

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031120\
 Data File : BF119350.D
 Acq On : 11 Mar 2020 16:56
 Operator : CG/JU
 Sample : L1864-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2

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-BF022020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.910	477	480	497	rBV	171084	246013	8.93%	1.184%
2	5.345	551	554	586	rBV	1243435	1649031	59.86%	7.935%
3	6.375	725	729	757	rBV	1288149	1709300	62.05%	8.225%
4	6.716	784	787	798	rBV	597397	526668	19.12%	2.534%
5	6.869	810	813	827	rBV	1426993	1391221	50.50%	6.695%
6	7.280	880	883	905	rBV	997178	1131406	41.07%	5.444%
7	7.998	1002	1005	1020	rBV	738182	697561	25.32%	3.357%
8	9.074	1184	1188	1199	rBV	2290907	2124948	77.14%	10.226%
9	9.474	1252	1256	1262	rBV	137742	146859	5.33%	0.707%
10	9.751	1299	1303	1314	rVB	1033982	904262	32.82%	4.351%
11	10.545	1434	1438	1453	rBV	1643387	1644008	59.68%	7.911%
12	11.239	1552	1556	1565	rBV	1028673	988208	35.87%	4.755%
13	11.763	1640	1645	1658	rBV2	125679	187481	6.81%	0.902%
14	12.463	1759	1764	1772	rBV	21630	35486	1.29%	0.171%
15	12.551	1775	1779	1786	rBV2	43130	63777	2.32%	0.307%
16	12.815	1820	1824	1828	rBV	2960276	2754812	100.00%	13.257%
17	13.386	1916	1921	1924	rBV2	41488	47789	1.73%	0.230%
18	13.715	1972	1977	1988	rBV	217205	381103	13.83%	1.834%
19	13.880	2001	2005	2015	rBV	849938	848700	30.81%	4.084%
20	14.027	2026	2030	2034	rBV	118338	112020	4.07%	0.539%
21	14.333	2078	2082	2085	rBV	210174	199316	7.24%	0.959%
22	14.627	2127	2132	2136	rBV	335248	332865	12.08%	1.602%
23	14.945	2180	2186	2191	rBV	372387	451927	16.41%	2.175%
24	15.309	2238	2248	2263	rBV2	607806	1263586	45.87%	6.081%
25	15.698	2308	2314	2320	rVB	331077	473603	17.19%	2.279%
26	16.162	2386	2393	2406	rVB	175144	330690	12.00%	1.591%
27	16.703	2479	2485	2492	rVB2	50414	102374	3.72%	0.493%
28	17.339	2590	2593	2602	rVB8	15459	35825	1.30%	0.172%

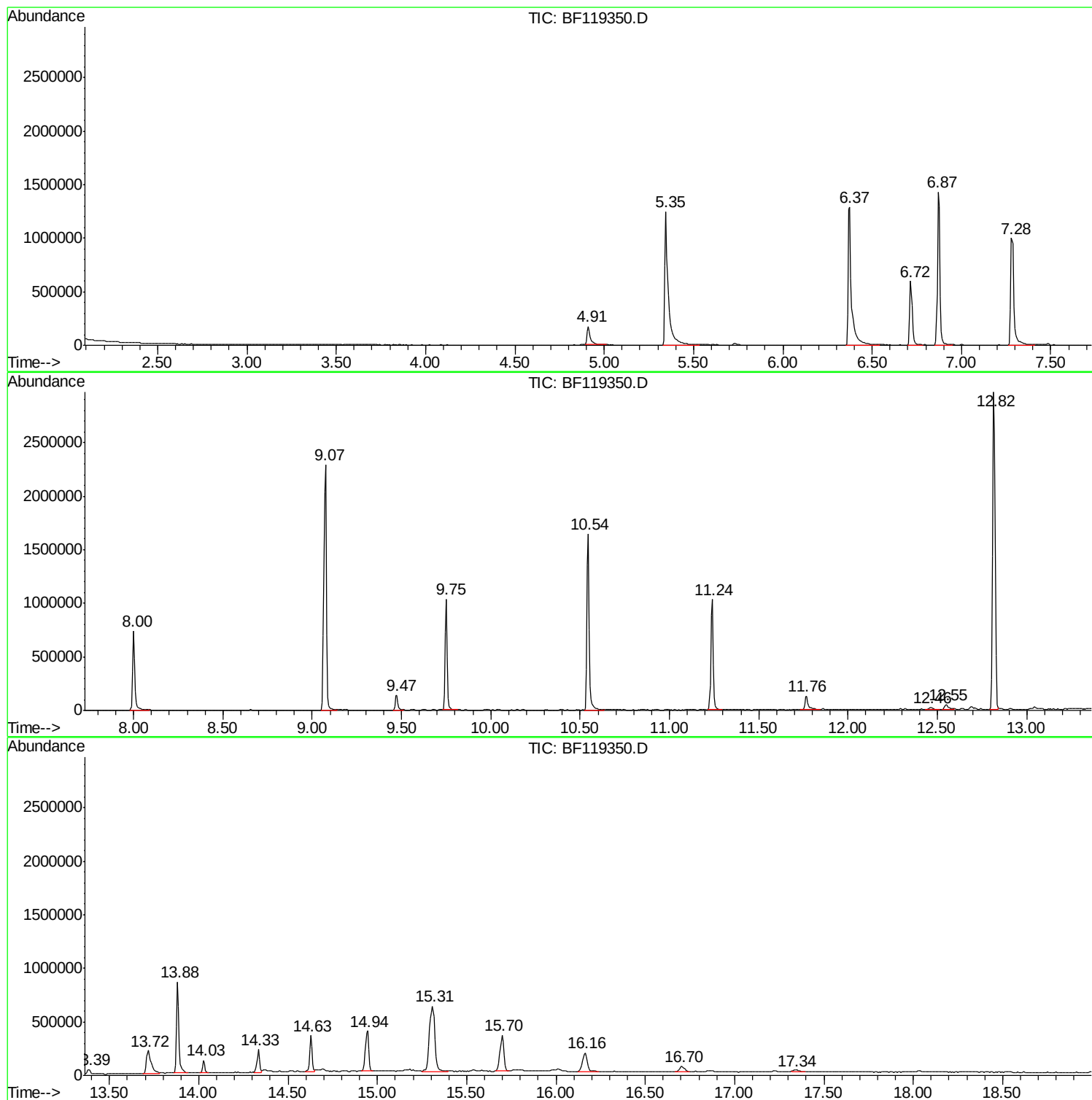
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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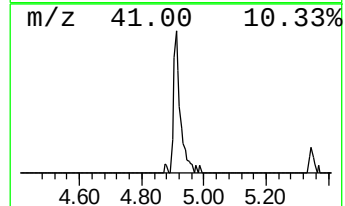
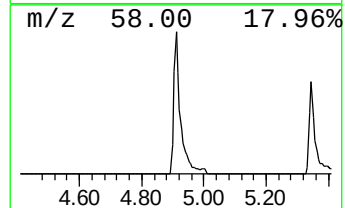
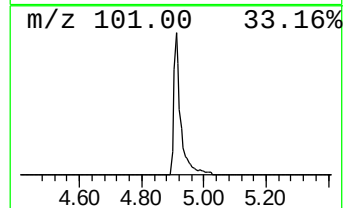
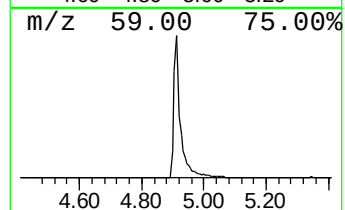
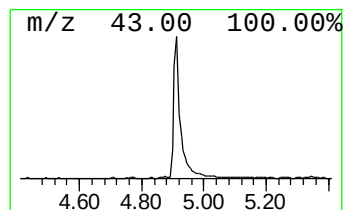
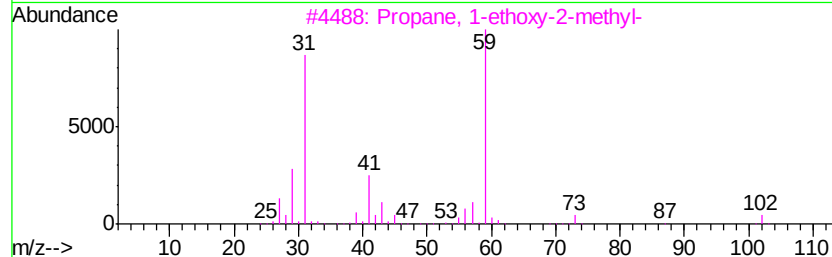
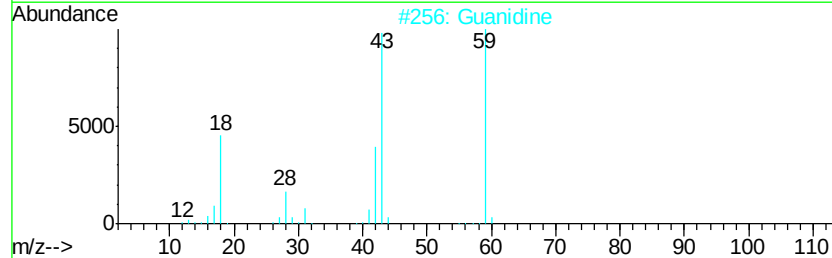
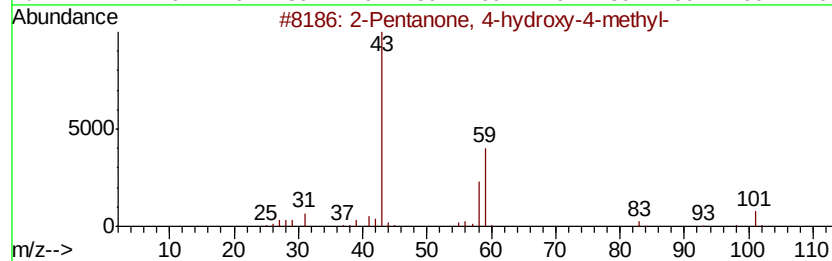
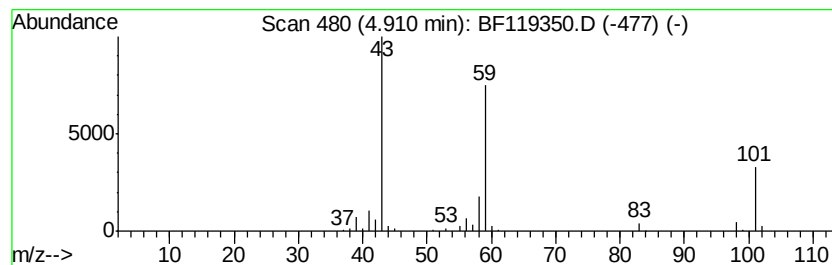
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.91	9.34 ng	246013	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Guanidine	59	CH5N3	000113-00-8	50
3			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38
4			2,3-Dimethyloxirane-2-carboxylic...	130	C6H10O3	1000306-64-4	36
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	35



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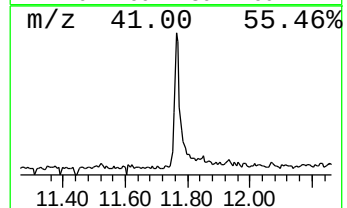
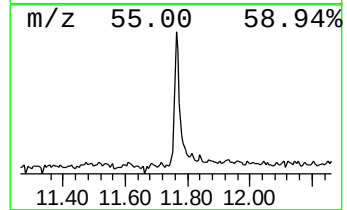
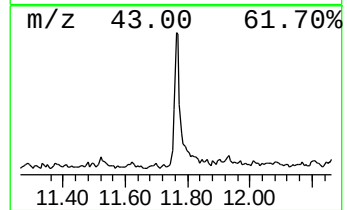
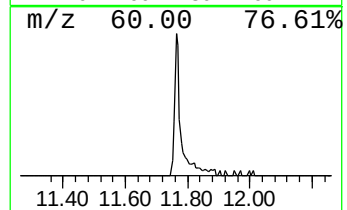
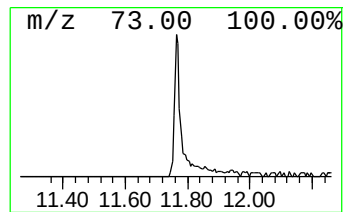
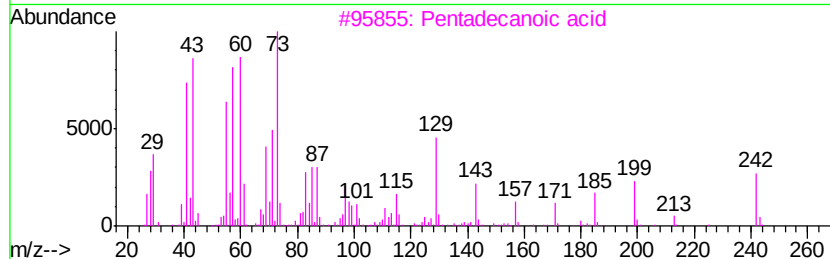
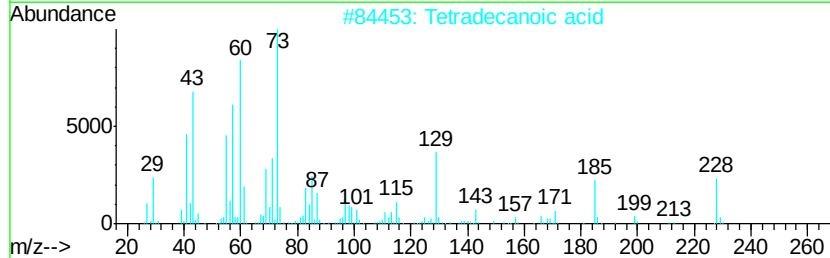
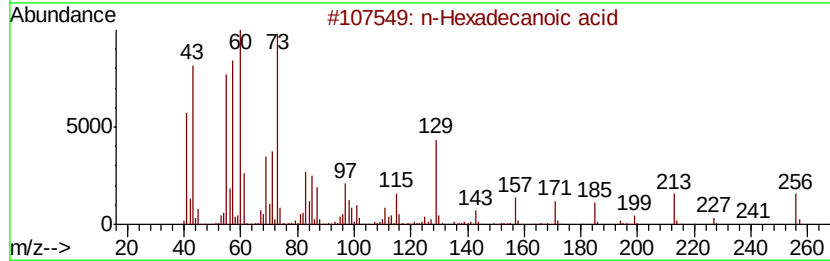
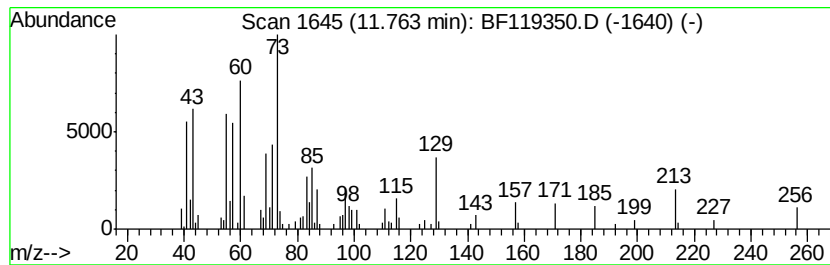
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.76	3.79 ng	187481	Phenanthrene-d10	11.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	Nonanoic acid	158	C9H18O2	000112-05-0	30



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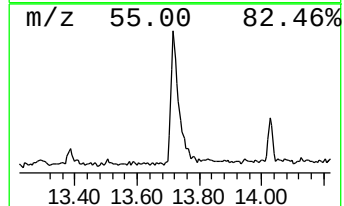
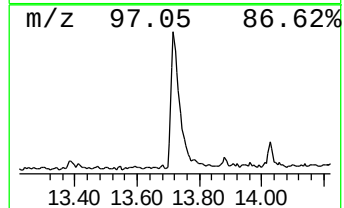
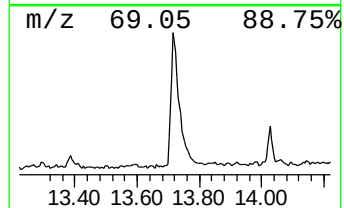
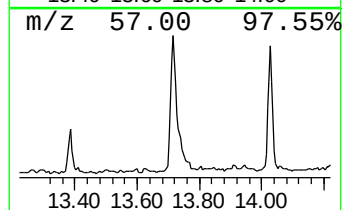
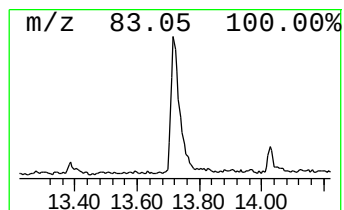
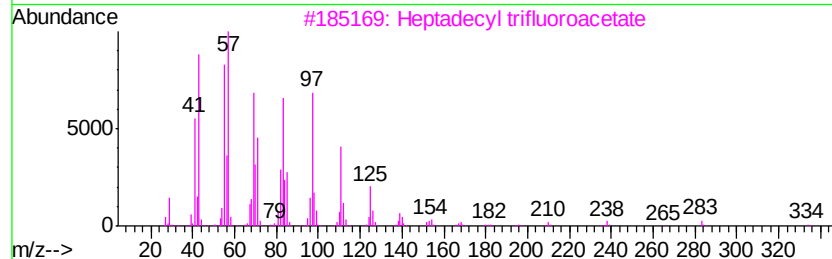
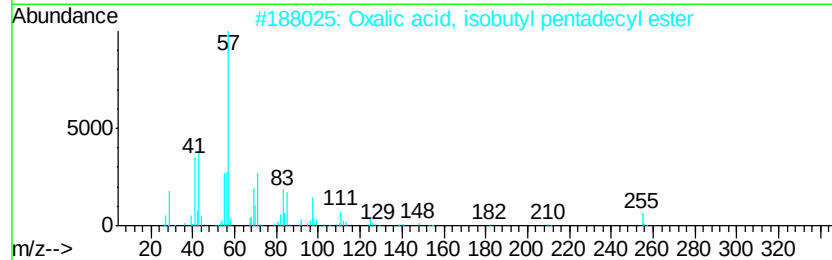
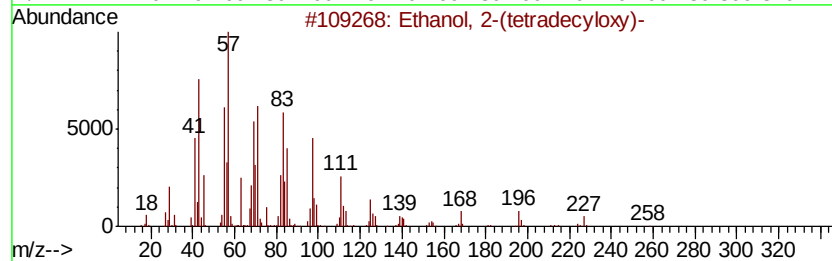
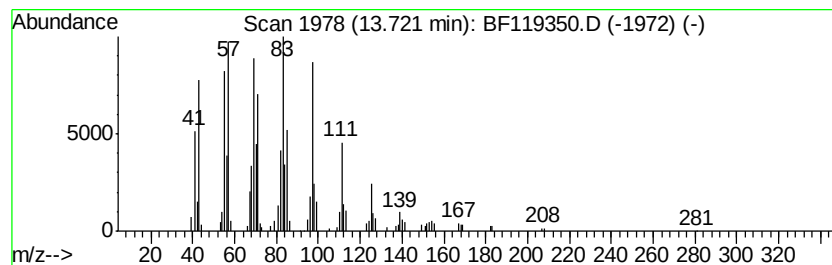
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Peak Number 4 Ethanol, 2-(tetradecyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	8.98 ng	381103	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2		Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	91
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
4		1-Octadecanol	270	C18H38O	000112-92-5	91
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	89



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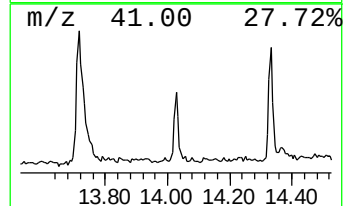
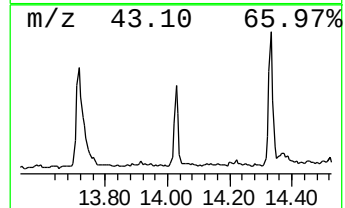
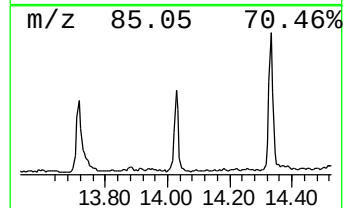
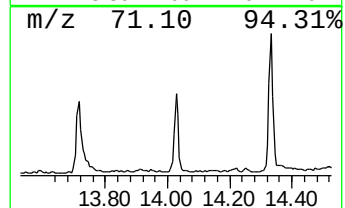
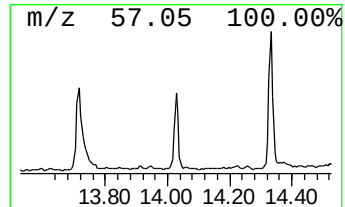
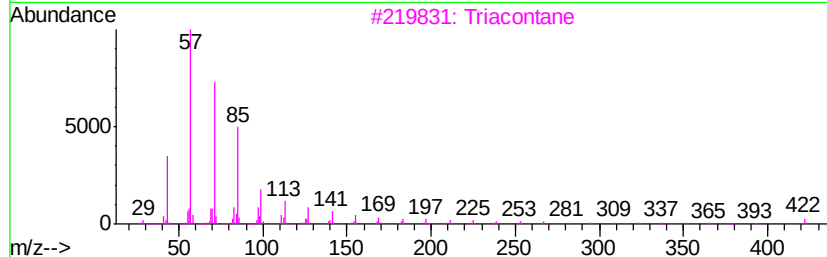
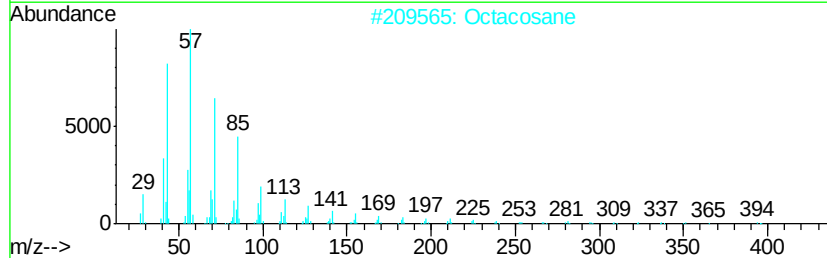
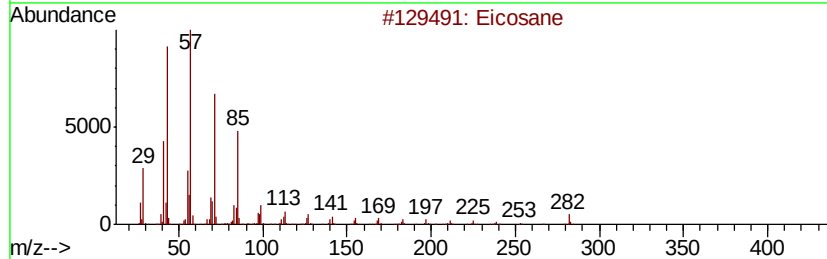
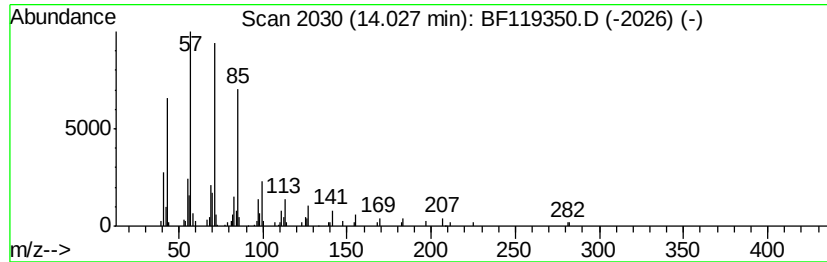
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Peak Number 5 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.03	2.64 ng	112020	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



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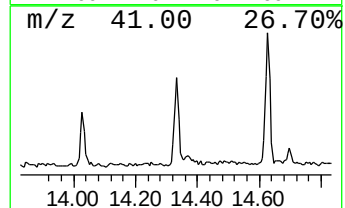
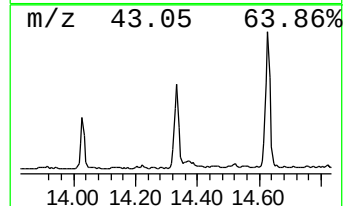
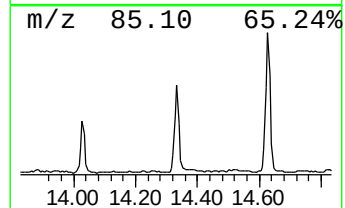
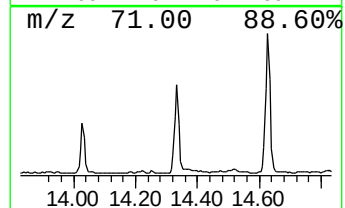
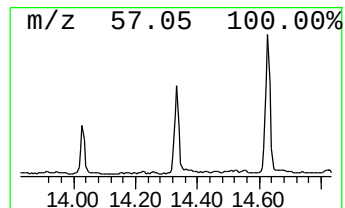
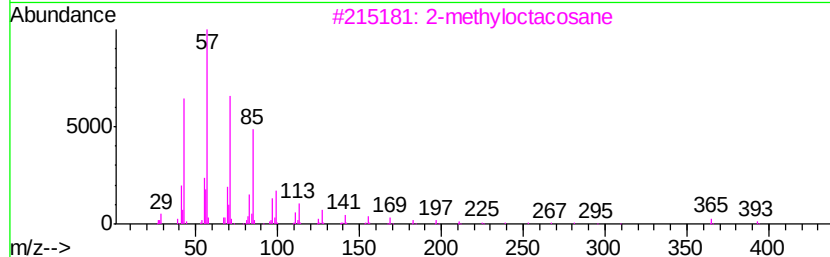
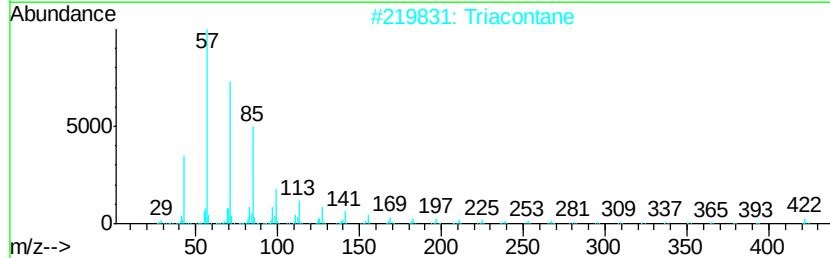
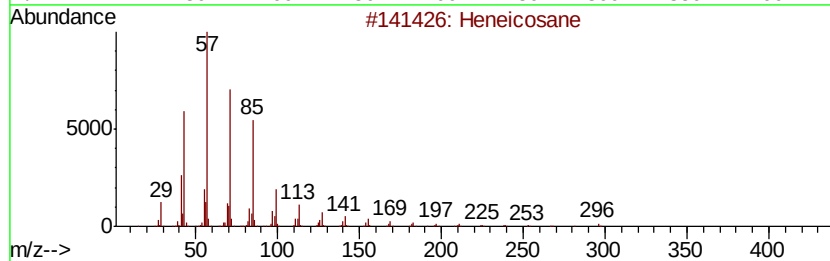
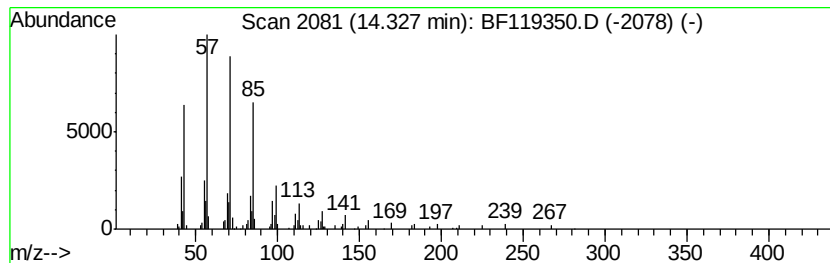
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	4.70 ng	199316	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Triacontane	422	C30H62	000638-68-6	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Eicosane	282	C20H42	000112-95-8	91
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



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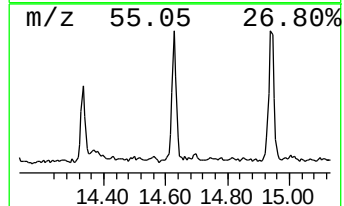
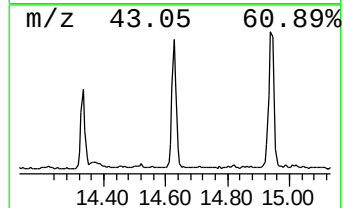
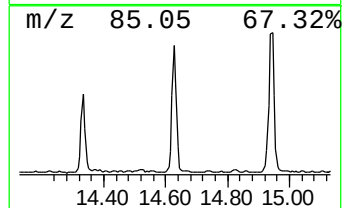
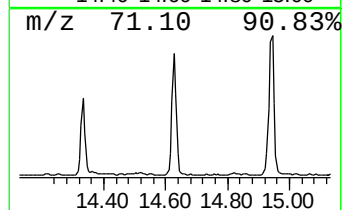
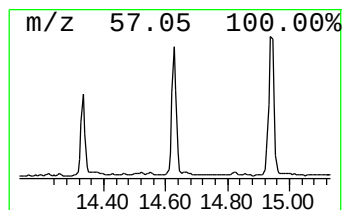
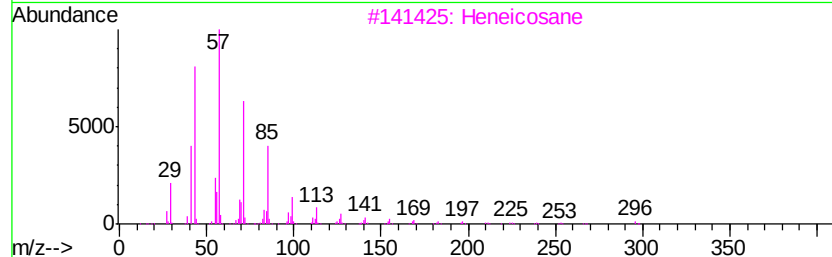
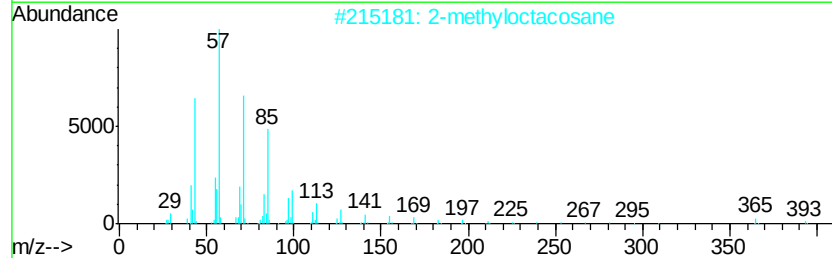
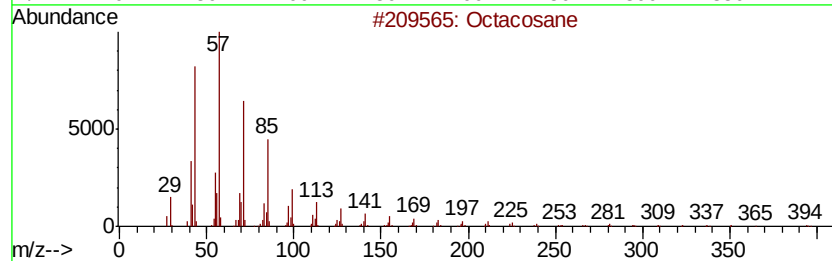
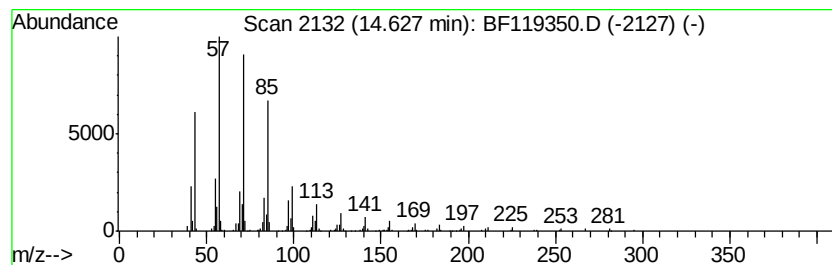
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	5.27 ng	332865	Perylene-d12	15.31

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	91
2	2-methyloctacosane		408	C29H60	1000376-72-8	91
3	Heneicosane		296	C21H44	000629-94-7	91
4	Heptacosane		380	C27H56	000593-49-7	91
5	Triacontane		422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119350.D
Acq On : 11 Mar 2020 16:56
Operator : CG/JU
Sample : L1864-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-2

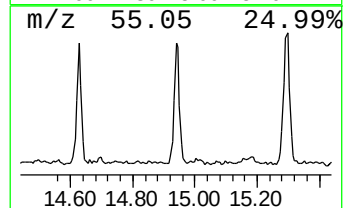
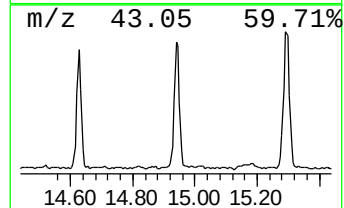
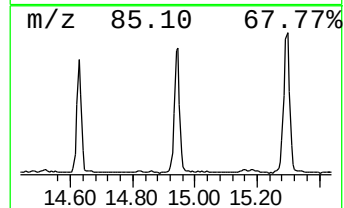
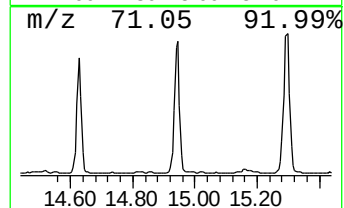
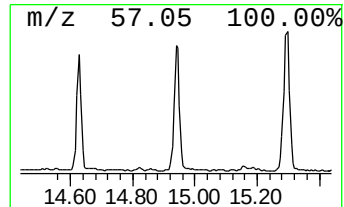
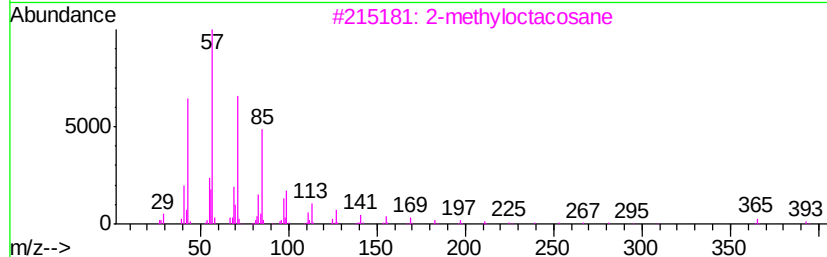
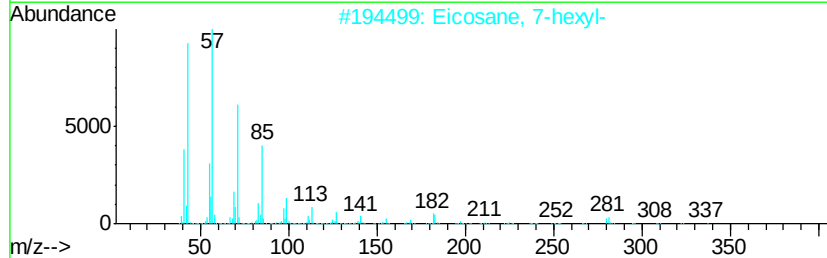
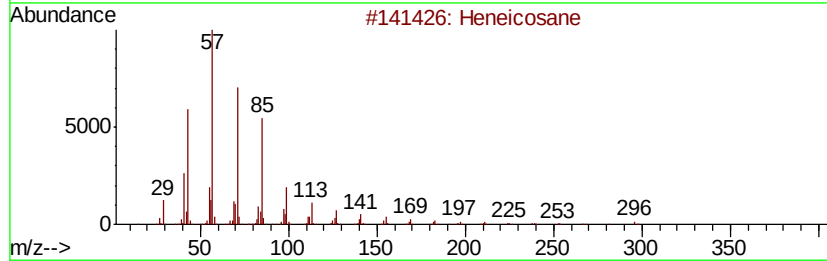
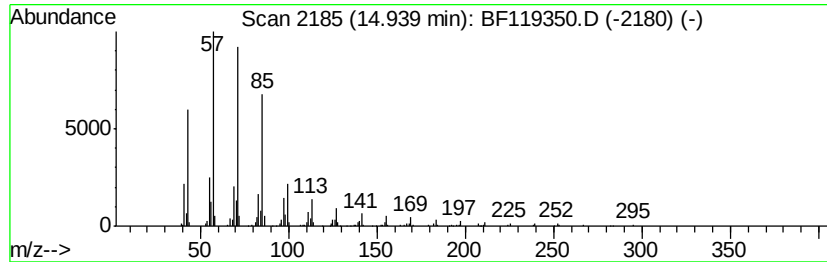
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane, 7-hexyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	7.15 ng	451927	Perylene-d12	15.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119350.D
Acq On : 11 Mar 2020 16:56
Operator : CG/JU
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Misc :
ALS Vial : 4 Sample Multiplier: 1

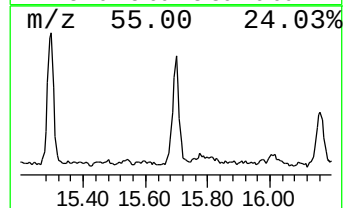
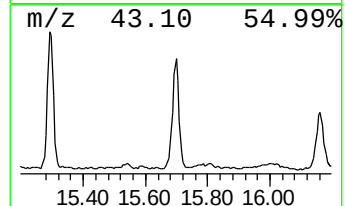
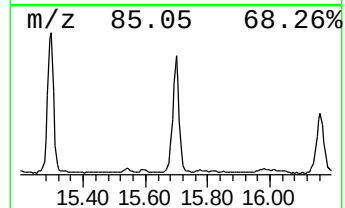
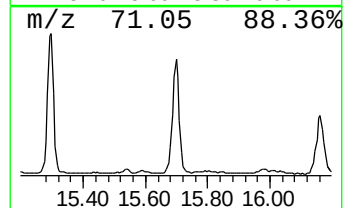
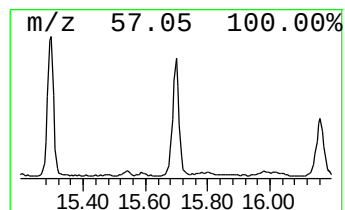
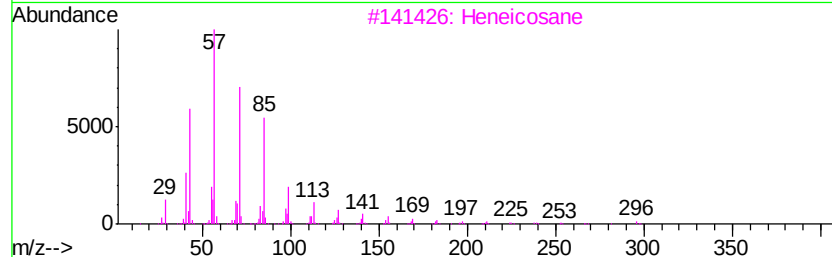
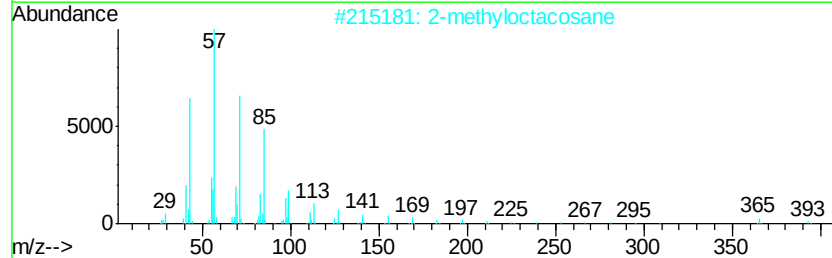
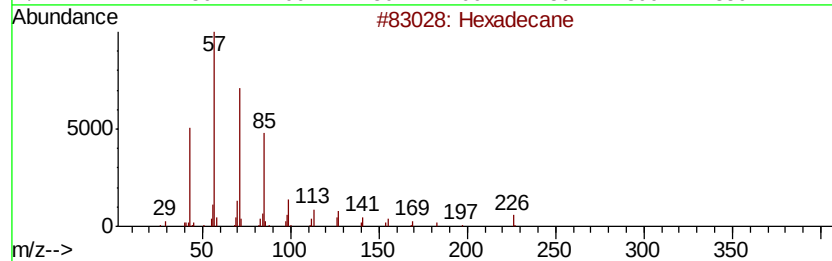
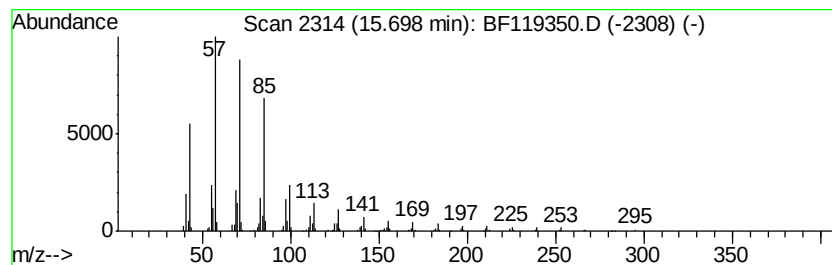
Instrument :
BNA_F
ClientSampleId :
SP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.70	7.50 ng	473603	Perylene-d12	15.31	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031120\
Data File : BF119350.D
Acq On : 11 Mar 2020 16:56
Operator : CG/JU
Sample : L1864-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-2

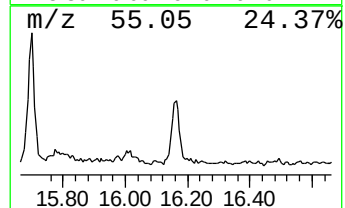
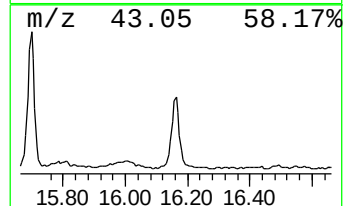
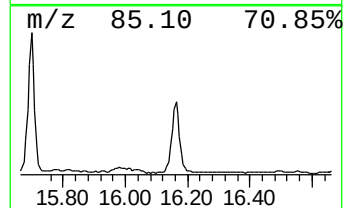
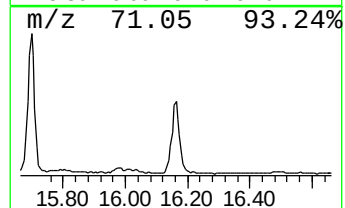
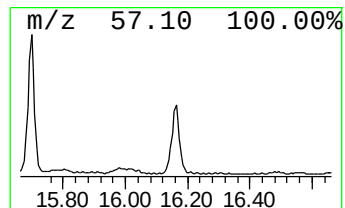
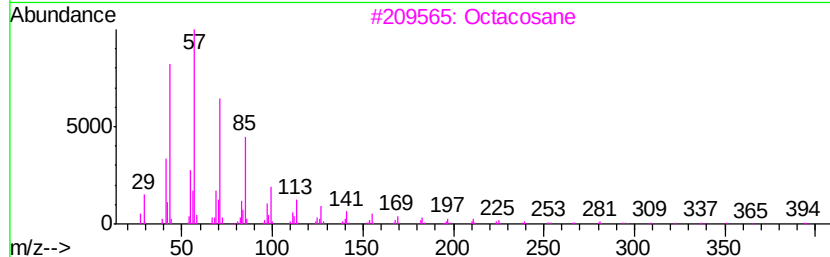
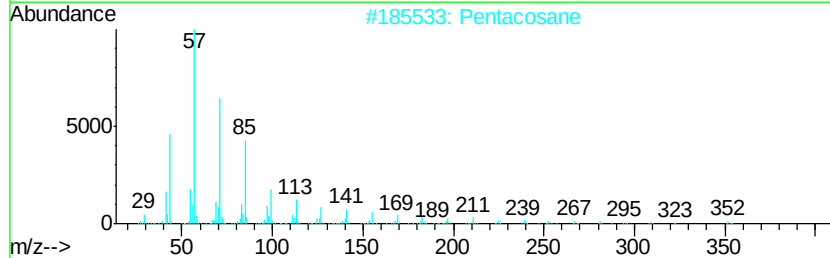
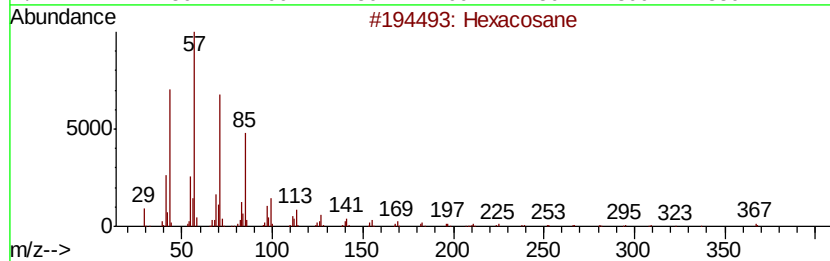
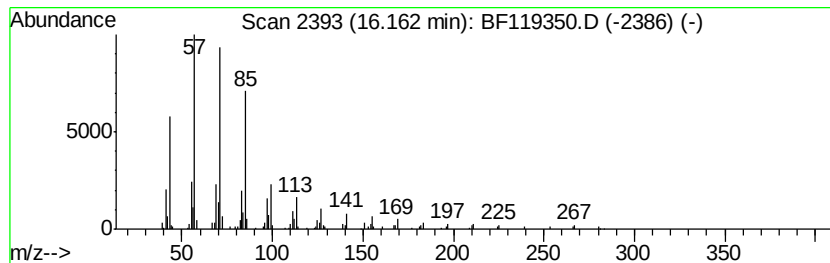
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	5.23 ng	330690	Perylene-d12	15.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Pentacosane	352	C25H52	000629-99-2	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031120\
Data File : BF119350.D
Acq On : 11 Mar 2020 16:56
Operator : CG/JU
Sample : L1864-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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
2-Pentanone, 4-hy...	4.91	9.3	ng	246013	1	6.72	526668	20.0
n-Hexadecanoic acid	11.76	3.8	ng	187481	4	11.24	988208	20.0
Ethanol, 2-(tetra...	13.72	9.0	ng	381103	5	13.88	848700	20.0
Eicosane	14.03	2.6	ng	112020	5	13.88	848700	20.0
Heneicosane	14.33	4.7	ng	199316	5	13.88	848700	20.0
Octacosane	14.63	5.3	ng	332865	6	15.31	1263590	20.0
Eicosane, 7-hexyl-	14.94	7.2	ng	451927	6	15.31	1263590	20.0
Hexadecane	15.70	7.5	ng	473603	6	15.31	1263590	20.0
Hexacosane	16.16	5.2	ng	330690	6	15.31	1263590	20.0