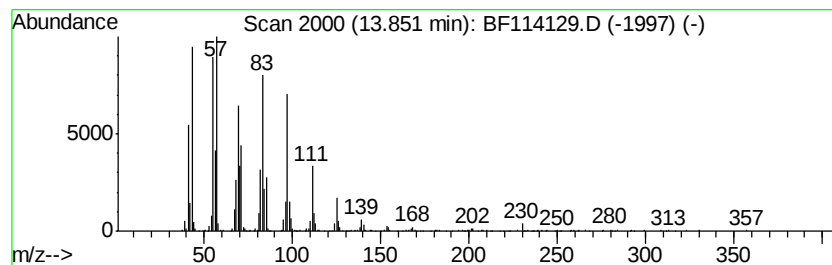
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Peak Number 6 Octacosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.85	11.29 ng	1217090	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosanol	410	C28H58O	000557-61-9	94
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3	1-Docosene	308	C22H44	001599-67-3	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114129.D
Acq On : 10 May 2019 19:15
Operator : JU/SJ
Sample : K2762-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

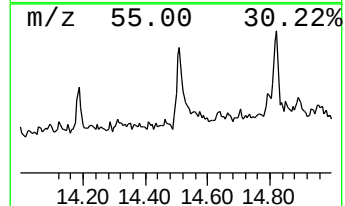
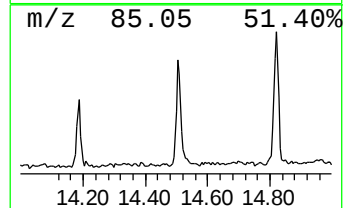
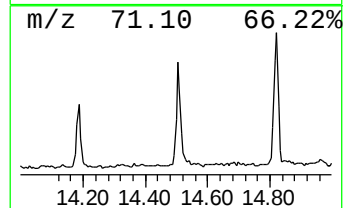
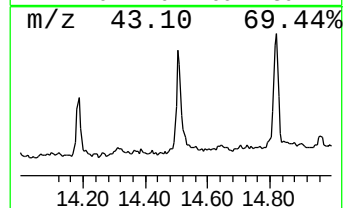
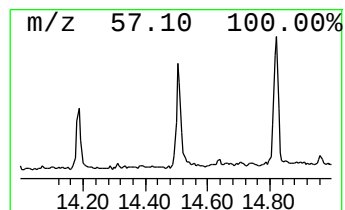
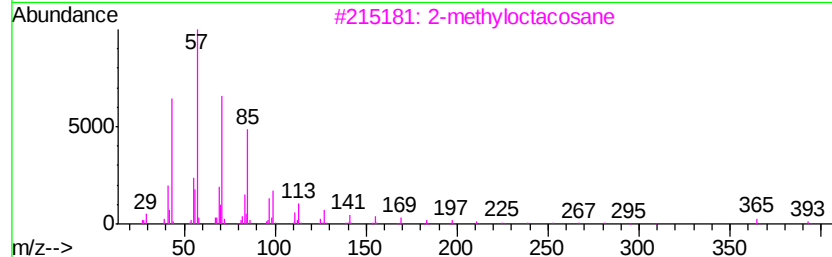
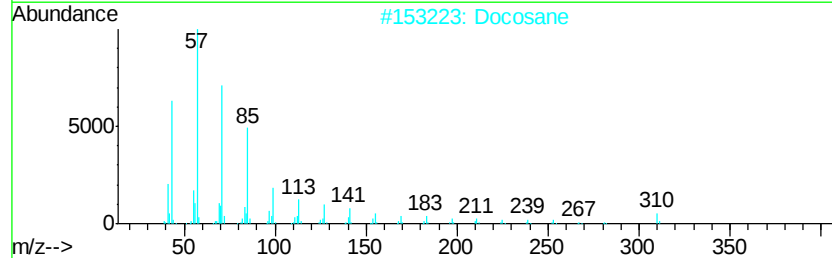
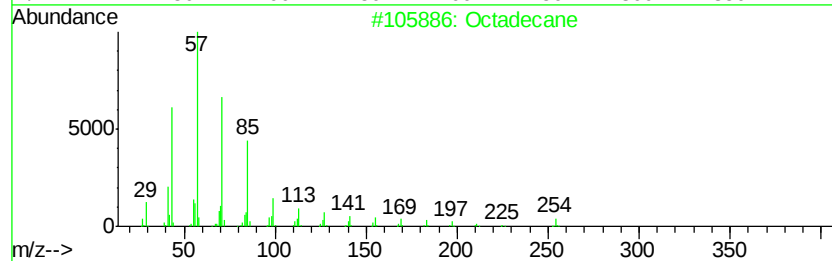
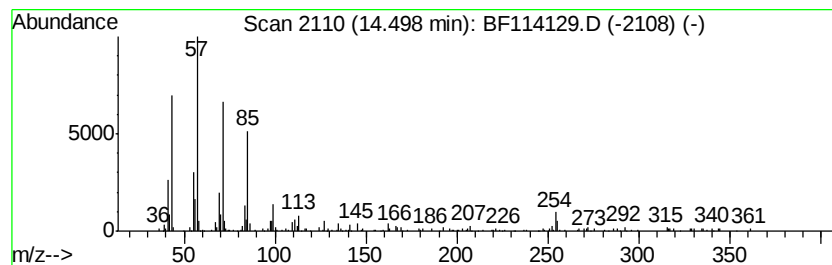
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BNA_F
ClientSampleId :
P021-SS001-1824-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	3.66 ng	394743	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octadecane	254 C18H38	000593-45-3	93
2	Docosane	310 C22H46	000629-97-0	91
3	2-methyloctacosane	408 C29H60	1000376-72-8	91
4	Heptacosane	380 C27H56	000593-49-7	91
5	Hexatriacontane	507 C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114129.D
Acq On : 10 May 2019 19:15
Operator : JU/SJ
Sample : K2762-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P021-SS001-1824-01

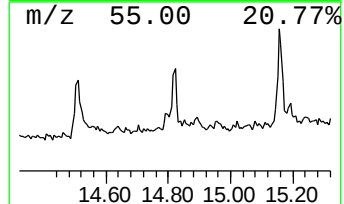
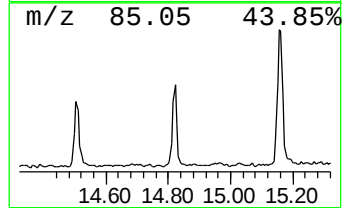
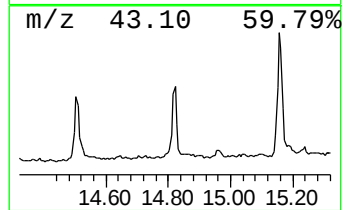
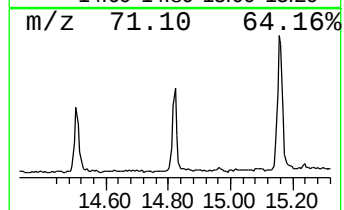
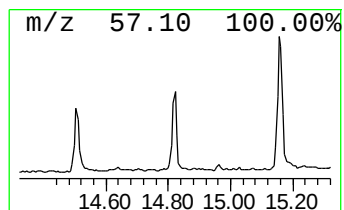
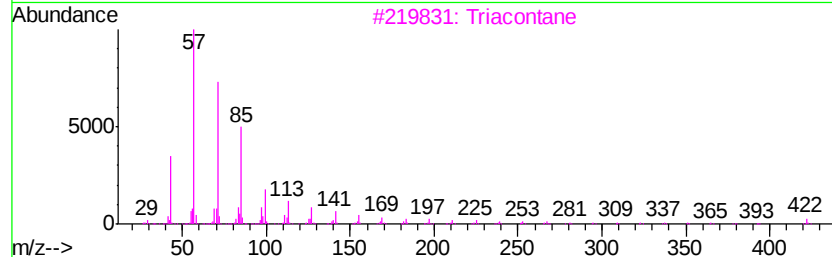
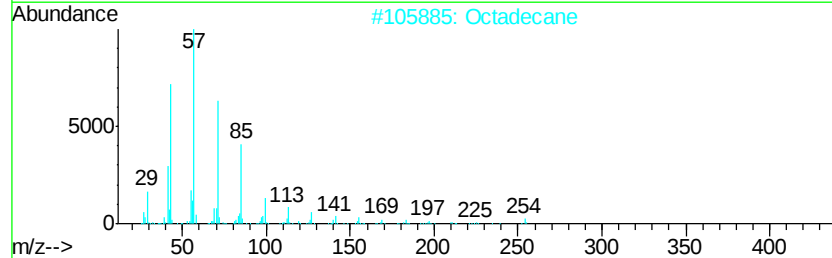
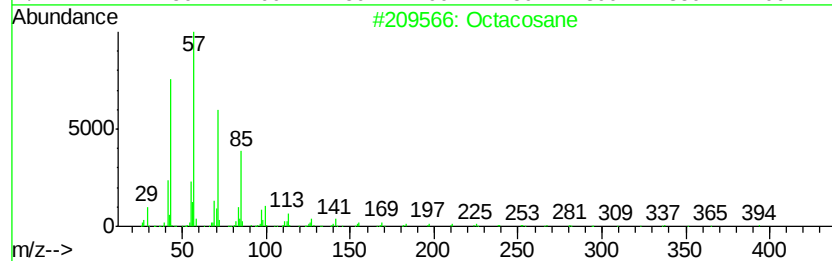
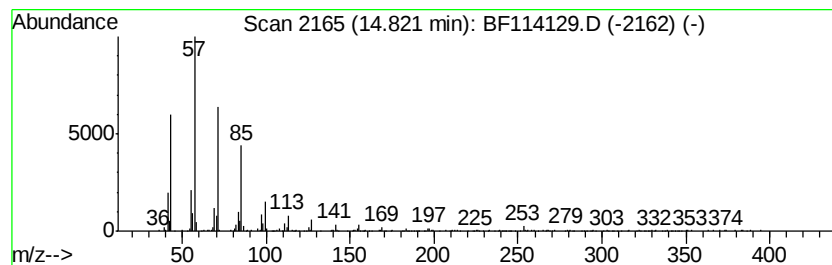
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	3.33 ng	341746	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	95
2		Octadecane	254	C18H38	000593-45-3	93
3		Triacontane	422	C30H62	000638-68-6	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114129.D
Acq On : 10 May 2019 19:15
Operator : JU/SJ
Sample : K2762-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P021-SS001-1824-01

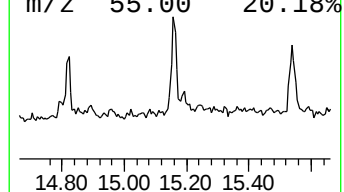
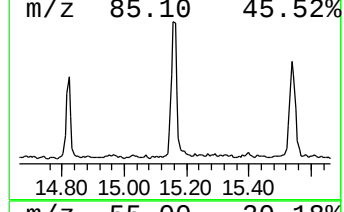
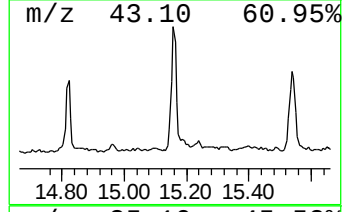
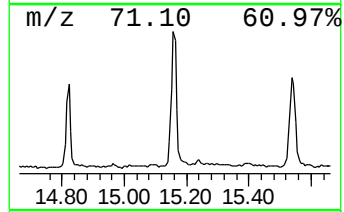
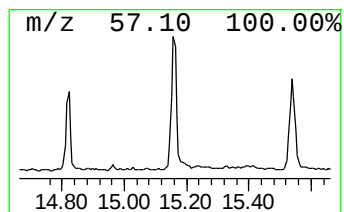
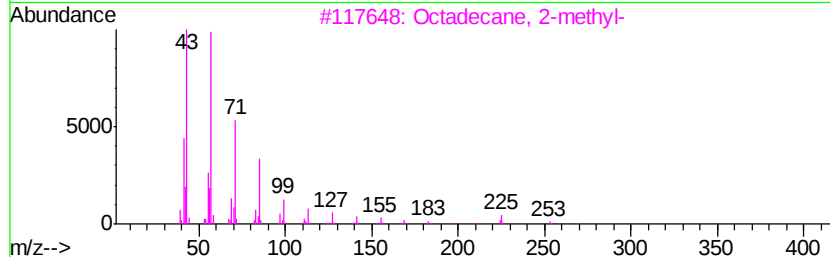
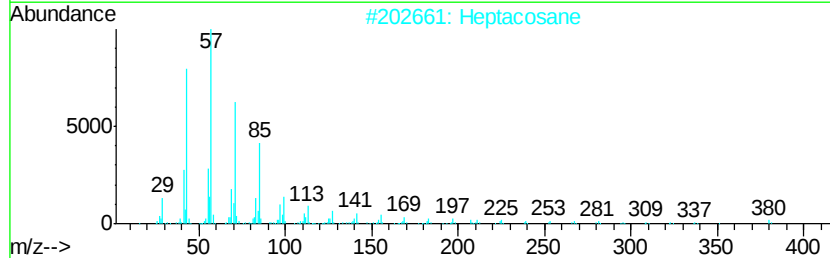
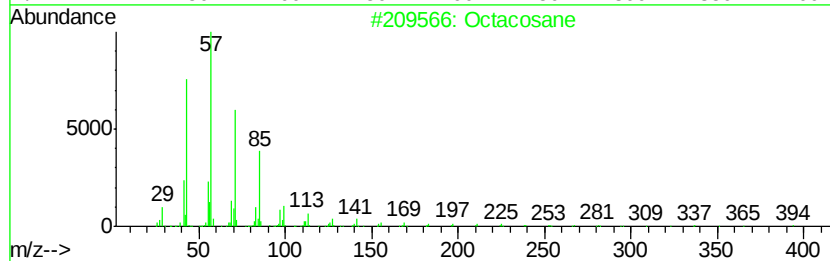
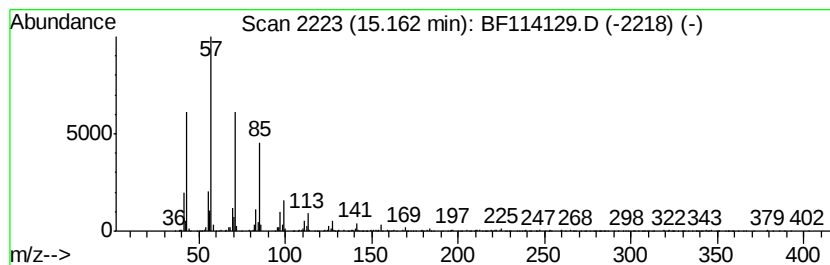
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	8.43 ng	864777	Perylene-d12	15.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
4		Nonadecane	268	C19H40	000629-92-5	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114129.D
Acq On : 10 May 2019 19:15
Operator : JU/SJ
Sample : K2762-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P021-SS001-1824-01

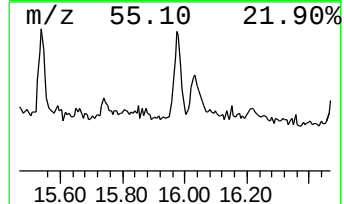
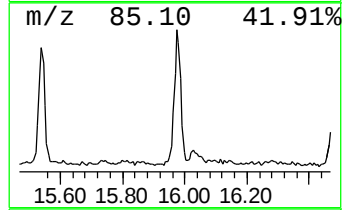
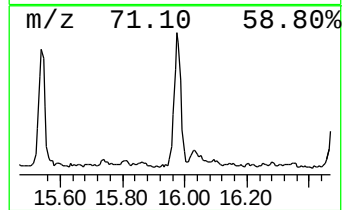
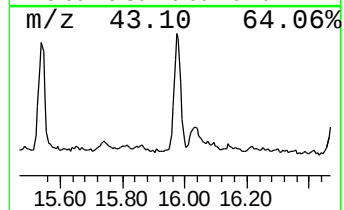
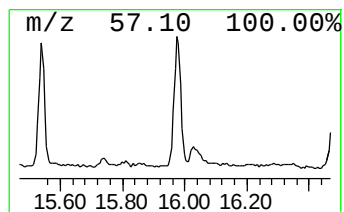
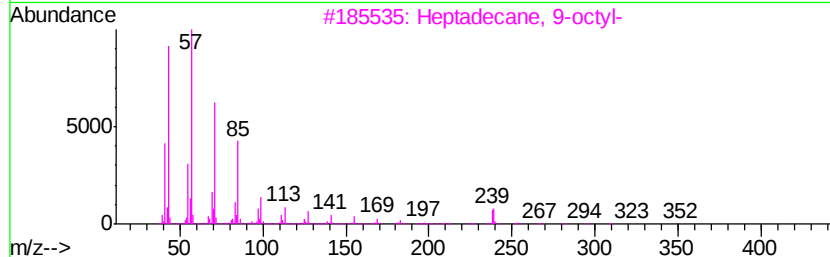
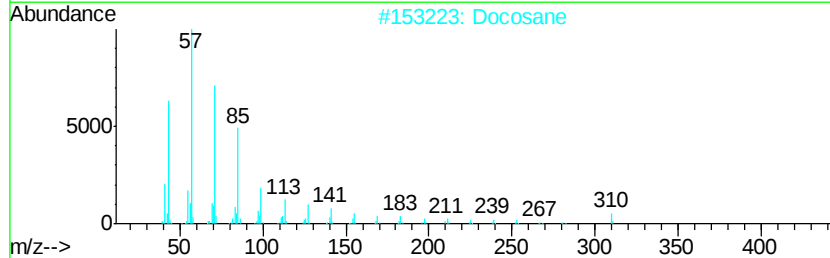
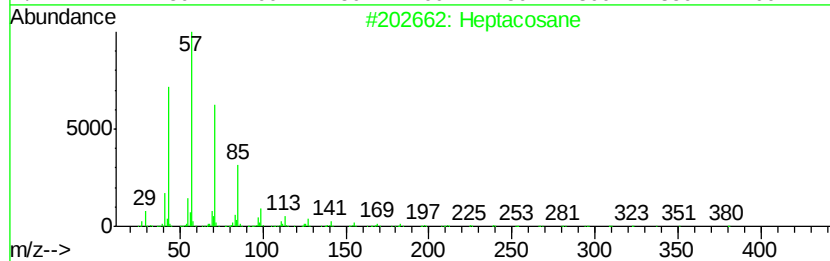
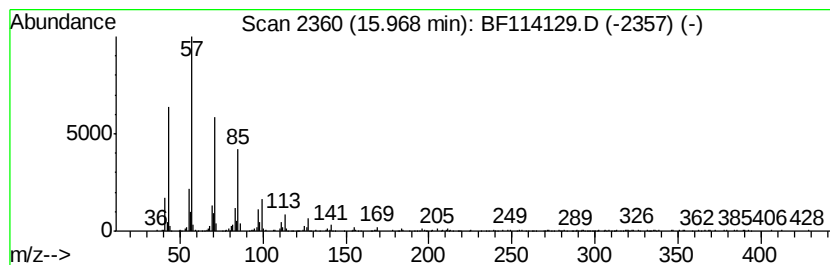
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	7.51 ng	771337	Perylene-d12	15.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	94
2			Docosane	310	C22H46	000629-97-0	93
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4			Eicosane	282	C20H42	000112-95-8	87
5			Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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Acq On : 10 May 2019 19:15
Operator : JU/SJ
Sample : K2762-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

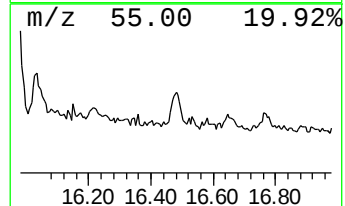
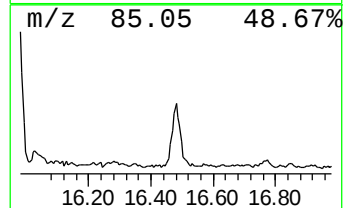
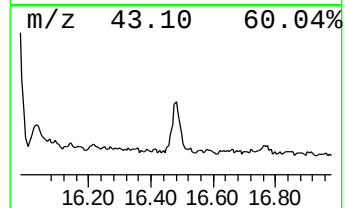
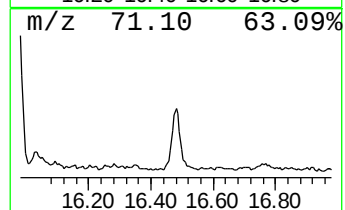
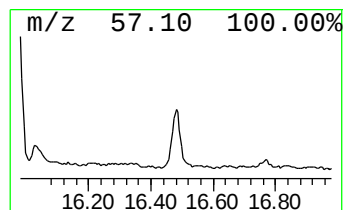
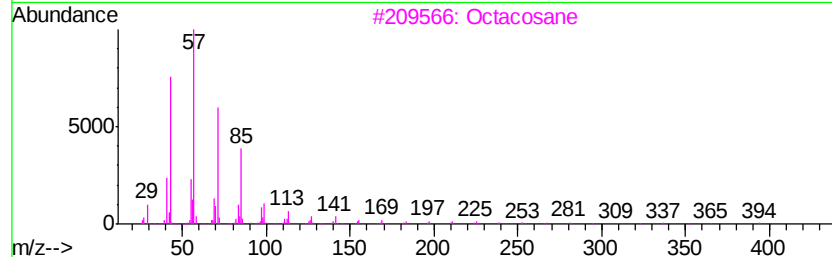
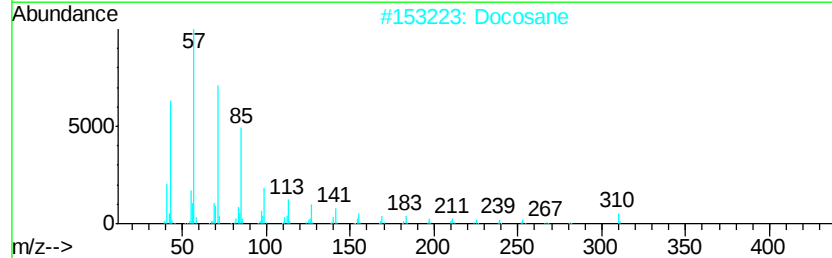
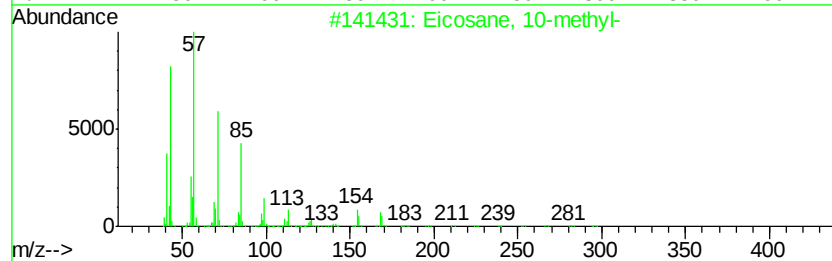
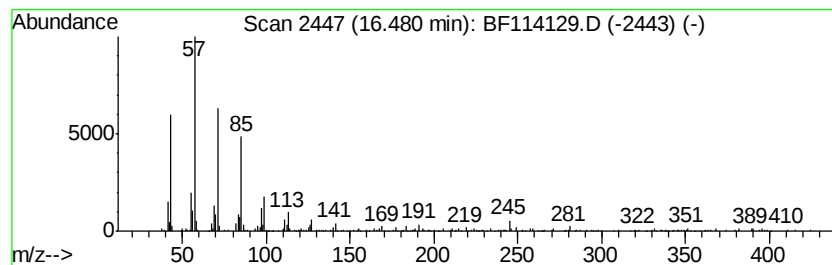
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane, 10-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.48	3.09 ng	317432	Perylene-d12	15.50		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93	
2	Docosane	310	C22H46	000629-97-0	91	
3	Octacosane	394	C28H58	000630-02-4	90	
4	Tetracosane	338	C24H50	000646-31-1	87	
5	Heneicosane	296	C21H44	000629-94-7	86	



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
Data File : BF114129.D
Acq On : 10 May 2019 19:15
Operator : JU/SJ
Sample : K2762-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P021-SS001-1824-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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
Butane, 2-methoxy...	2.13	18.1	ng	1891440	1	6.85	2094130	20.0
unknown6.62	6.62	68.8	ng	7209300	1	6.85	2094130	20.0
n-Hexadecanoic acid	11.90	8.4	ng	1004340	4	11.37	2396070	20.0
Octadecanoic acid	12.67	2.4	ng	285574	4	11.37	2396070	20.0
Octacosanol	13.85	11.3	ng	1217090	5	14.02	2155460	20.0
Octadecane	14.50	3.7	ng	394743	5	14.02	2155460	20.0
Triacontane	14.82	3.3	ng	341746	6	15.50	2052830	20.0
Octadecane, 2-met...	15.16	8.4	ng	864777	6	15.50	2052830	20.0
Eicosane	15.97	7.5	ng	771337	6	15.50	2052830	20.0
Eicosane, 10-methyl-	16.48	3.1	ng	317432	6	15.50	2052830	20.0