

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
 Data File : BF124261.D
 Acq On : 11 Jun 2021 15:17
 Operator : JU/CG
 Sample : M2631-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 0225-FIELD-BLANK-1

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-BF060321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124261.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.045	500	503	510	rVB	21630	25454	1.42%	0.260%
2	5.469	571	575	584	rVB	527160	496396	27.77%	5.076%
3	6.475	743	746	757	rVB	301948	296838	16.61%	3.036%
4	6.669	776	779	783	rBV2	23164	22492	1.26%	0.230%
5	6.839	805	808	811	rBV	320628	242467	13.56%	2.480%
6	7.404	899	904	907	rBV	845729	845196	47.28%	8.643%
7	8.122	1022	1026	1030	rBV	416430	330353	18.48%	3.378%
8	9.204	1205	1210	1213	rBV	1977347	1669248	93.38%	17.071%
9	9.592	1273	1276	1279	rBV	171672	122753	6.87%	1.255%
10	9.880	1321	1325	1328	rBV	531793	412179	23.06%	4.215%
11	10.674	1455	1460	1463	rBV	1416966	1210663	67.73%	12.381%
12	11.368	1574	1578	1581	rBV	515197	428360	23.96%	4.381%
13	11.892	1664	1667	1671	rBV	76084	60555	3.39%	0.619%
14	12.086	1697	1700	1705	rVB	45342	42325	2.37%	0.433%
15	12.774	1813	1817	1819	rBV	221423	193467	10.82%	1.979%
16	12.821	1823	1825	1827	rBV	22867	21159	1.18%	0.216%
17	12.845	1827	1829	1833	rVB2	58825	47327	2.65%	0.484%
18	12.957	1843	1848	1851	rBV	1916237	1787610	100.00%	18.281%
19	13.180	1881	1886	1889	rBV	35511	33074	1.85%	0.338%
20	13.480	1933	1937	1941	rBV	186185	176793	9.89%	1.808%
21	13.521	1941	1944	1947	rVV	49984	50646	2.83%	0.518%
22	13.545	1947	1948	1955	rVB2	31613	28430	1.59%	0.291%
23	13.733	1978	1980	1985	rBV5	13939	23517	1.32%	0.240%
24	13.857	1997	2001	2011	rVB4	62214	117397	6.57%	1.201%
25	14.009	2024	2027	2031	rVB	323474	285146	15.95%	2.916%
26	14.168	2049	2054	2057	rBV	57503	53005	2.97%	0.542%
27	14.474	2102	2106	2110	rBV	59616	57886	3.24%	0.592%
28	14.780	2156	2158	2162	rVB	51834	54647	3.06%	0.559%
29	14.945	2178	2186	2188	rBV8	10731	22079	1.24%	0.226%
30	15.121	2213	2216	2221	rVB	52795	60245	3.37%	0.616%
31	15.492	2274	2279	2285	rVB	273671	363430	20.33%	3.717%
32	15.939	2351	2355	2363	rVB2	44028	71894	4.02%	0.735%
33	16.450	2437	2442	2447	rVB4	30474	54892	3.07%	0.561%
34	17.045	2537	2543	2548	rBV4	24311	49783	2.78%	0.509%
35	17.574	2627	2633	2634	rBV6	12457	20701	1.16%	0.212%

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ClientSampleId :
0225-FIELD-BLANK-1

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-BF060321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

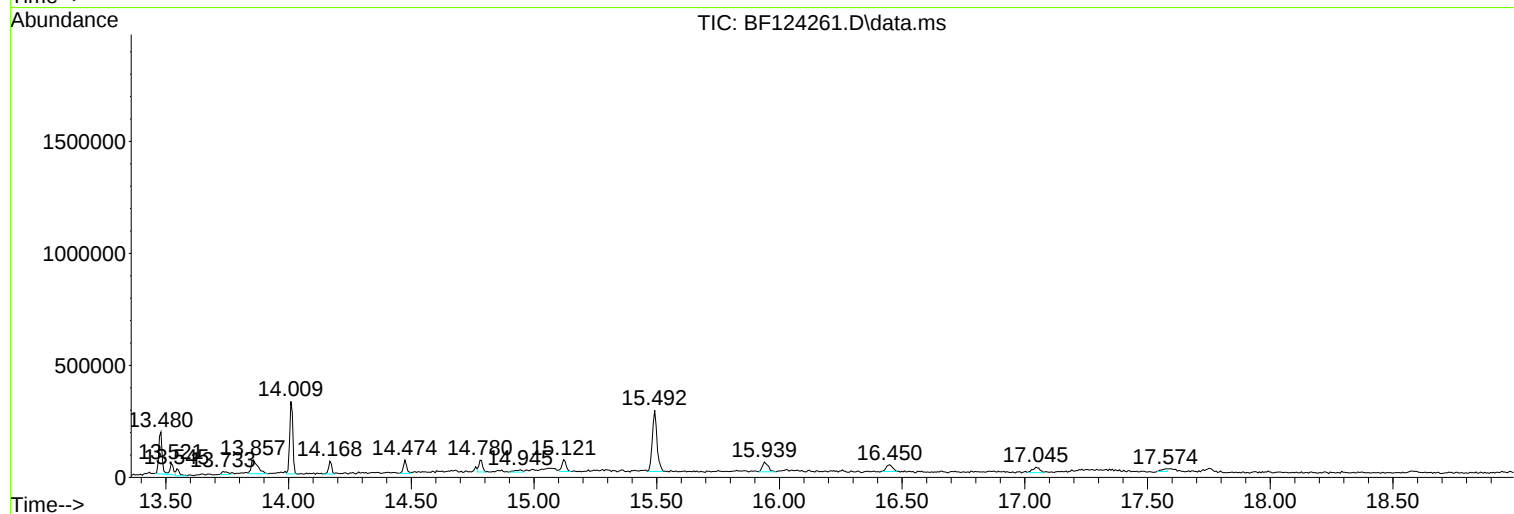
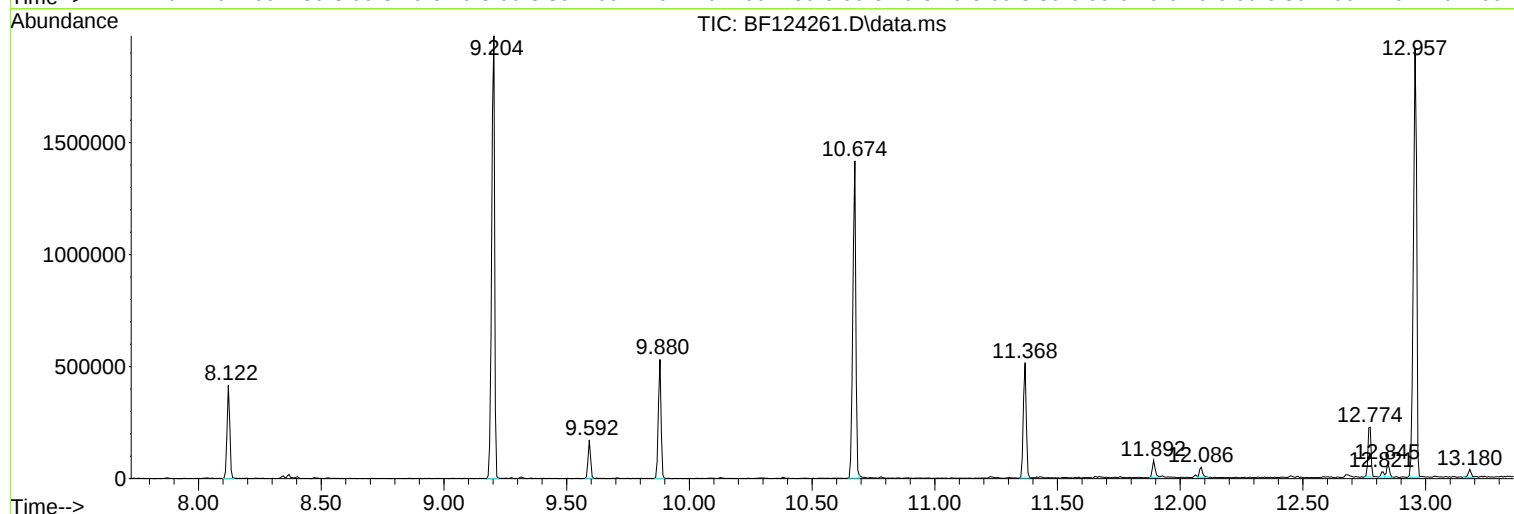
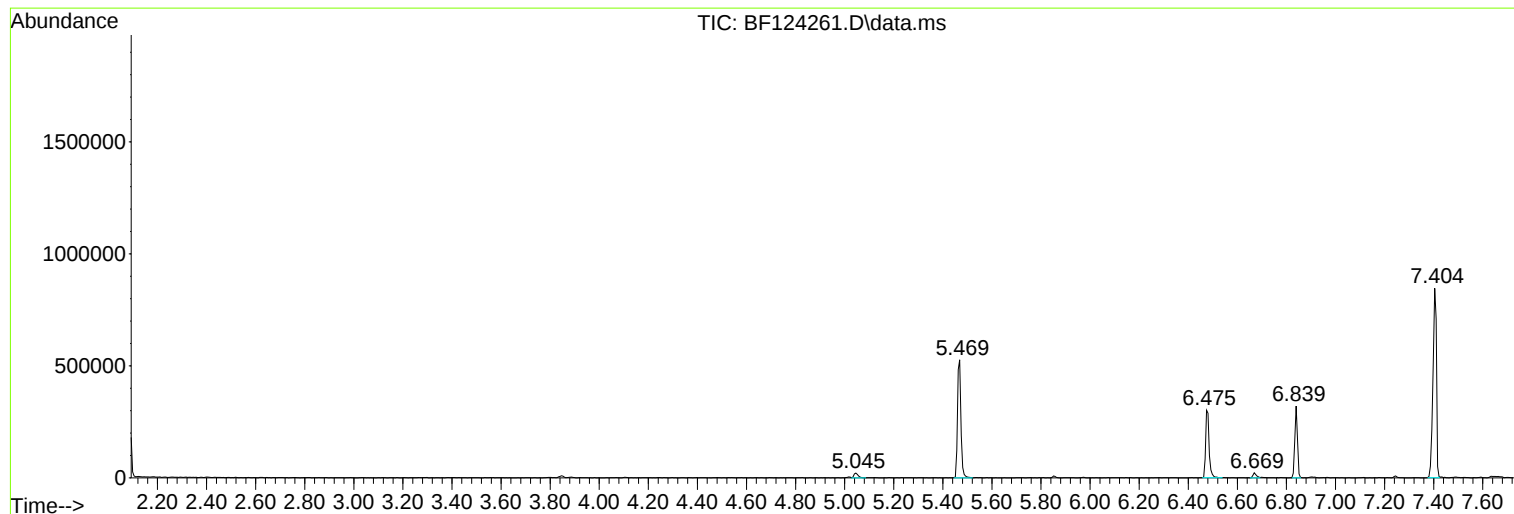
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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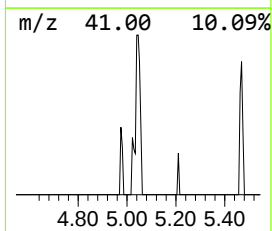
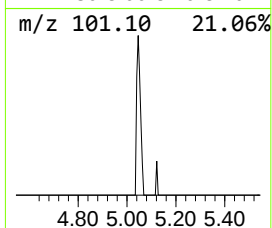
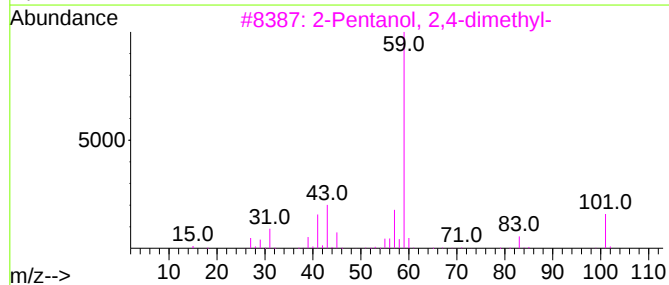
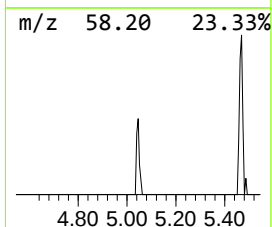
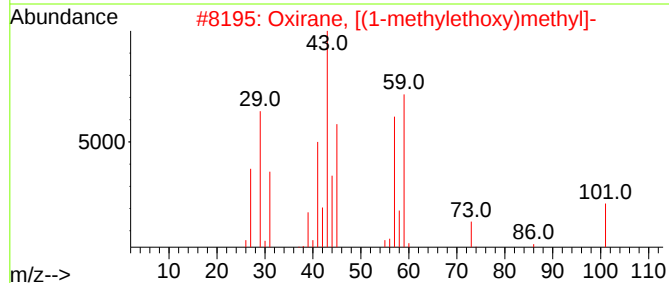
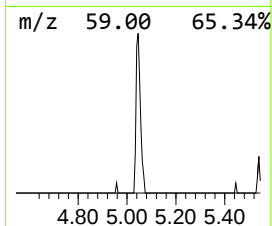
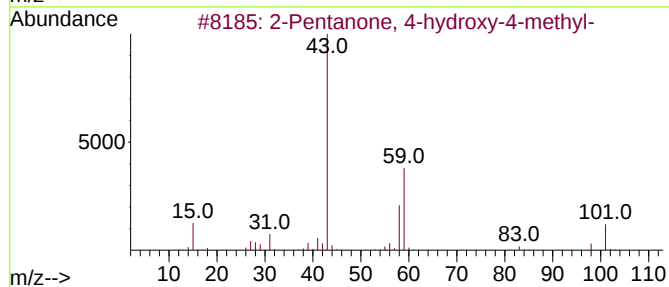
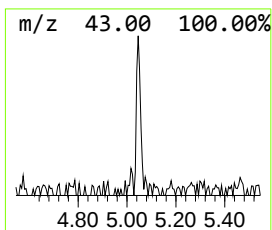
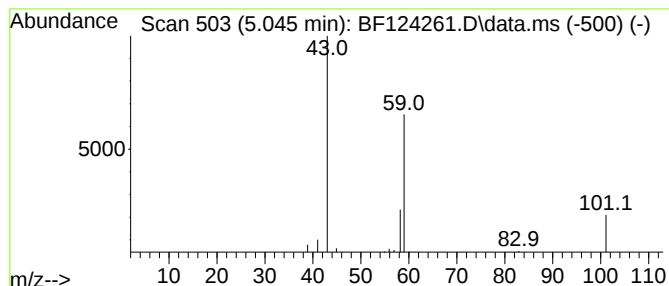
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.045	2.10 ng	25454	1,4-Dichlorobenzene-d4	6.839

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
2	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	43		
3	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	42		
4	3-Hexanol	102	C6H14O	000623-37-0	9		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9		



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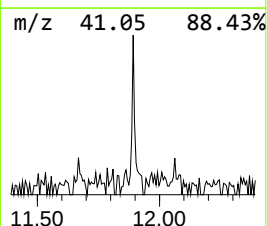
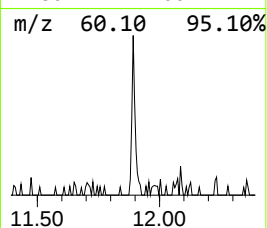
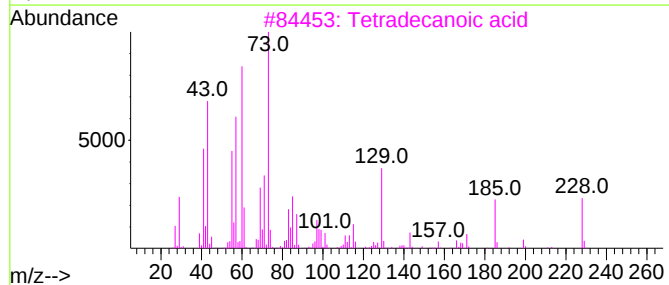
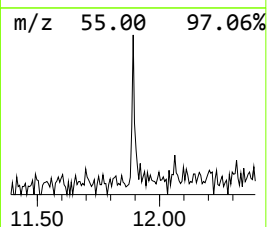
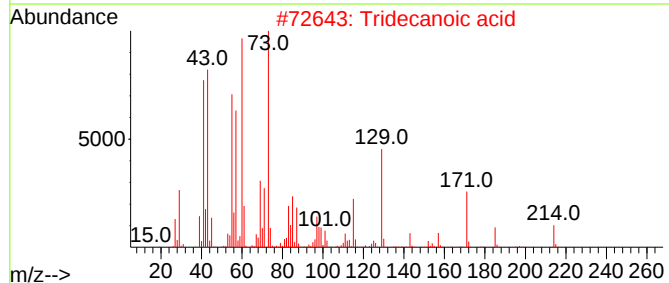
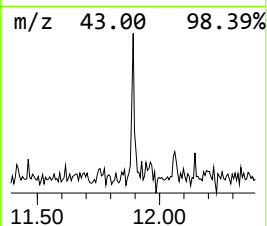
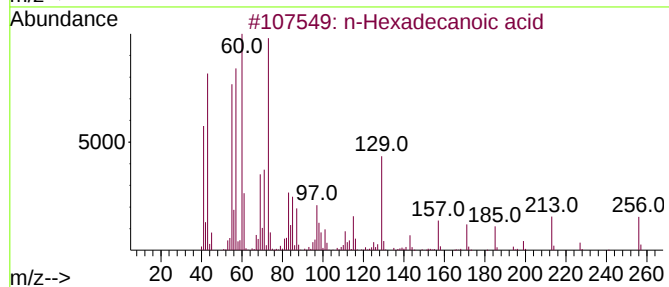
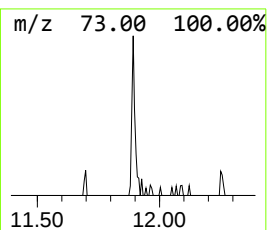
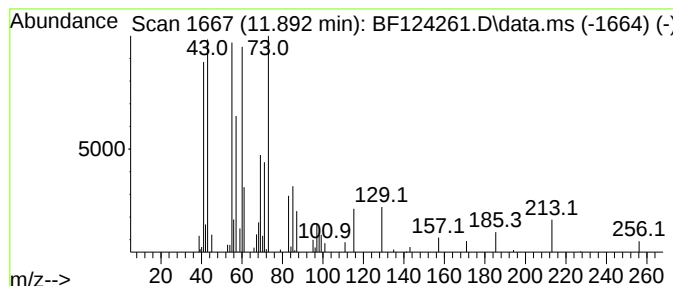
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.892	2.83 ng	60555	Phenanthrene-d10	11.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
2			Tridecanoic acid	214	C13H26O2	000638-53-9	53
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	53
4			n-Decanoic acid	172	C10H20O2	000334-48-5	53
5			Nonanoic acid	158	C9H18O2	000112-05-0	38



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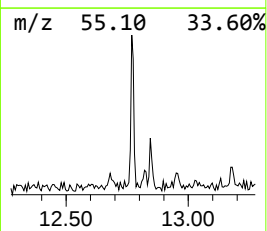
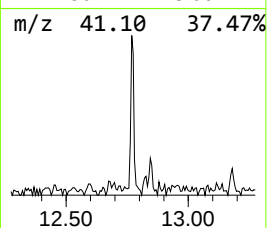
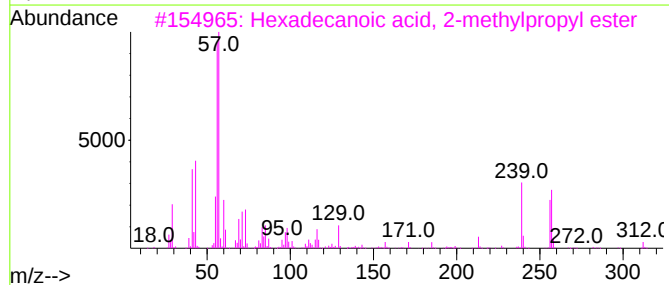
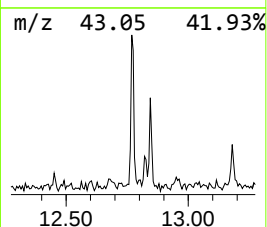
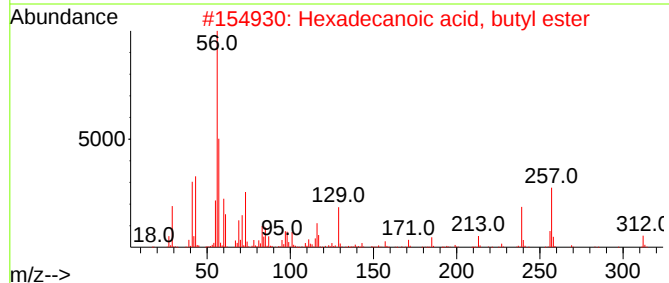
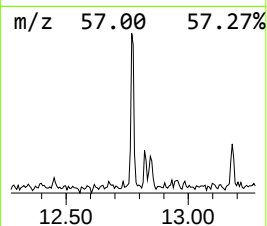
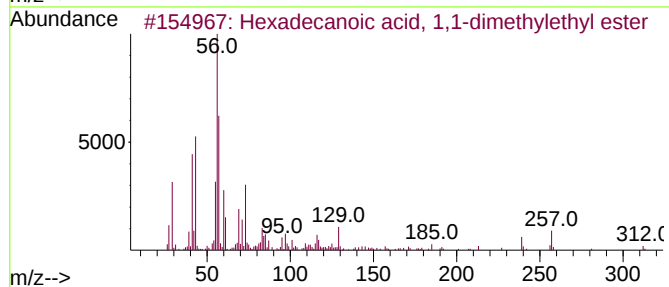
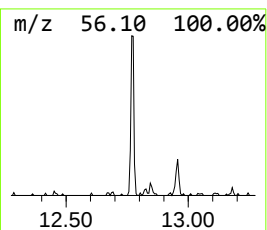
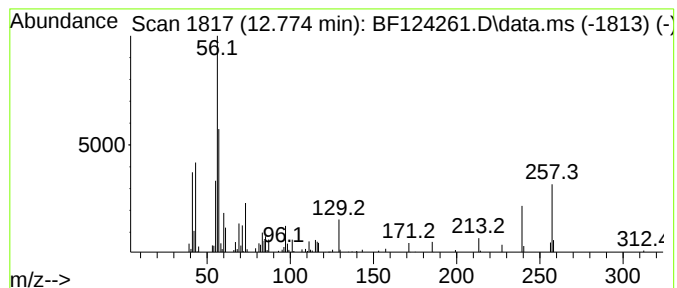
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexadecanoic acid, 1,1-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.774	13.57 ng	193467	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	93
2			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
3			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	74
4			Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	38
5			Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	35



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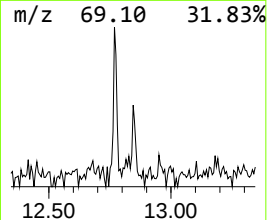
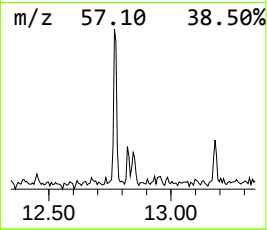
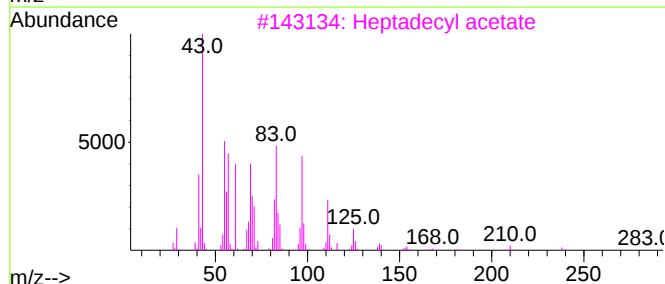
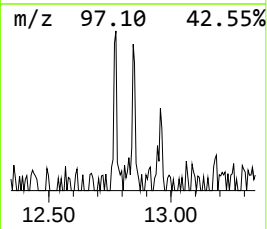
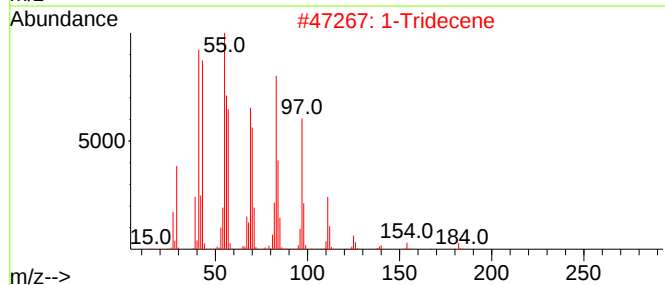
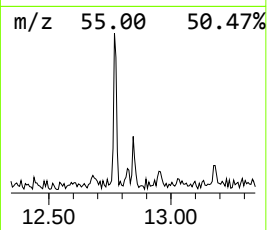
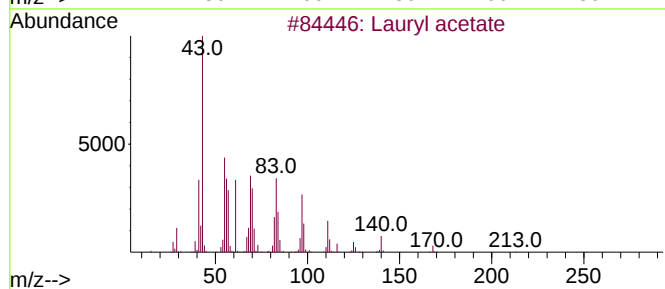
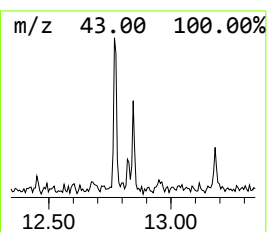
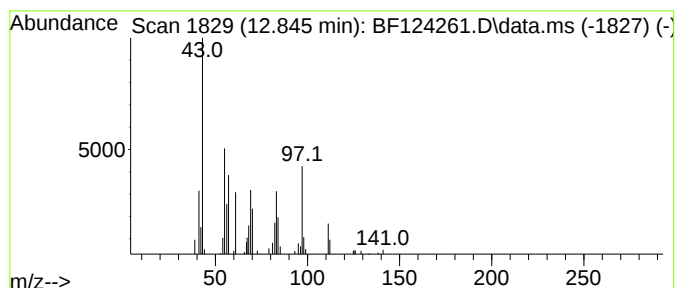
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Lauryl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.845	3.32 ng	47327	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lauryl acetate	228	C14H28O2	000112-66-3	80
2			1-Tridecene	182	C13H26	002437-56-1	58
3			Heptadecyl acetate	298	C19H38O2	000822-20-8	58
4			1-Dodecanol	186	C12H26O	000112-53-8	53
5			1-Hexadecanol	242	C16H34O	036653-82-4	52



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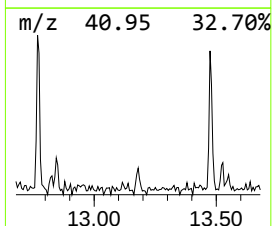
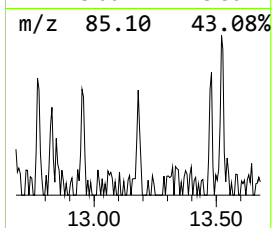
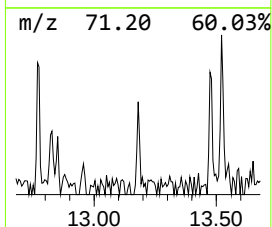
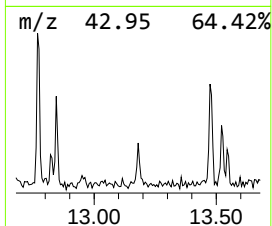
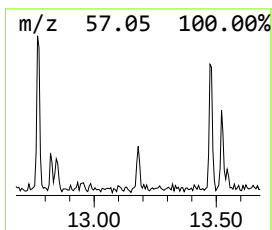
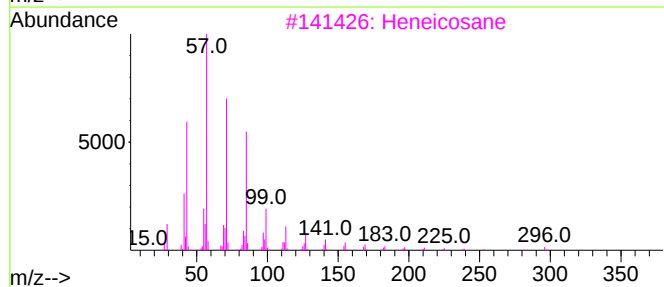
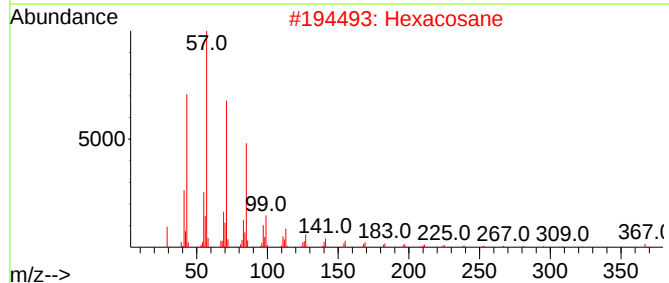
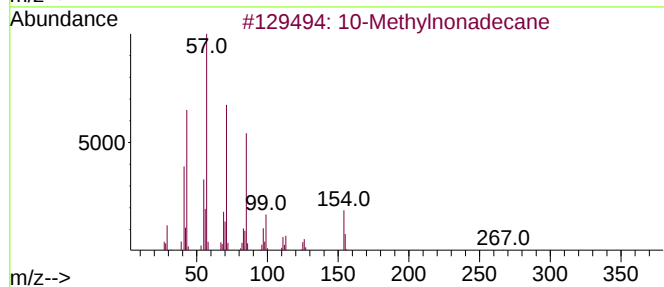
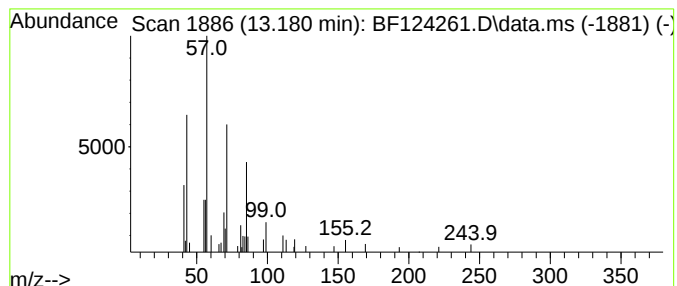
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 10-Methylnonadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.180	2.32 ng	33074	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			10-Methylnonadecane	282	C20H42	056862-62-5	72
2			Hexacosane	366	C26H54	000630-01-3	59
3			Heneicosane	296	C21H44	000629-94-7	59
4			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	59
5			Heptacosane	380	C27H56	000593-49-7	59



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Acq On : 11 Jun 2021 15:17
Operator : JU/CG
Sample : M2631-01
Misc :
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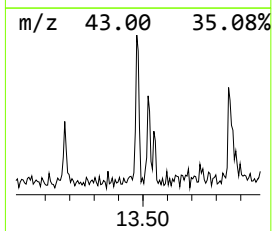
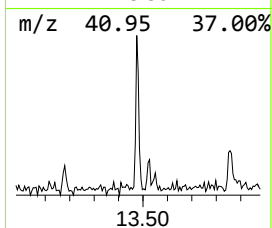
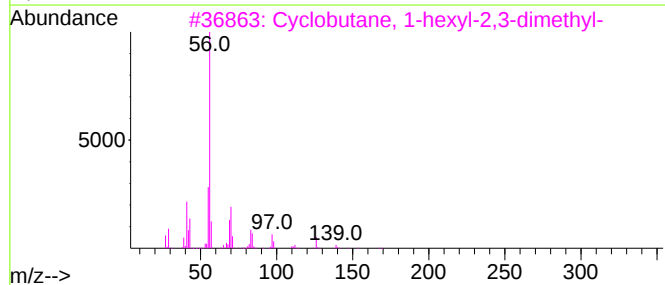
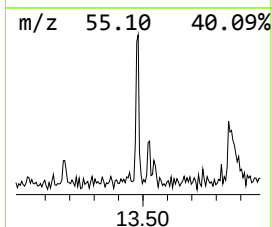
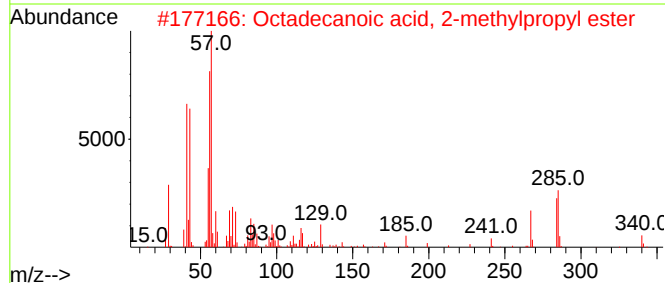
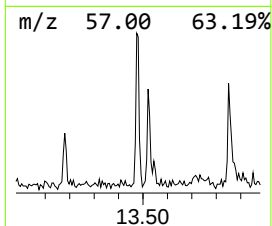
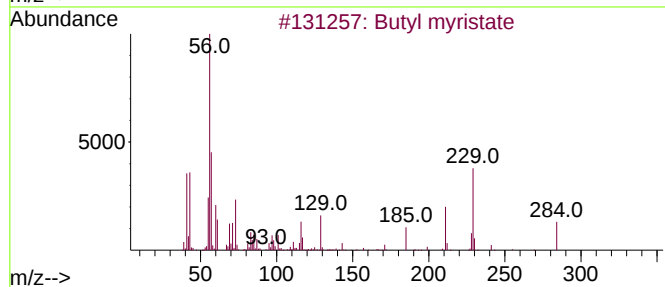
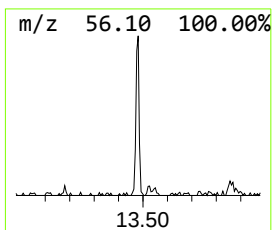
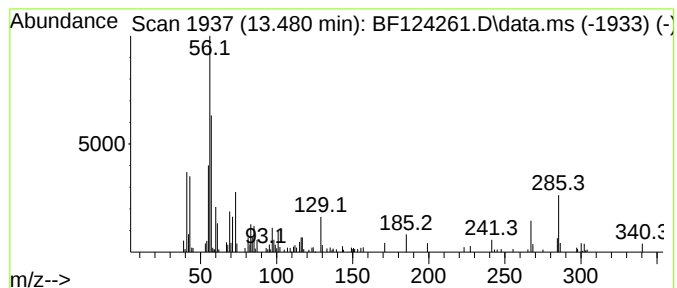
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Butyl myristate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.480	12.40 ng	176793	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butyl myristate	284	C18H36O2	000110-36-1	58
2			Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	41
3			Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	35
4			1-Heptene, 2-methyl-	112	C8H16	015870-10-7	35
5			1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
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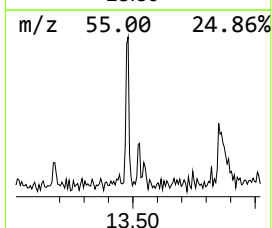
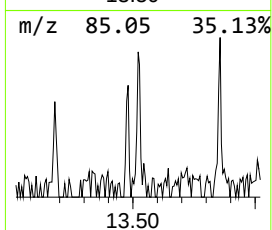
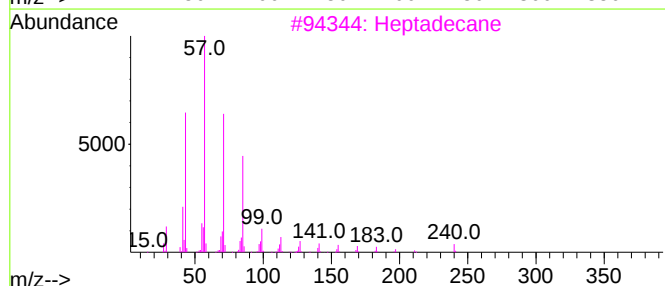
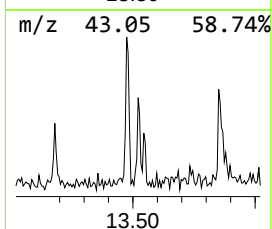
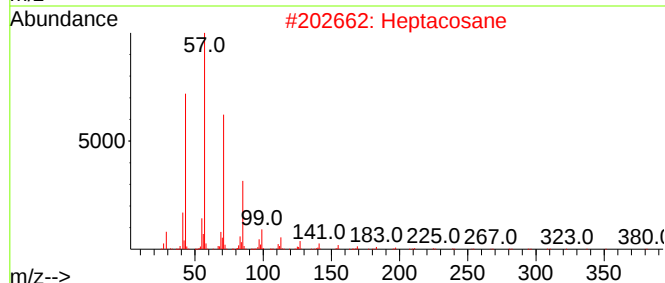
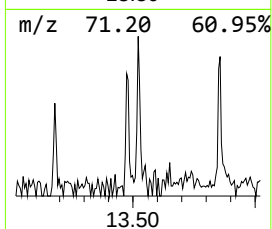
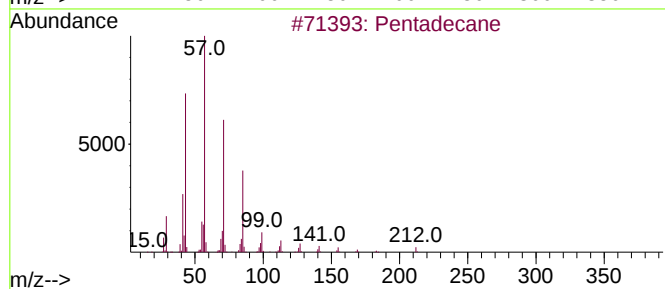
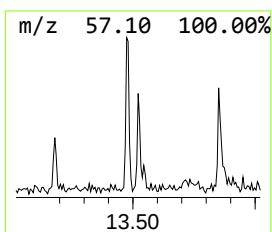
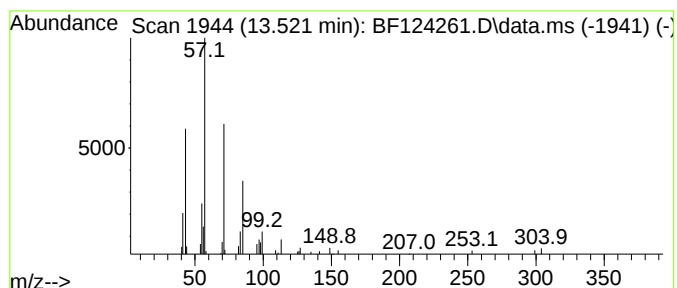
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.521	3.55 ng	50646	Chrysene-d12	14.009

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane	212	C15H32	000629-62-9	72
2		Heptacosane	380	C27H56	000593-49-7	72
3		Heptadecane	240	C17H36	000629-78-7	72
4		Octacosane	394	C28H58	000630-02-4	72
5		Tetracosane	338	C24H50	000646-31-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
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Acq On : 11 Jun 2021 15:17
Operator : JU/CG
Sample : M2631-01
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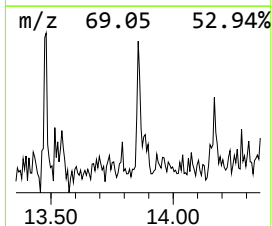
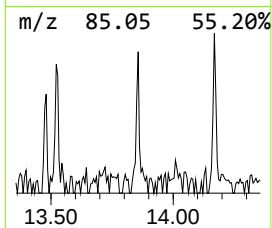
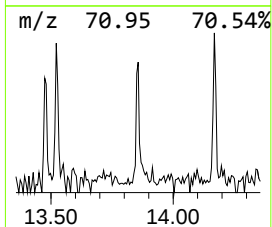
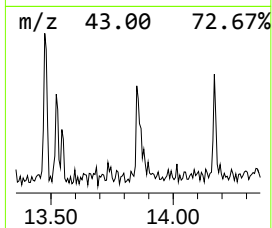
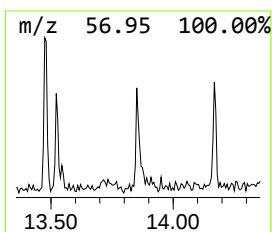
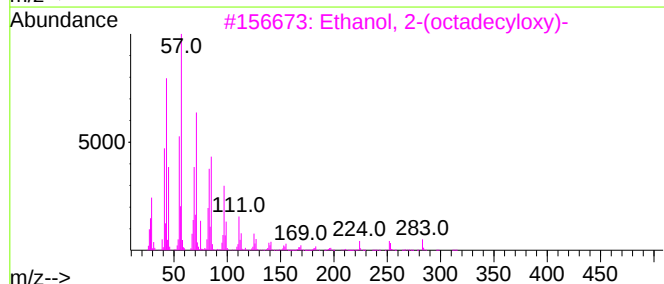
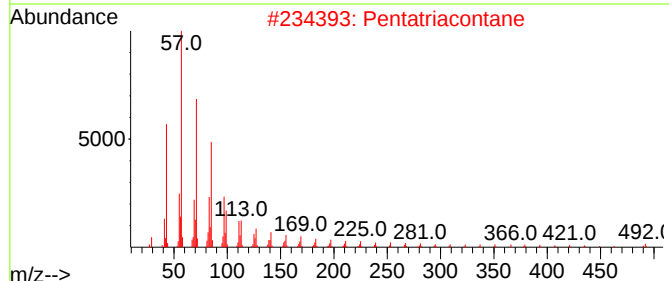
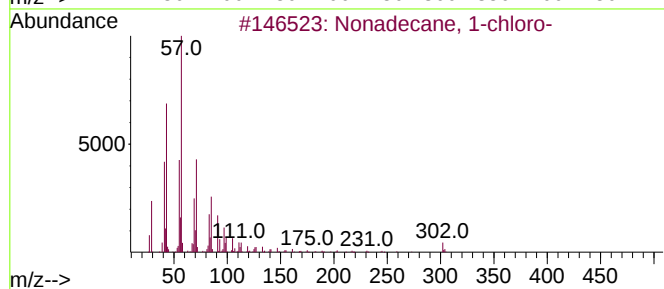
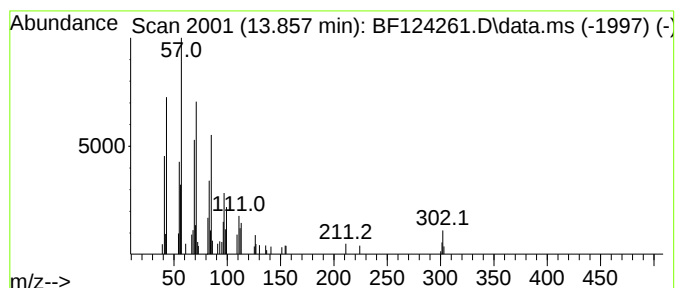
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane, 1-chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.857	8.23 ng	117397	Chrysene-d12	14.009	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	93
2	Pentatriacontane	493	C35H72	000630-07-9	72
3	Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	64
4	Heptadecane	240	C17H36	000629-78-7	58
5	Hexadecane	226	C16H34	000544-76-3	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
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Acq On : 11 Jun 2021 15:17
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Instrument :
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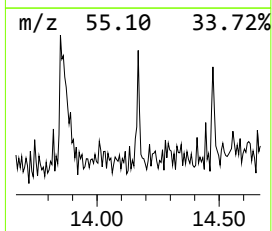
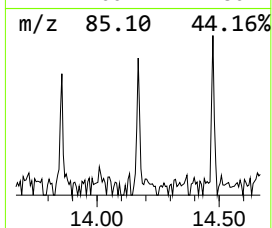
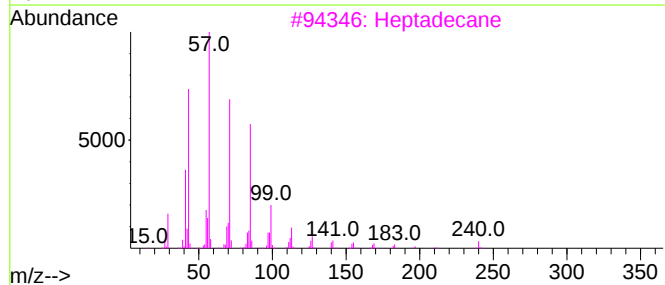
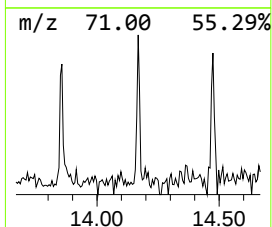
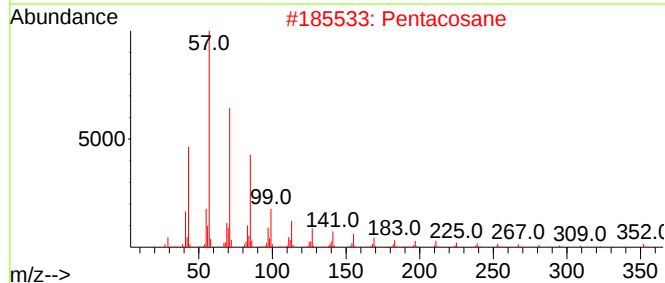
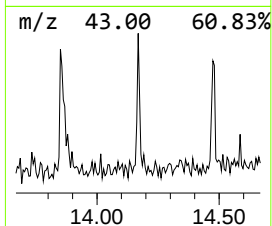
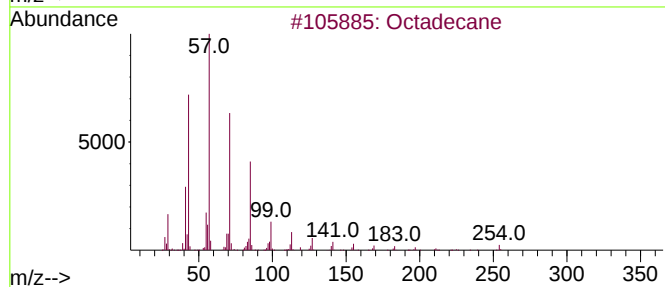
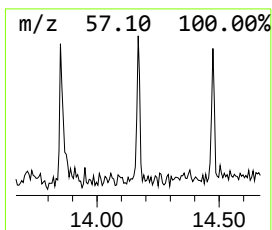
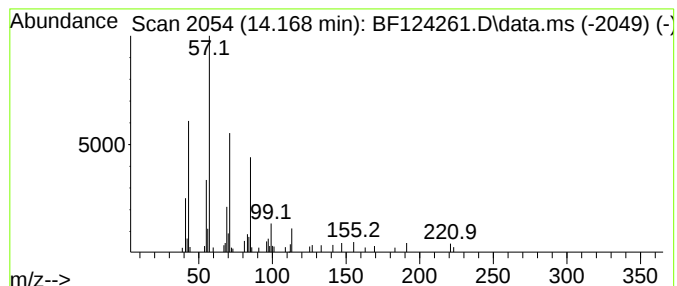
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.168	3.72 ng	53005	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	86
2			Pentacosane	352	C25H52	000629-99-2	80
3			Heptadecane	240	C17H36	000629-78-7	80
4			Hexadecane	226	C16H34	000544-76-3	72
5			Hexacosane	366	C26H54	000630-01-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
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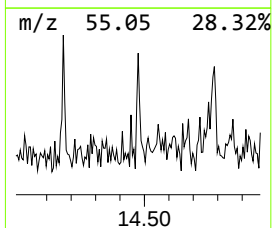
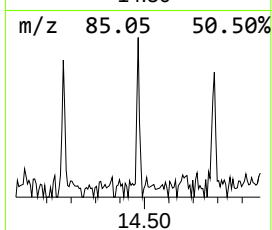
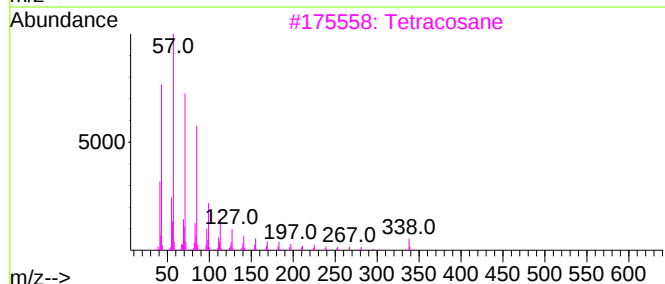
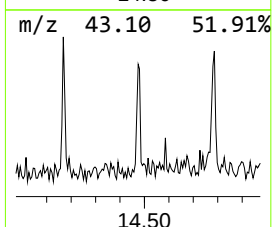
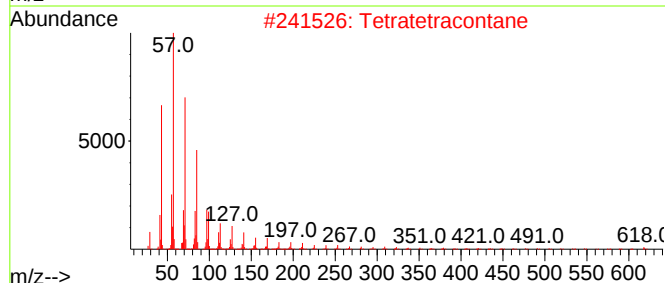
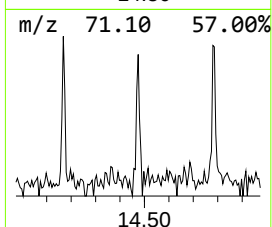
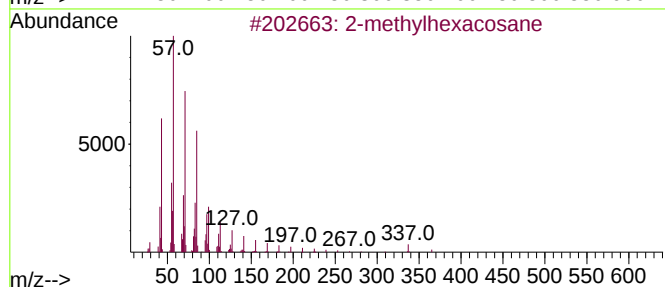
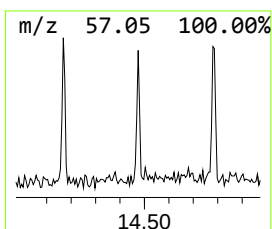
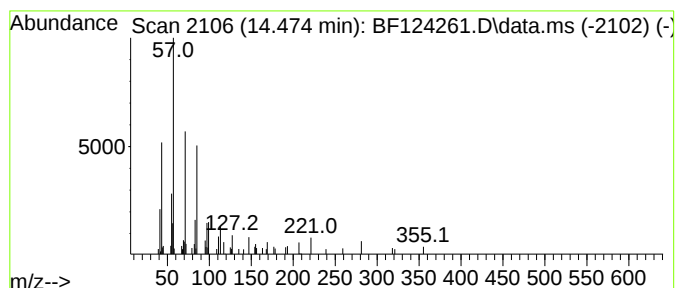
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 2-methylhexacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.474	4.06 ng	57886	Chrysene-d12	14.009

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-methylhexacosane	380	C27H56	1000376-72-7	86
2		Tetratetracontane	619	C44H90	007098-22-8	86
3		Tetracosane	338	C24H50	000646-31-1	86
4		Hexacosane	366	C26H54	000630-01-3	78
5		2-Bromotetradecane	276	C14H29Br	074036-95-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
Acq On : 11 Jun 2021 15:17
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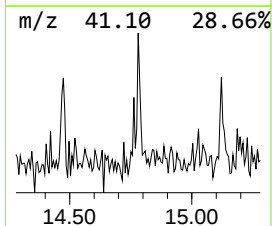
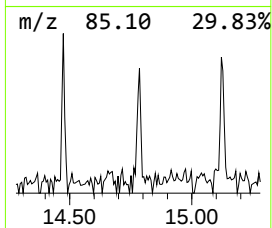
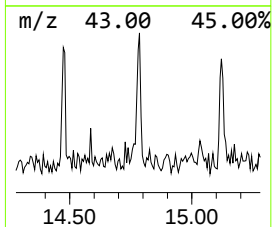
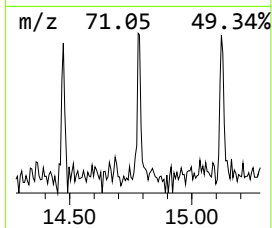
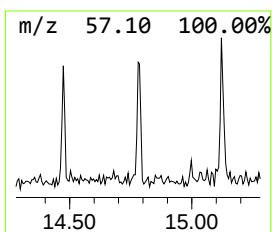
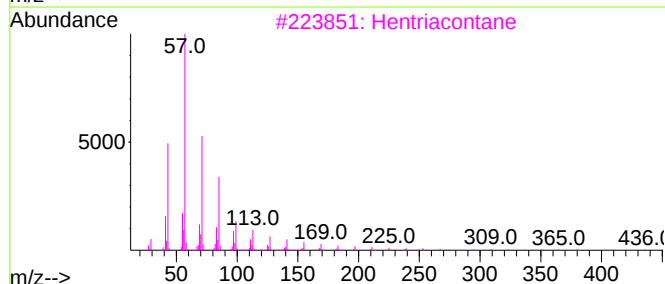
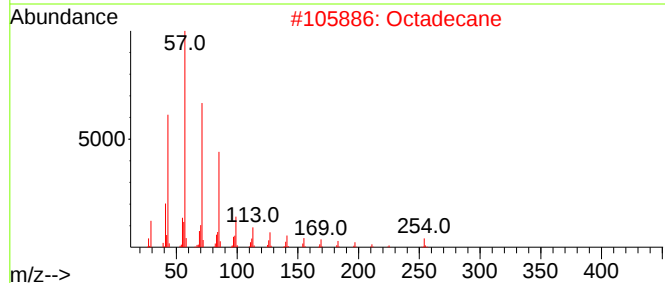
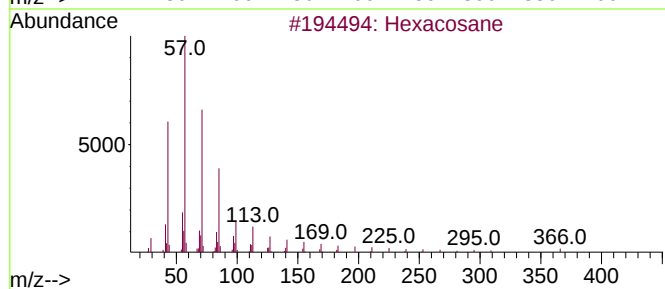
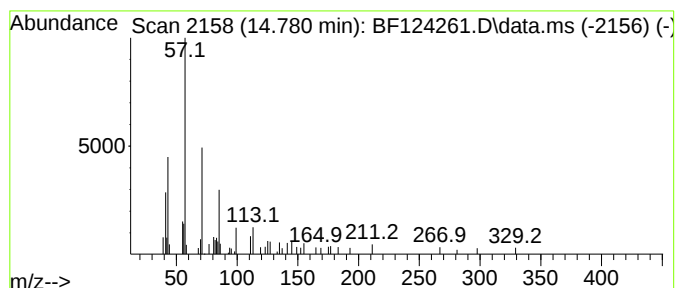
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	3.01 ng	54647	Perylene-d12	15.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	86
2			Octadecane	254	C18H38	000593-45-3	86
3			Hentriacontane	437	C31H64	000630-04-6	86
4			Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	80
5			Pentacosane	352	C25H52	000629-99-2	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
Acq On : 11 Jun 2021 15:17
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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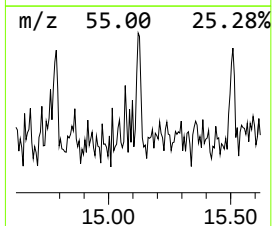
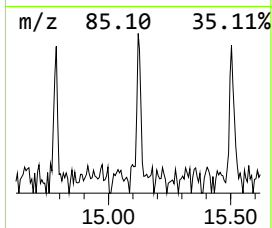
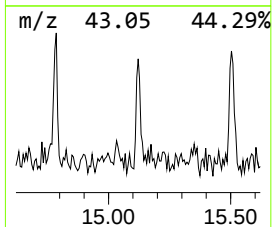
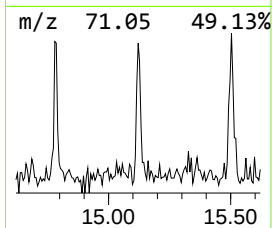
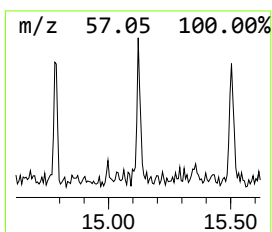
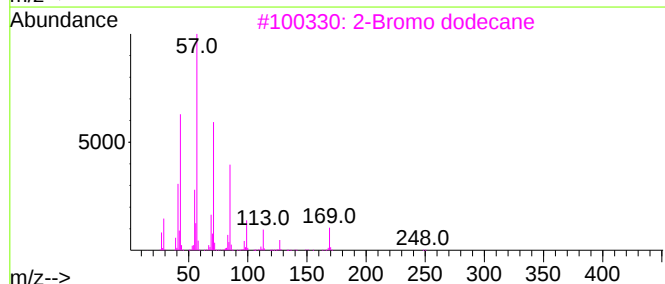
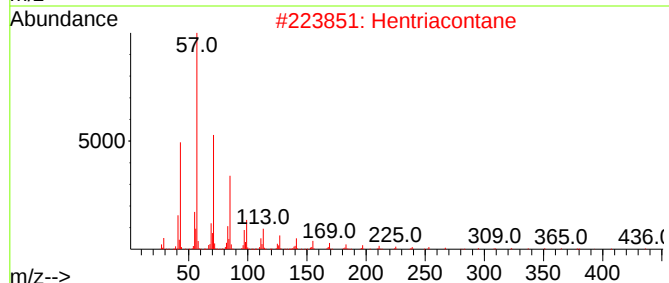
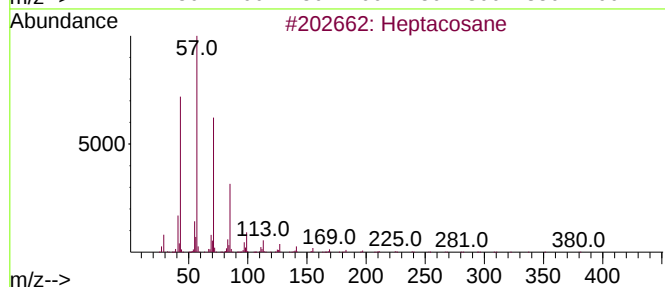
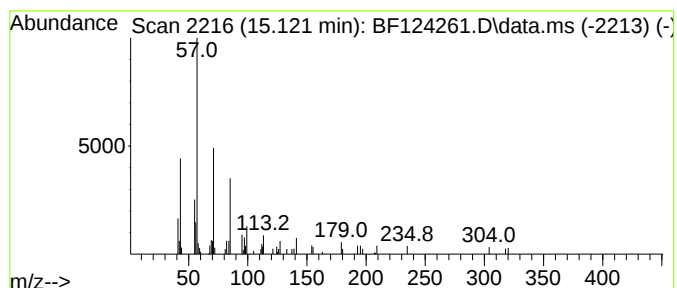
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.121	3.32 ng	60245	Perylene-d12	15.492

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	64
2		Hentriacontane	437	C31H64	000630-04-6	59
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	59
4		Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	59
5		Tridecane	184	C13H28	000629-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
Acq On : 11 Jun 2021 15:17
Operator : JU/CG
Sample : M2631-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

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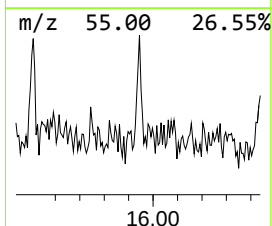
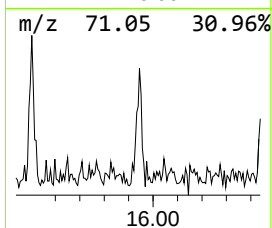
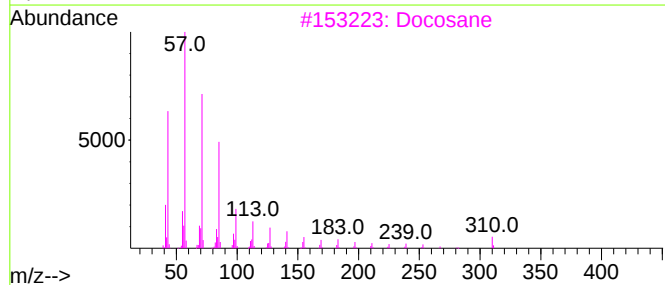
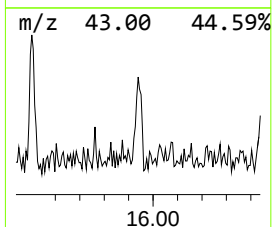
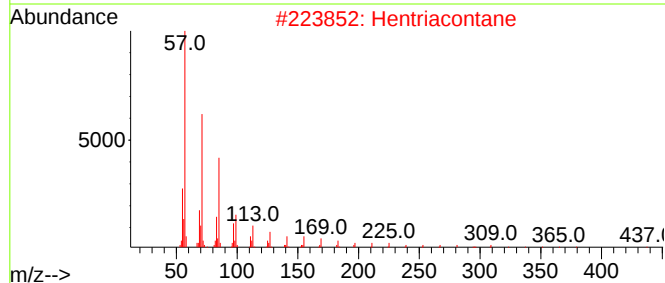
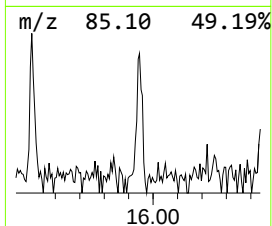
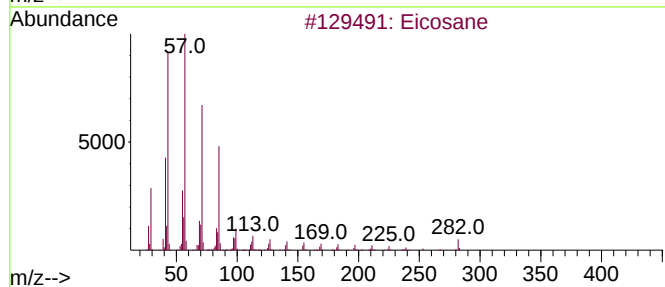
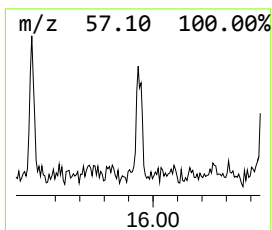
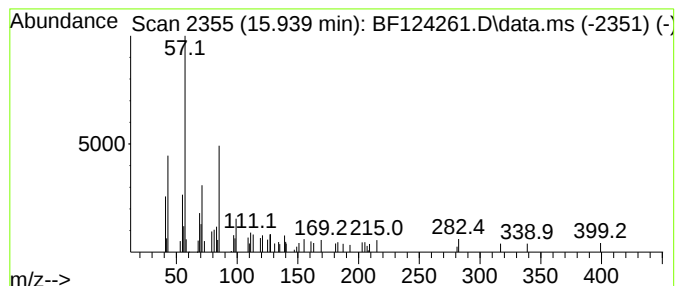
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.939	3.96 ng	71894	Perylene-d12	15.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	60
2			Hentriacontane	437	C31H64	000630-04-6	58
3			Docosane	310	C22H46	000629-97-0	49
4			Heptacosane	380	C27H56	000593-49-7	47
5			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
Acq On : 11 Jun 2021 15:17
Operator : JU/CG
Sample : M2631-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
0225-FIELD-BLANK-1

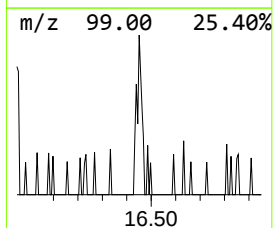
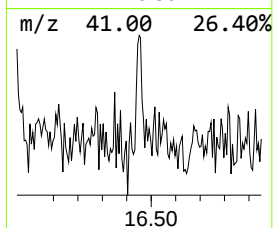
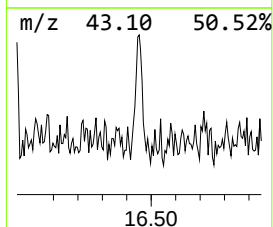
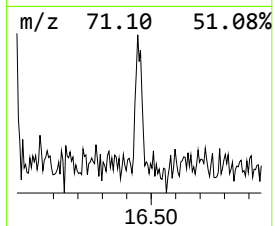
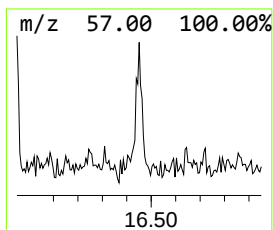
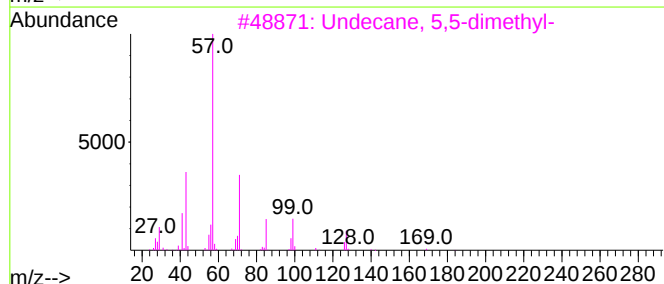
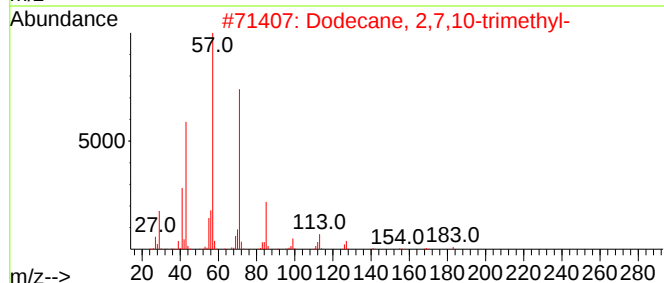
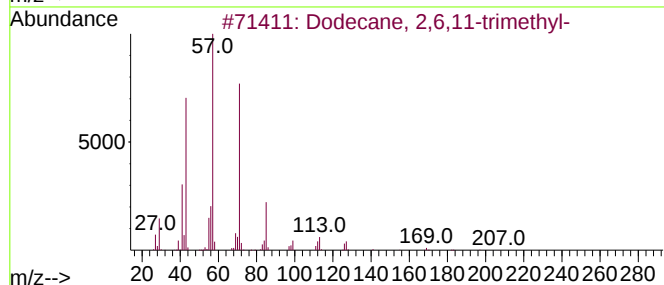
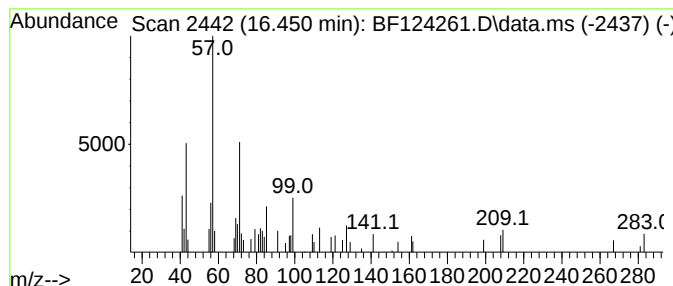
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Dodecane, 2,6,11-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.451	3.02 ng	54892	Perylene-d12	15.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	59
2			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	59
3			Undecane, 5,5-dimethyl-	184	C13H28	017312-73-1	59
4			Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	50
5			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
Acq On : 11 Jun 2021 15:17
Operator : JU/CG
Sample : M2631-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

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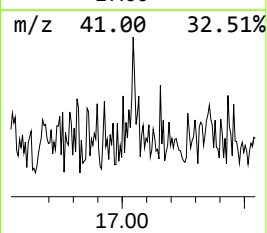
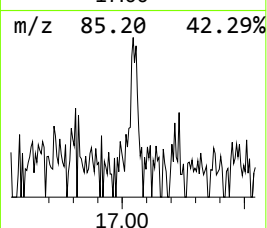
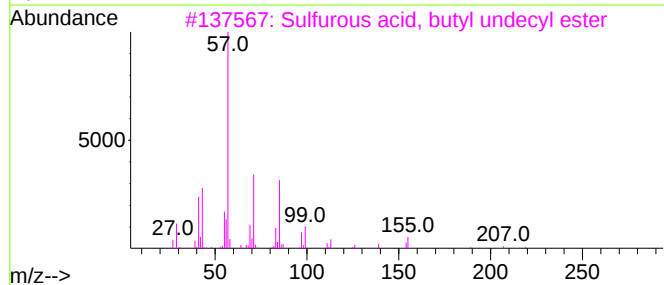
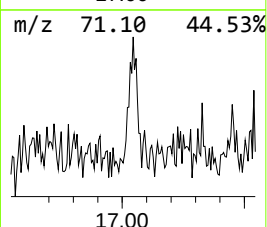
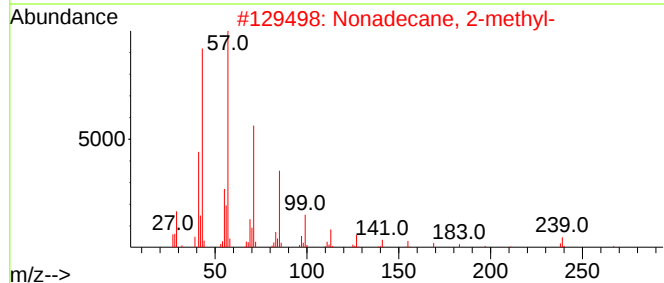
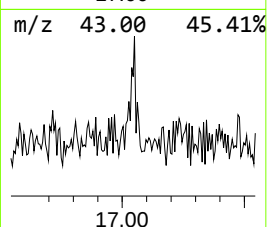
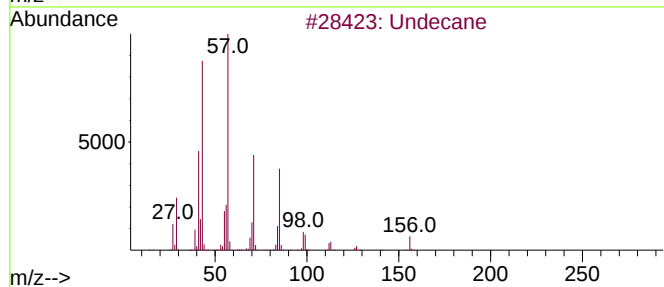
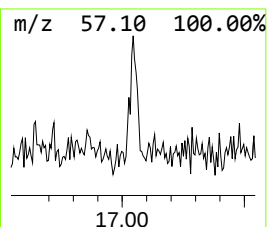
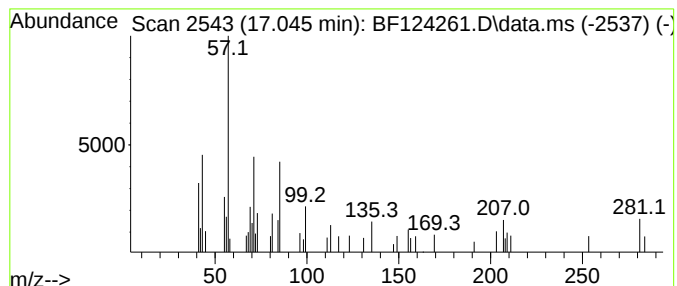
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Undecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.045	2.74 ng	49783	Perylene-d12	15.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	55
2	Nonadecane, 2-methyl-			282	C20H42	001560-86-7	50
3	Sulfurous acid, butyl undecyl ester			292	C15H32O3S	1000309-17-8	47
4	Heptacosane			380	C27H56	000593-49-7	47
5	2-Bromo dodecane			248	C12H25Br	013187-99-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124261.D
Acq On : 11 Jun 2021 15:17
Operator : JU/CG
Sample : M2631-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.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-...	5.045	2.1	ng	25454	1	6.839	242467	20.0
n-Hexadecanoic ...	11.892	2.8	ng	60555	4	11.368	428360	20.0
Hexadecanoic ac...	12.774	13.6	ng	193467	5	14.009	285146	20.0
Lauryl acetate	12.845	3.3	ng	47327	5	14.009	285146	20.0
10-Methylnonade...	13.180	2.3	ng	33074	5	14.009	285146	20.0
Butyl myristate	13.480	12.4	ng	176793	5	14.009	285146	20.0
Pentadecane	13.521	3.5	ng	50646	5	14.009	285146	20.0
Nonadecane, 1-c...	13.857	8.2	ng	117397	5	14.009	285146	20.0
Octadecane	14.168	3.7	ng	53005	5	14.009	285146	20.0
2-methylhexacosane	14.474	4.1	ng	57886	5	14.009	285146	20.0
Hexacosane	14.780	3.0	ng	54647	6	15.492	363430	20.0
Heptacosane	15.121	3.3	ng	60245	6	15.492	363430	20.0
Eicosane	15.939	4.0	ng	71894	6	15.492	363430	20.0
Dodecane, 2,6,1...	16.451	3.0	ng	54892	6	15.492	363430	20.0
Undecane	17.045	2.7	ng	49783	6	15.492	363430	20.0