

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
 Data File : BF120249.D
 Acq On : 30 Apr 2020 19:45
 Operator : MA/SJ
 Sample : L2452-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 224

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-BF040820.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	5.257	535	539	554	rBV	3379680	3307089	74.73%	9.374%
2	6.310	713	718	740	rVB	3124092	3470554	78.43%	9.837%
3	6.492	746	749	758	rVB	44684	46430	1.05%	0.132%
4	6.645	772	775	779	rBV	752097	678929	15.34%	1.924%
5	6.804	798	802	805	rBV	3254547	2829561	63.94%	8.021%
6	7.222	868	873	875	rBV	2345799	2300281	51.98%	6.520%
7	7.934	990	994	997	rBV	1172820	885236	20.00%	2.509%
8	8.039	1009	1012	1023	rBV	353529	344320	7.78%	0.976%
9	9.016	1174	1178	1181	rBV	4433535	4146110	93.69%	11.752%
10	9.416	1242	1246	1249	rBV	234761	200744	4.54%	0.569%
11	9.692	1289	1293	1296	rBV	1170562	980442	22.16%	2.779%
12	10.486	1423	1428	1431	rBV	2830024	2612665	59.04%	7.406%
13	11.175	1541	1545	1548	rBV	1275787	1050819	23.75%	2.979%
14	11.727	1634	1639	1647	rBV	1335820	1375321	31.08%	3.898%
15	12.510	1768	1772	1785	rBV	892540	951692	21.51%	2.698%
16	12.616	1785	1790	1793	rVB	49441	61393	1.39%	0.174%
17	12.774	1811	1817	1820	rBV	4292012	4425187	100.00%	12.543%
18	13.351	1906	1915	1918	rBV3	41104	63782	1.44%	0.181%
19	13.627	1958	1962	1966	rBV2	35008	46872	1.06%	0.133%
20	13.674	1966	1970	1980	rVV	583762	741483	16.76%	2.102%
21	13.816	1989	1994	1997	rBV2	1334121	1288682	29.12%	3.653%
22	13.998	2021	2025	2032	rBV	125776	137651	3.11%	0.390%
23	14.304	2072	2077	2084	rBV	234337	307972	6.96%	0.873%
24	14.474	2100	2106	2111	rBV3	34133	67257	1.52%	0.191%
25	14.598	2123	2127	2130	rBV	315508	303196	6.85%	0.859%
26	14.910	2176	2180	2184	rBV	358976	385902	8.72%	1.094%
27	15.221	2228	2233	2236	rBV	926933	1031007	23.30%	2.922%
28	15.257	2236	2239	2247	rVB	335614	409160	9.25%	1.160%
29	15.657	2302	2307	2312	rBV	244272	338821	7.66%	0.960%
30	15.710	2312	2316	2323	rBV3	86936	166607	3.76%	0.472%
31	16.115	2380	2385	2389	rBV	124181	219568	4.96%	0.622%
32	16.657	2471	2477	2485	rBV3	45936	104170	2.35%	0.295%

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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-BF040820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

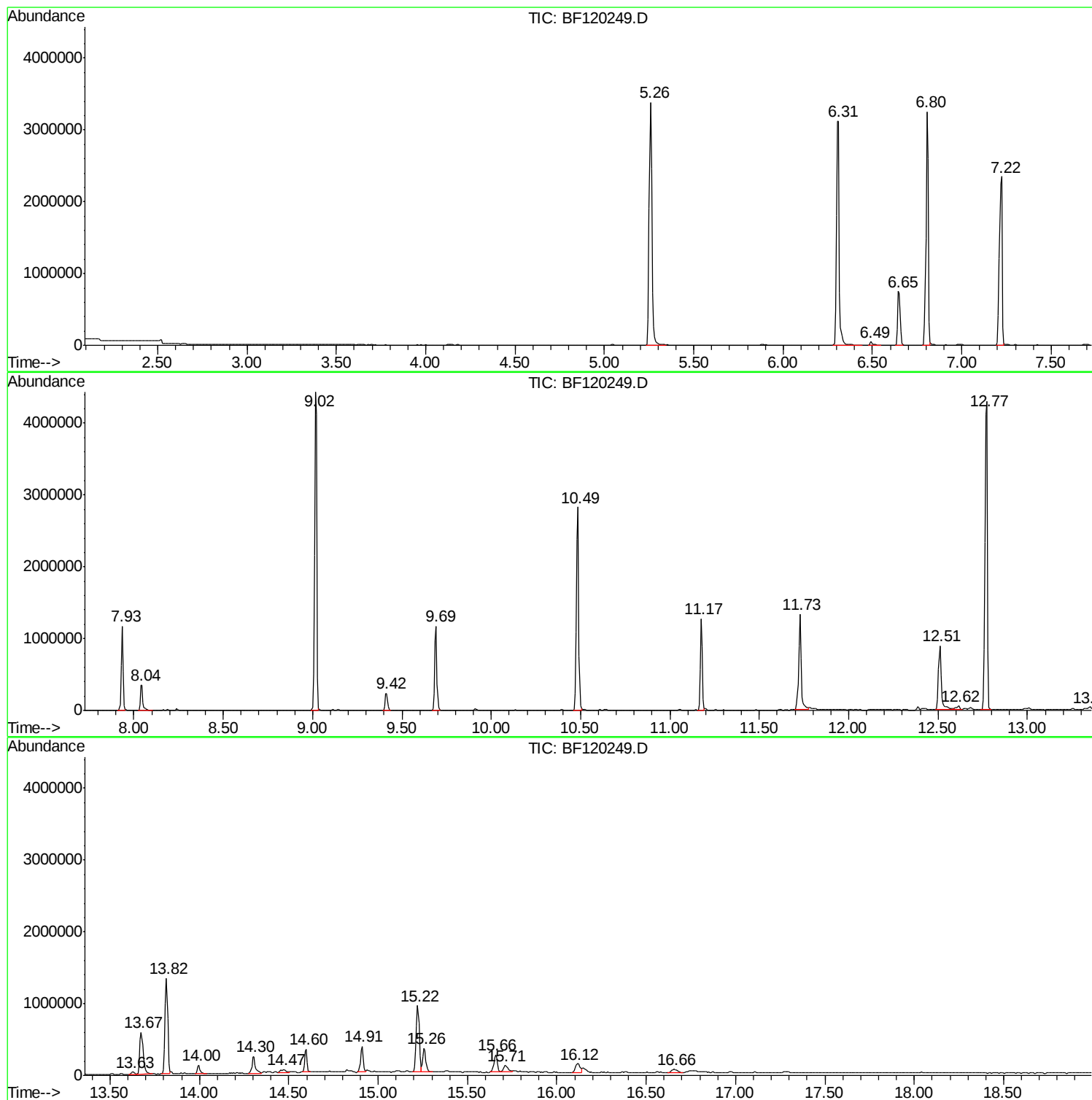
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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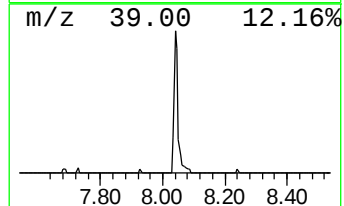
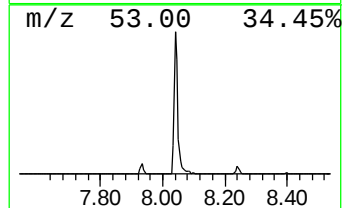
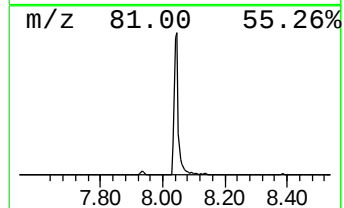
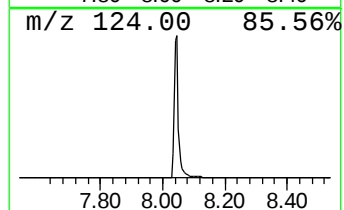
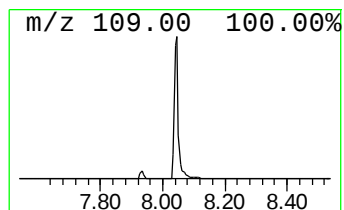
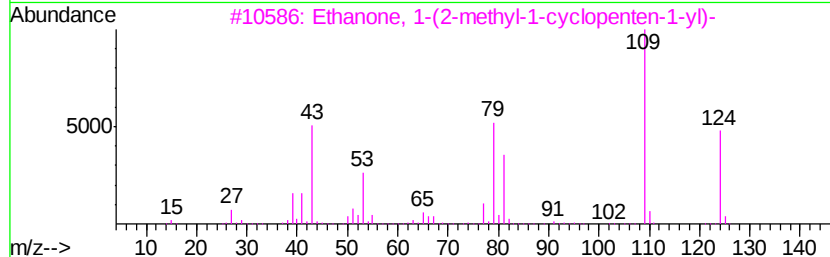
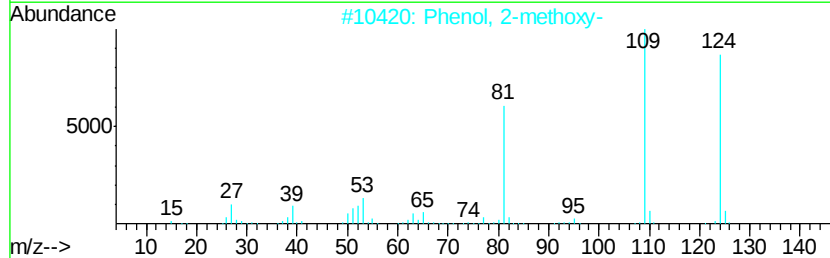
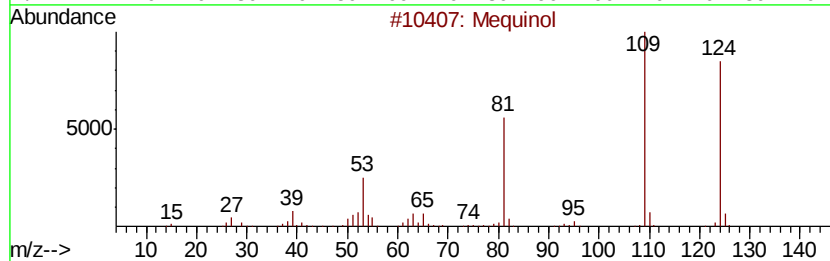
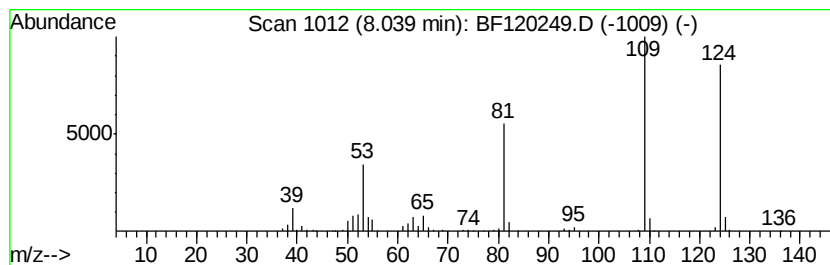
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Mequinol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	7.78 ng	344320	Naphthalene-d8	7.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mequinol		124	C7H8O2	000150-76-5	95
2	Phenol, 2-methoxy-		124	C7H8O2	000090-05-1	94
3	Ethanone, 1-(2-methyl-1-cyclopenten-1-yl)-		124	C8H12O	003168-90-9	91
4	2-Acetyl-5-methylfuran		124	C7H8O2	001193-79-9	87
5	2-Cyclopenten-1-one, 3,4,5-trimethyl-		124	C8H12O	055683-21-1	78



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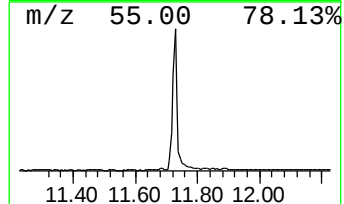
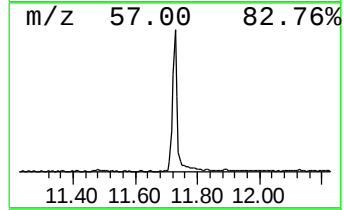
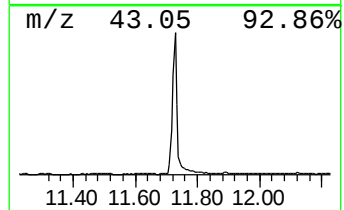
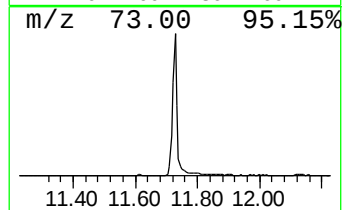
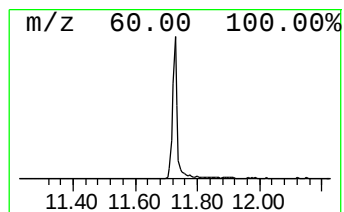
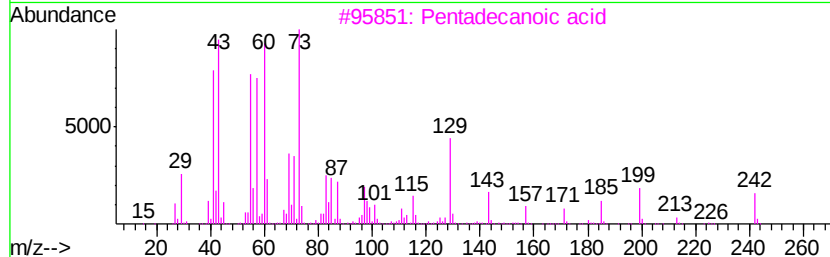
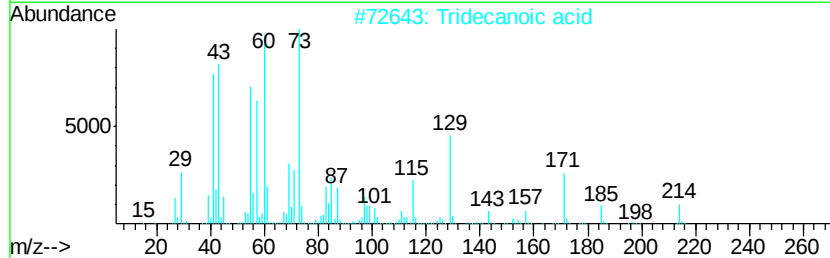
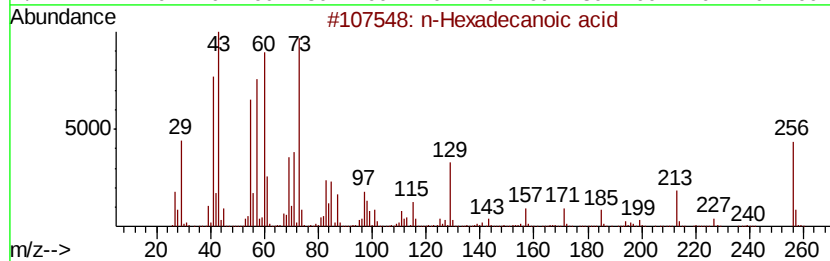
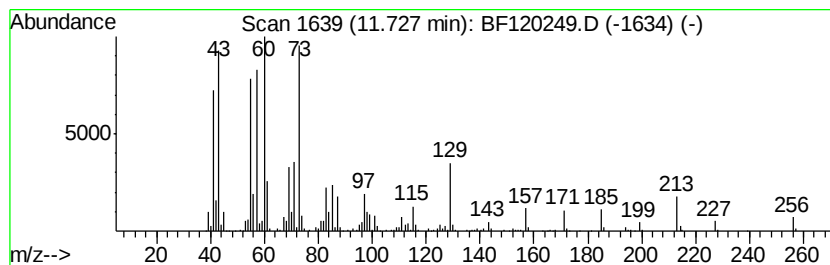
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	26.18 ng	1375320	Phenanthrene-d10	11.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Dodecanoic acid	200	C12H24O2	000143-07-7	53



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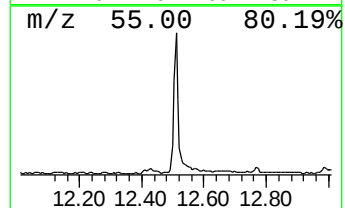
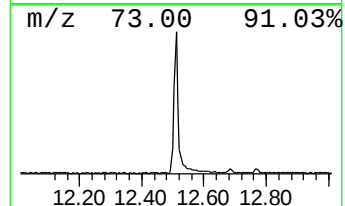
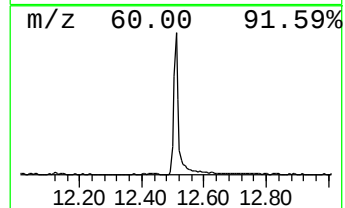
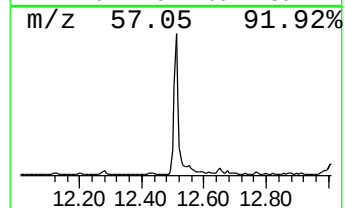
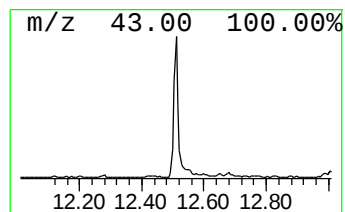
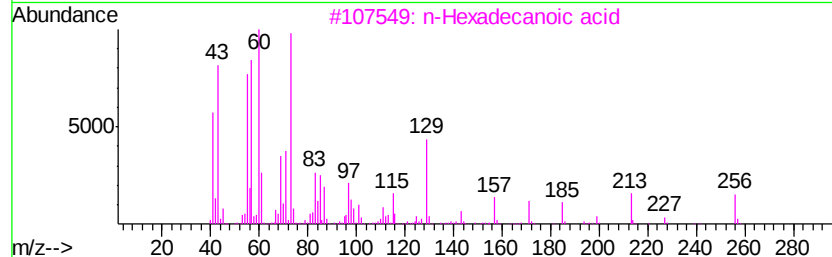
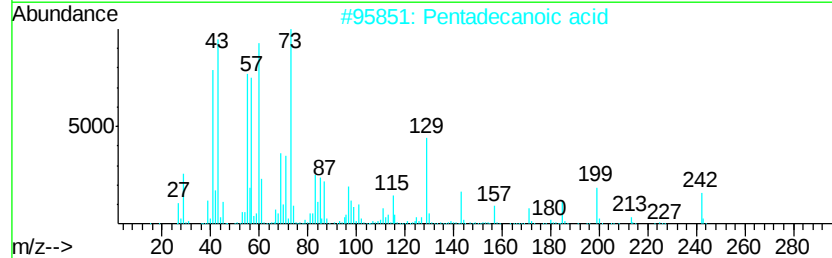
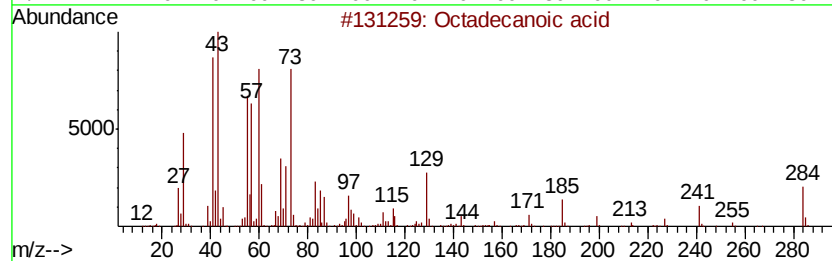
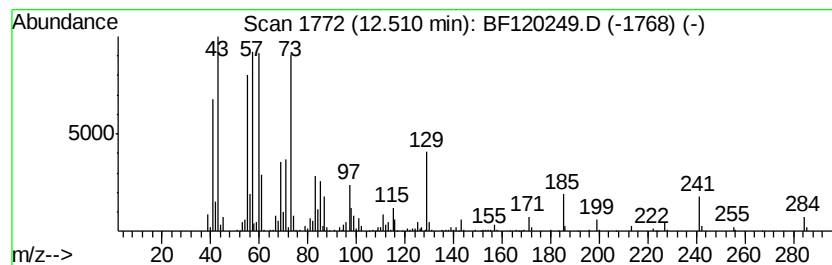
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	14.77 ng	951692	Chrysene-d12	13.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	72



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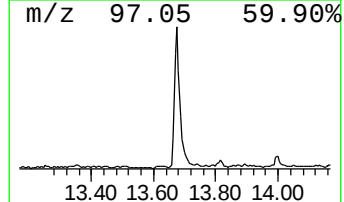
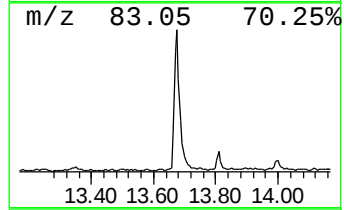
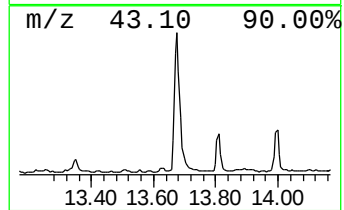
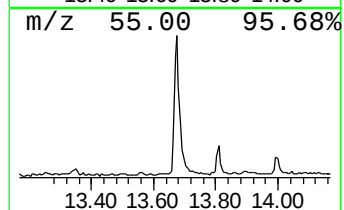
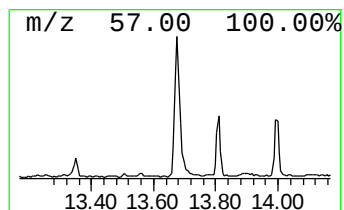
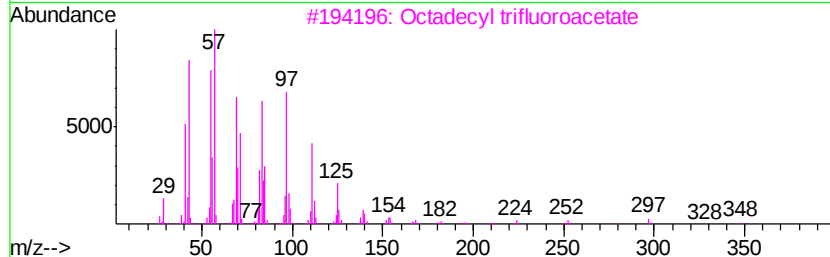
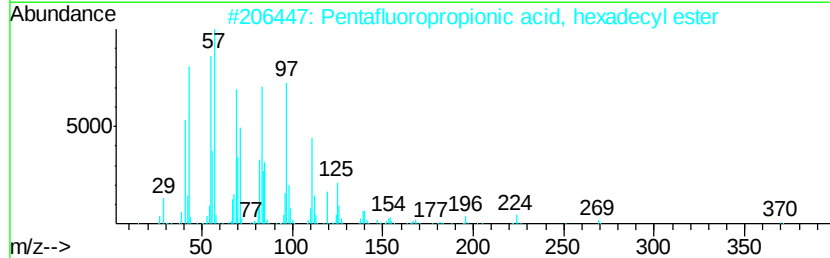
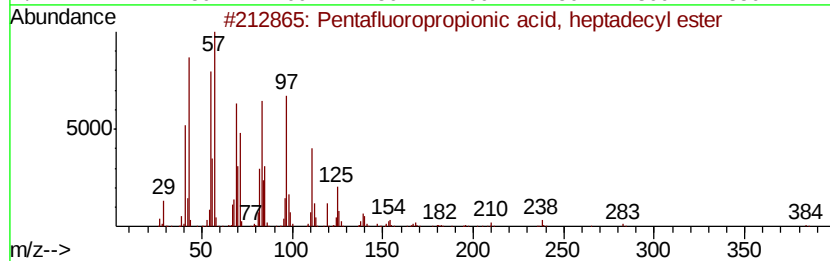
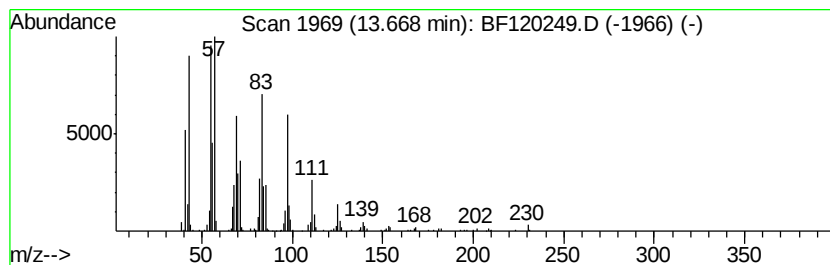
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Pentafluoropropionic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	11.51 ng	741483	Chrysene-d12	13.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	94
2		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
3		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
4		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
5		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91



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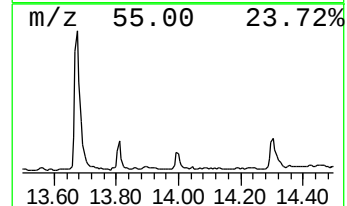
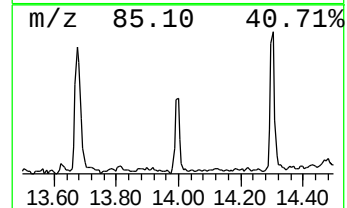
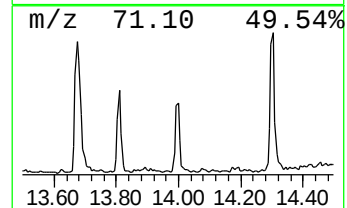
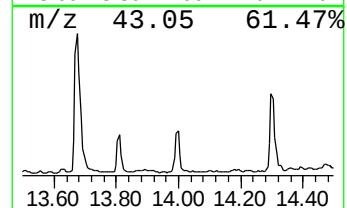
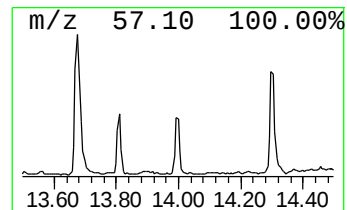
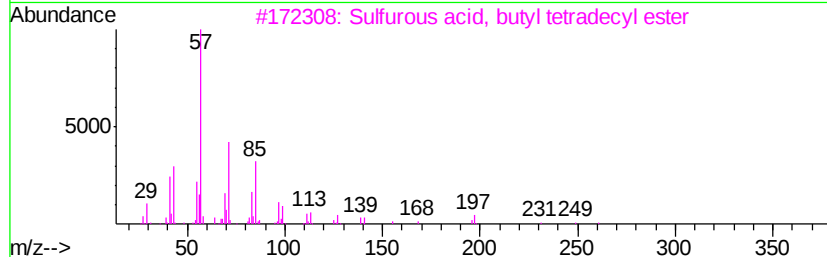
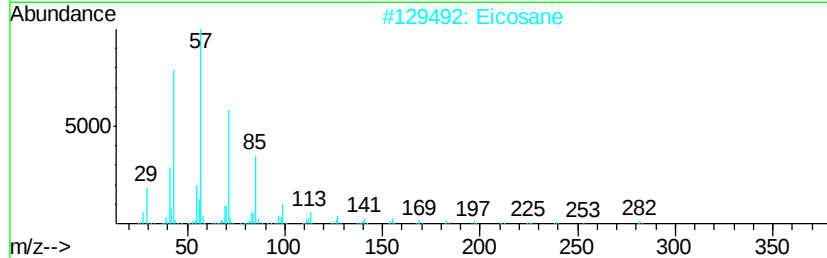
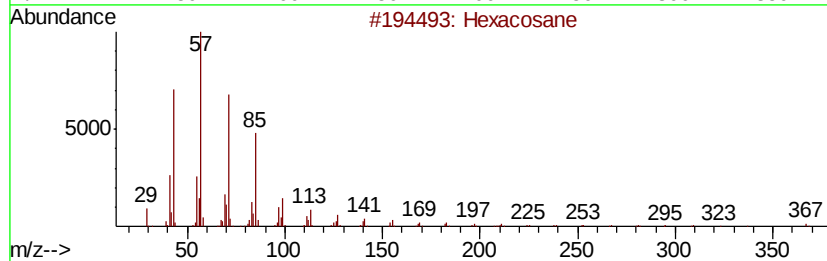
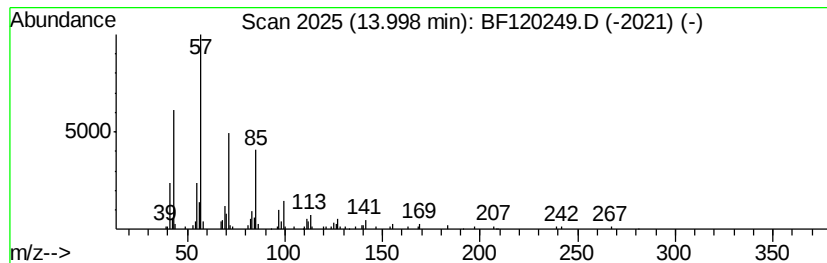
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	2.14 ng	137651	Chrysene-d12	13.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	90
2	Eicosane		282	C20H42	000112-95-8	90
3	Sulfurous acid, butyl tetradecyl...		334	C18H38O3S	1000309-18-1	90
4	Hexadecane		226	C16H34	000544-76-3	87
5	Pentadecane		212	C15H32	000629-62-9	87



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ALS Vial : 18 Sample Multiplier: 1

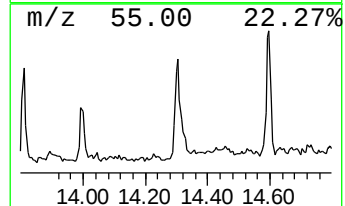
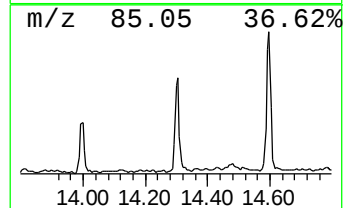
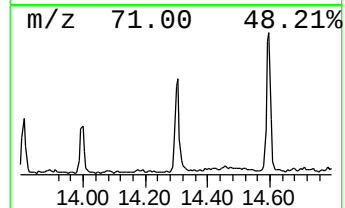
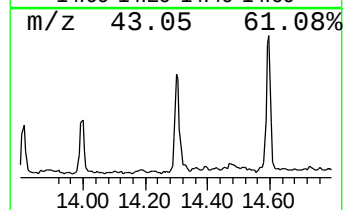
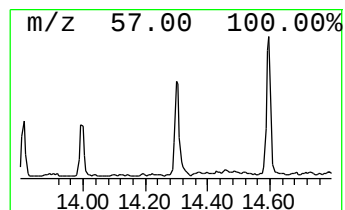
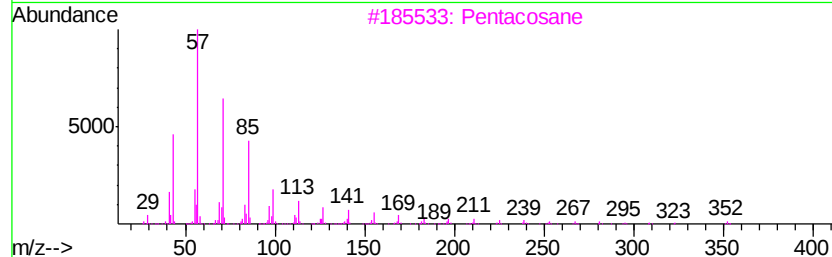
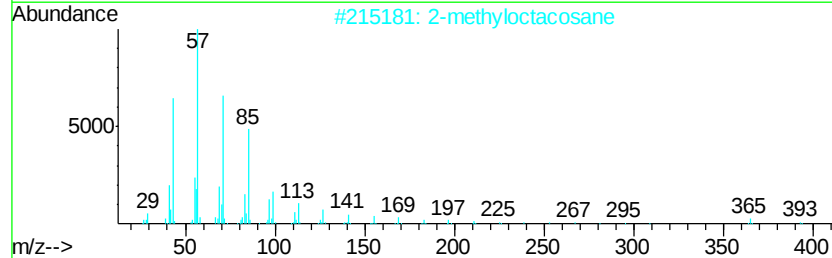
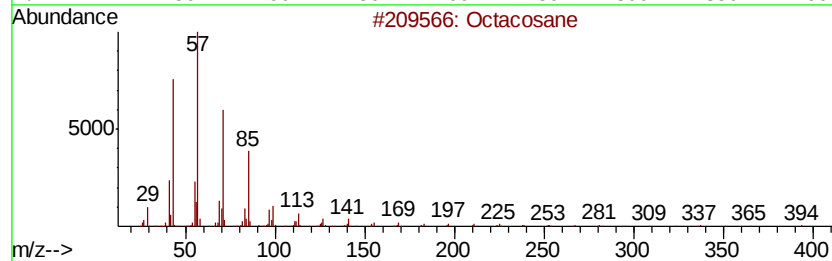
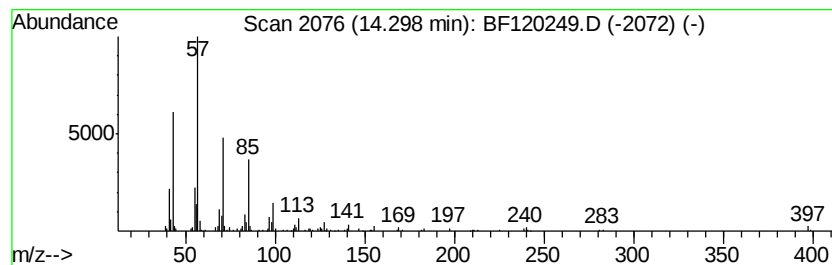
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	4.78 ng	307972	Chrysene-d12	13.82
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	Pentacosane	352 C25H52	000629-99-2	91
4	Docosane, 11-butyl-	366 C26H54	013475-76-8	91
5	Tetratetracontane	619 C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120249.D
Acq On : 30 Apr 2020 19:45
Operator : MA/SJ
Sample : L2452-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

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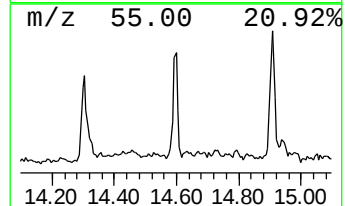
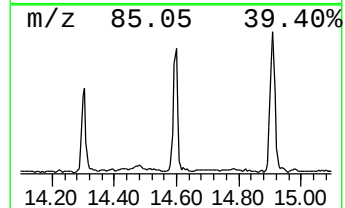
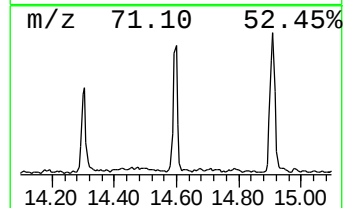
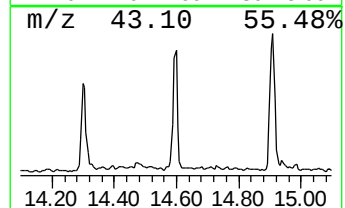
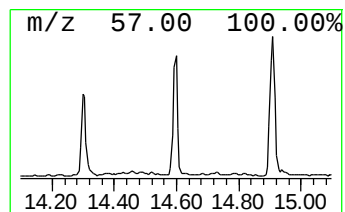
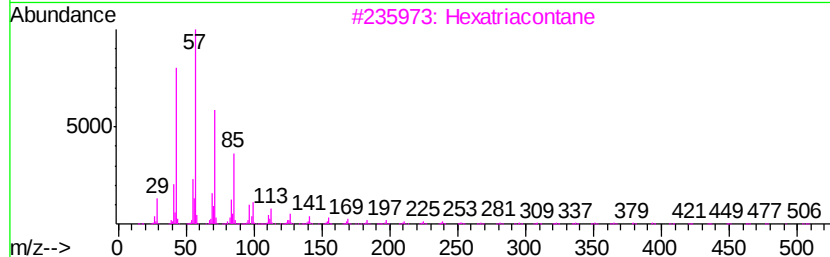
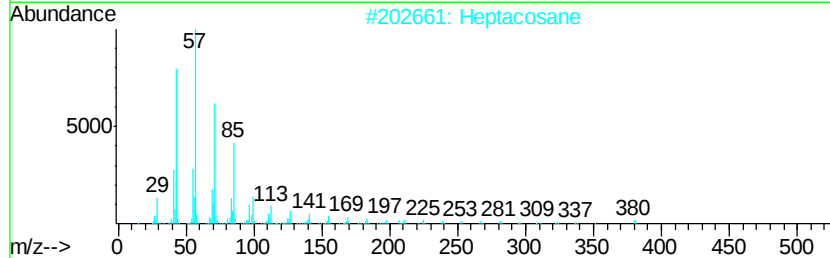
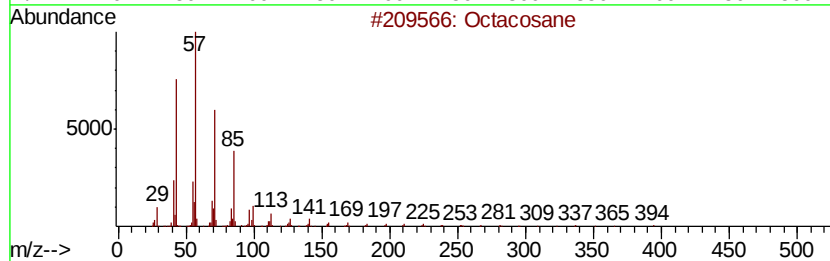
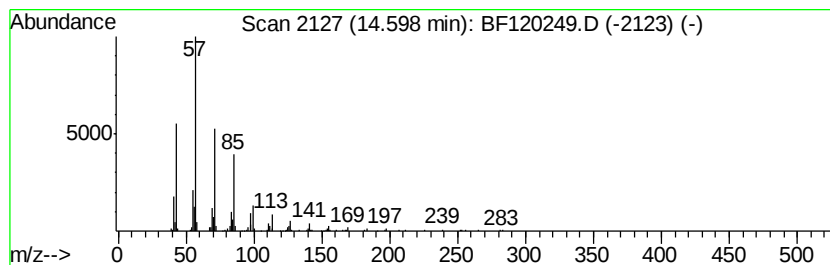
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	5.88 ng	303196	Perylene-d12	15.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Hexatriacontane	507	C36H74	000630-06-8	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120249.D
Acq On : 30 Apr 2020 19:45
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Sample : L2452-03
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ALS Vial : 18 Sample Multiplier: 1

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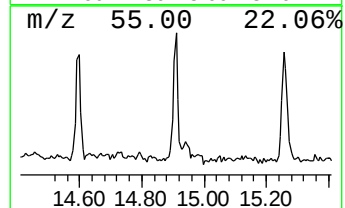
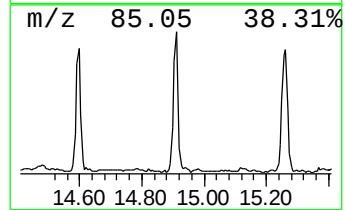
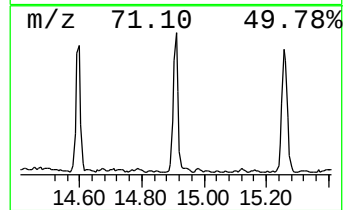
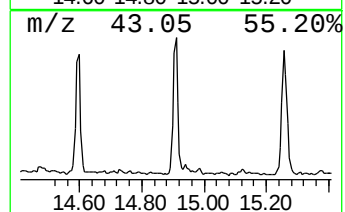
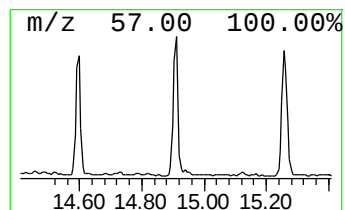
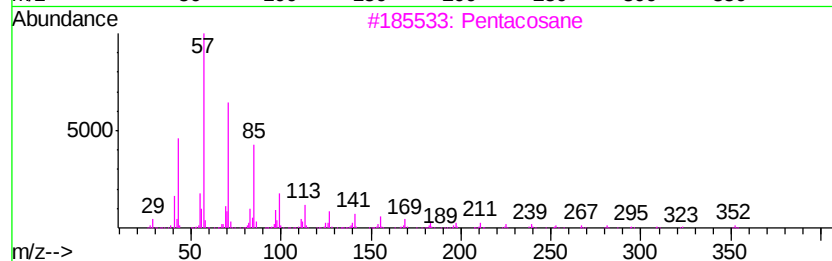
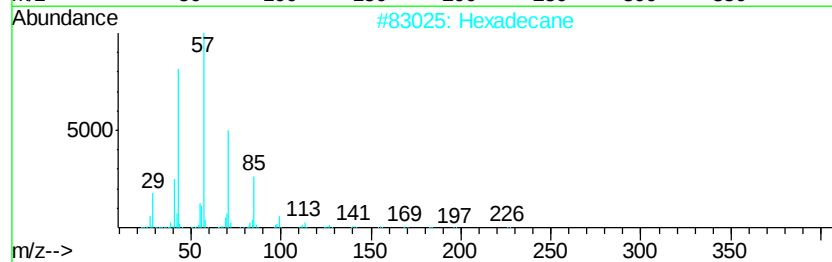
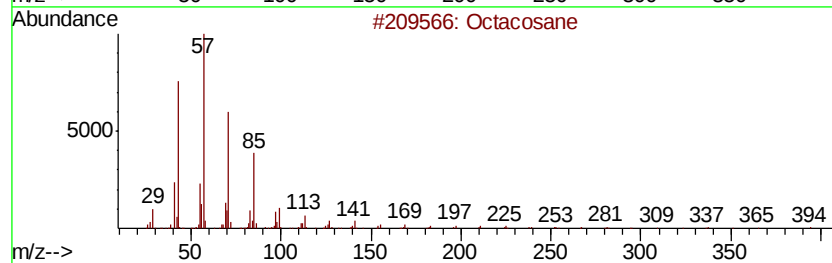
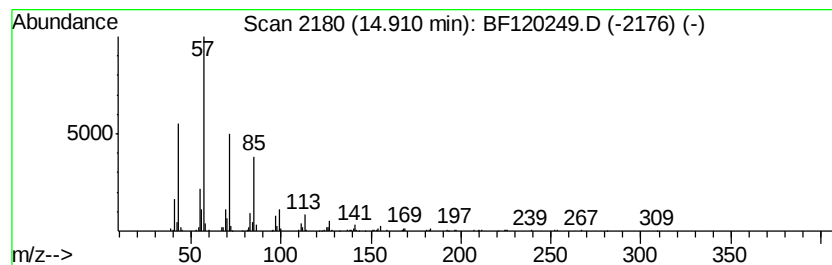
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	7.49 ng	385902	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Hexadecane	226	C16H34	000544-76-3	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120249.D
Acq On : 30 Apr 2020 19:45
Operator : MA/SJ
Sample : L2452-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

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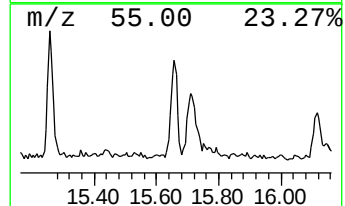
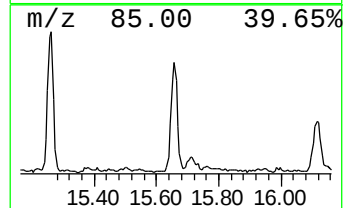
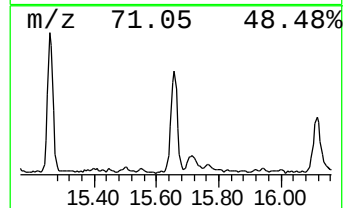
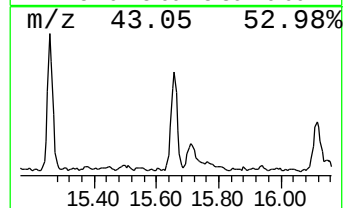
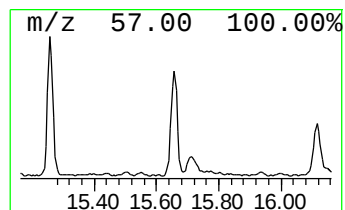
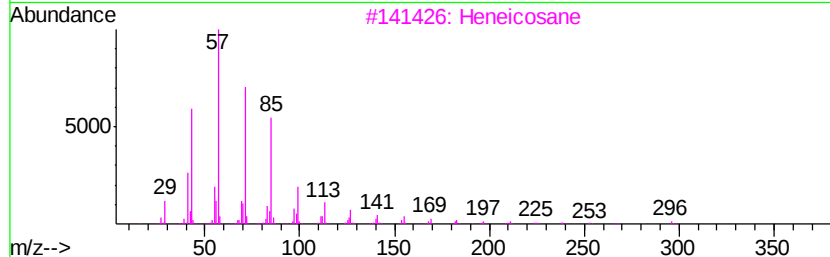
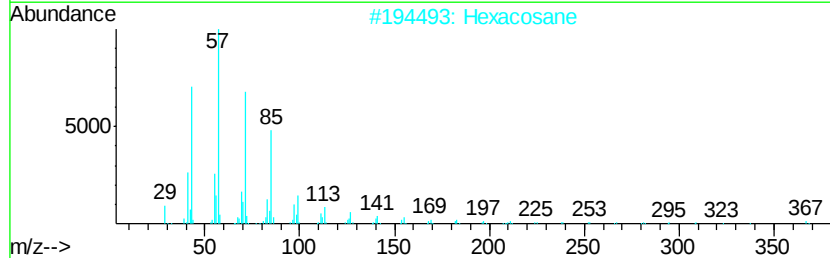
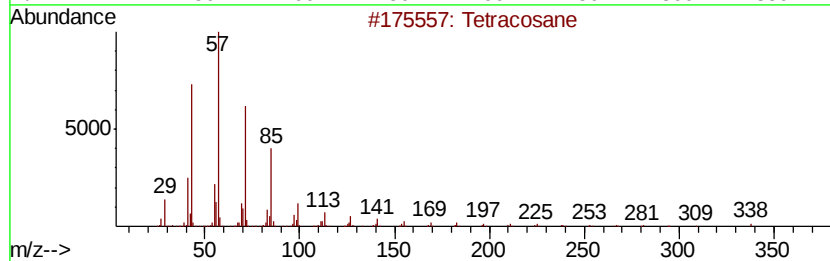
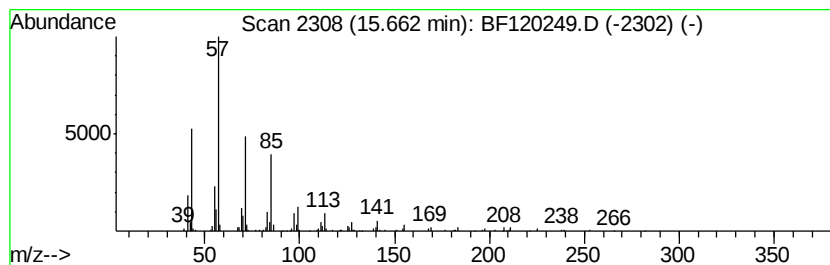
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	6.57 ng	338821	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Hexacosane	366	C26H54	000630-01-3	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Heptadecane	240	C17H36	000629-78-7	90
5			Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
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Misc :
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Instrument :
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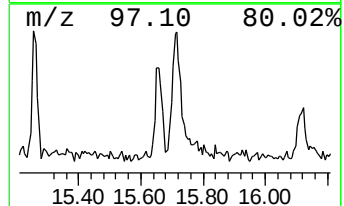
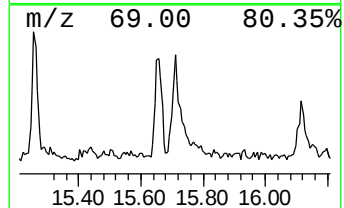
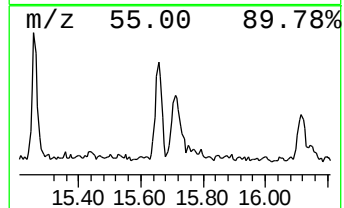
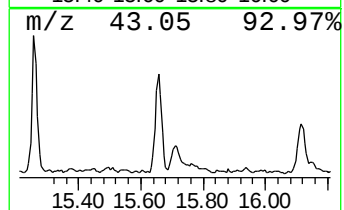
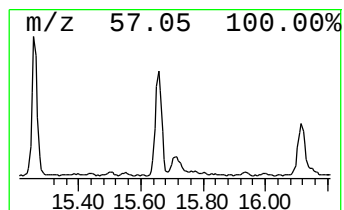
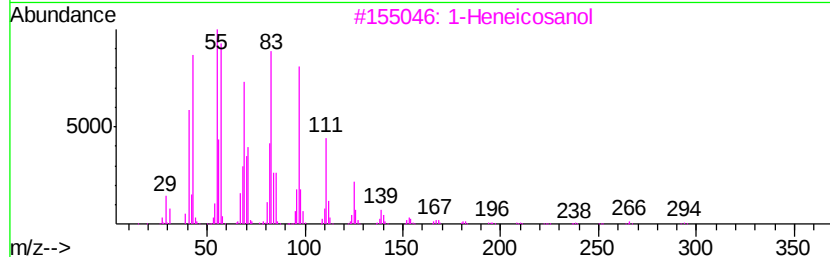
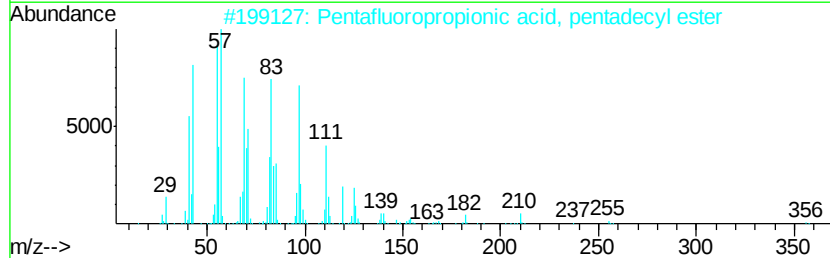
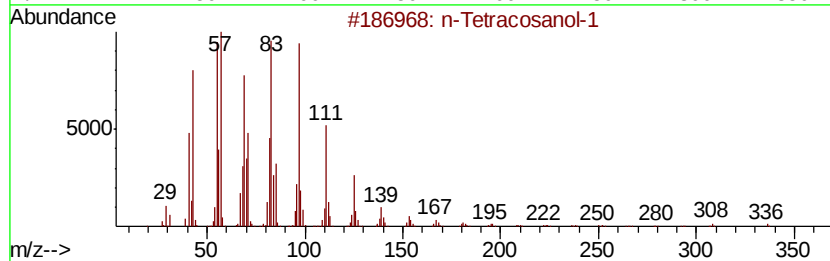
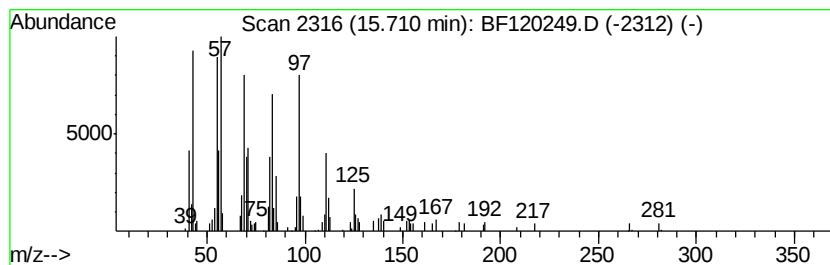
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Tetracosanol-1 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.71	3.23 ng	166607	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	94
4		Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
5		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
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Instrument :
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ClientSampleId :
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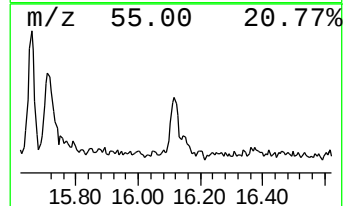
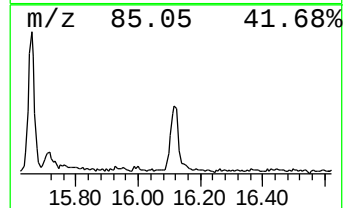
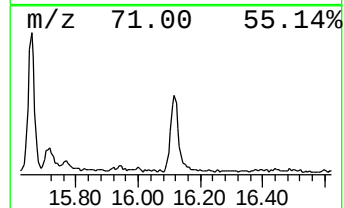
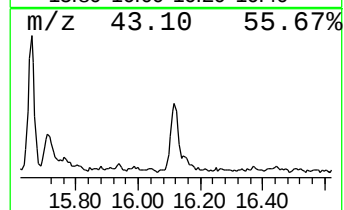
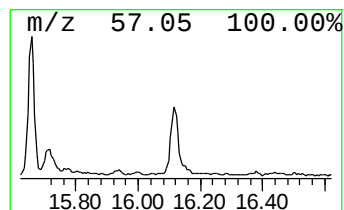
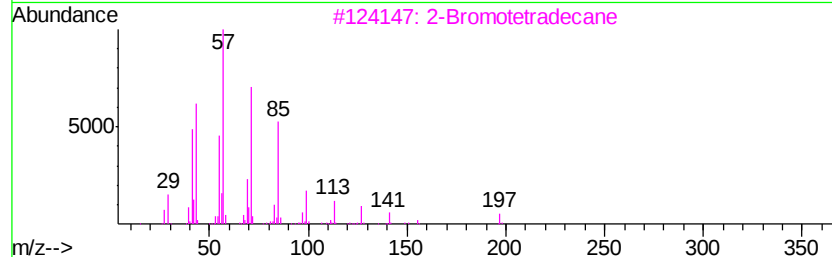
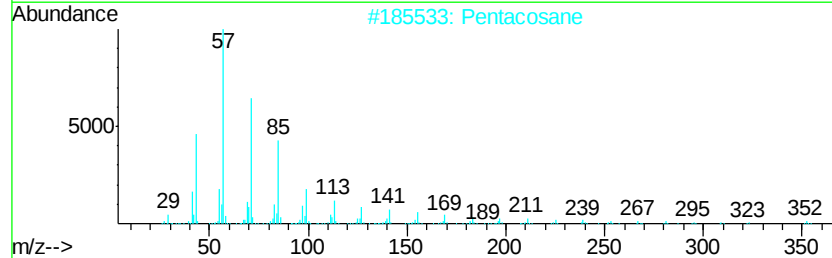
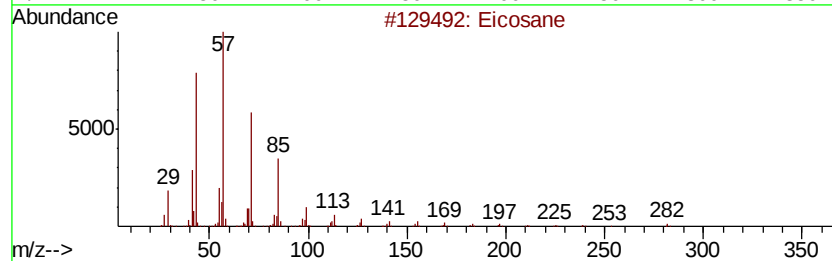
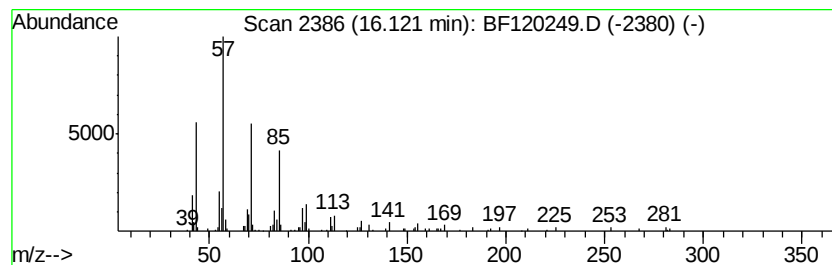
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	4.26 ng	219568	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	87
2		Pentacosane	352	C25H52	000629-99-2	87
3		2-Bromotetradecane	276	C14H29Br	074036-95-6	87
4		Hexatriacontane	507	C36H74	000630-06-8	86
5		Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120249.D
Acq On : 30 Apr 2020 19:45
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Sample : L2452-03
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Instrument :
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ClientSampleId :
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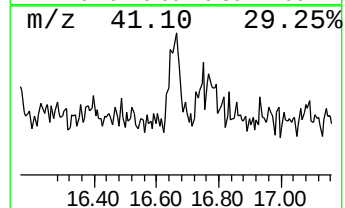
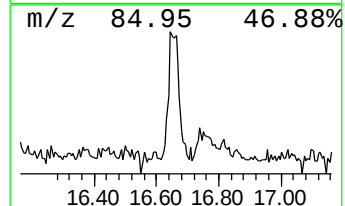
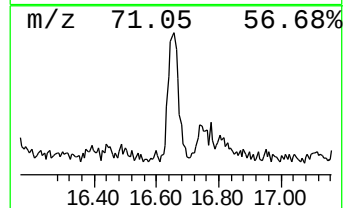
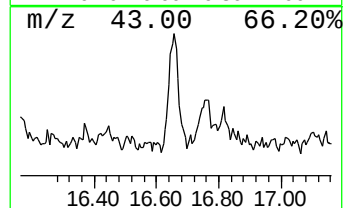
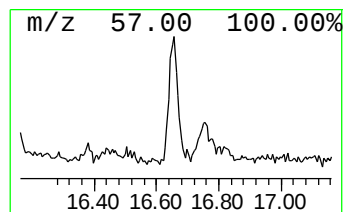
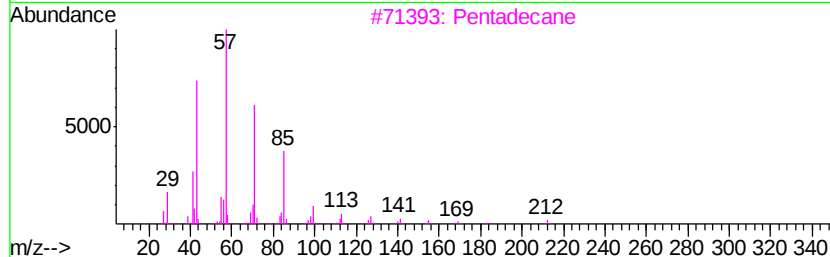
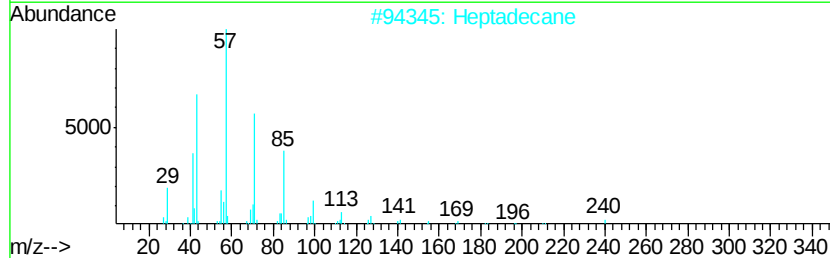
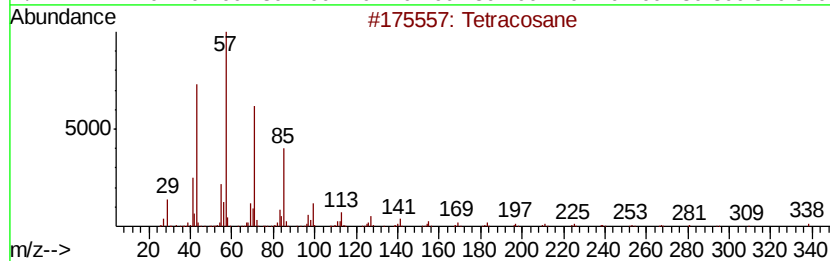
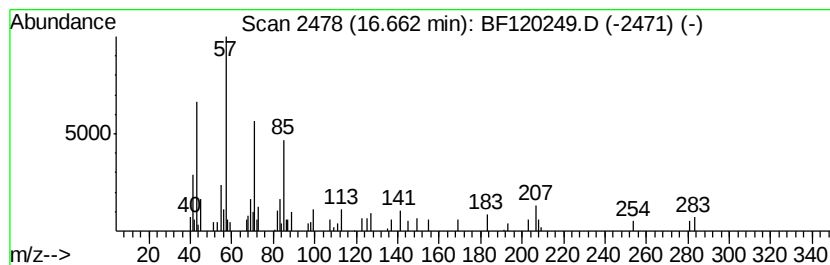
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.66	2.02 ng	104170	Perylene-d12	15.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	74
2			Heptadecane	240	C17H36	000629-78-7	72
3			Pentadecane	212	C15H32	000629-62-9	72
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	58
5			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF043020\
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 Sample : L2452-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 224

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.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
Mequinol	8.04	7.8 ng		344320	2	7.93	885236	20.0
n-Hexadecanoic acid	11.73	26.2 ng		1375320	4	11.17	1050820	20.0
Octadecanoic acid	12.51	14.8 ng		951692	5	13.82	1288680	20.0
Pentafluoropropio...	13.67	11.5 ng		741483	5	13.82	1288680	20.0
Hexacosane	14.00	2.1 ng		137651	5	13.82	1288680	20.0
Octacosane	14.30	4.8 ng		307972	5	13.82	1288680	20.0
Heptacosane	14.60	5.9 ng		303196	6	15.22	1031010	20.0
Hexadecane	14.91	7.5 ng		385902	6	15.22	1031010	20.0
Tetracosane	15.66	6.6 ng		338821	6	15.22	1031010	20.0
n-Tetracosanol-1	15.71	3.2 ng		166607	6	15.22	1031010	20.0
Eicosane	16.12	4.3 ng		219568	6	15.22	1031010	20.0
Heptadecane	16.66	2.0 ng		104170	6	15.22	1031010	20.0