

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012720\
 Data File : BF118718.D
 Acq On : 27 Jan 2020 14:09
 Operator : CG/JU
 Sample : L1241-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-16

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-BF012020.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.205	10	20	27	rVB	1747158	2532479	22.67%	2.888%
2	5.110	511	514	522	rVB	91904	138510	1.24%	0.158%
3	5.493	573	579	582	rBV	7347558	8261568	73.97%	9.423%
4	6.492	742	749	752	rBV	6557550	8520298	76.29%	9.718%
5	6.857	804	811	815	rBV	2012931	1750218	15.67%	1.996%
6	7.022	834	839	842	rVB	6854114	7209293	64.55%	8.222%
7	7.422	900	907	910	rBV	5149776	5637258	50.47%	6.429%
8	8.139	1024	1029	1033	rBV	2551337	2232767	19.99%	2.547%
9	9.222	1206	1213	1216	rBV	9409954	10888565	97.49%	12.419%
10	9.333	1224	1232	1240	rVB2	469432	445865	3.99%	0.509%
11	9.604	1273	1278	1281	rBV	1419377	1137551	10.19%	1.297%
12	9.892	1322	1327	1331	rBV	2973772	2731430	24.46%	3.115%
13	10.475	1422	1426	1435	rBV	305183	259517	2.32%	0.296%
14	10.686	1455	1462	1465	rBV	6989087	7393565	66.20%	8.433%
15	11.375	1575	1579	1583	rBV	2975222	2741888	24.55%	3.127%
16	11.445	1586	1591	1594	rVB2	171712	149579	1.34%	0.171%
17	11.604	1614	1618	1624	rBV	154471	163329	1.46%	0.186%
18	11.904	1665	1669	1679	rBV	863022	957634	8.57%	1.092%
19	12.422	1753	1757	1763	rBV	1907790	1757466	15.74%	2.004%
20	12.621	1784	1791	1794	rBV	336947	414334	3.71%	0.473%
21	12.686	1799	1802	1807	rBV	200397	229896	2.06%	0.262%
22	12.969	1844	1850	1853	rBV	10067869	11168779	100.00%	12.738%
23	13.169	1881	1884	1887	rBV	299954	304420	2.73%	0.347%
24	13.810	1989	1993	1996	rVB	95850	119293	1.07%	0.136%
25	13.857	1996	2001	2012	rBV3	1005833	1358408	12.16%	1.549%
26	14.016	2023	2028	2031	rBV	2153643	2134544	19.11%	2.434%
27	14.180	2052	2056	2062	rBV	353598	343104	3.07%	0.391%
28	14.486	2104	2108	2117	rVB	550295	521014	4.66%	0.594%
29	14.792	2156	2160	2166	rVB	634388	633092	5.67%	0.722%
30	14.868	2169	2173	2178	rVB	1117265	1094367	9.80%	1.248%
31	15.133	2213	2218	2225	rBV	574212	700622	6.27%	0.799%
32	15.474	2271	2276	2279	rBV	1410681	1786708	16.00%	2.038%
33	15.510	2279	2282	2297	rVB	579084	810004	7.25%	0.924%
34	15.945	2351	2356	2362	rBV	375983	602651	5.40%	0.687%

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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

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35	16.457	2436	2443	2452	rbV2	197960	415923	3.72%	0.474%
36	17.051	2539	2544	2550	rvB2	65981	133078	1.19%	0.152%

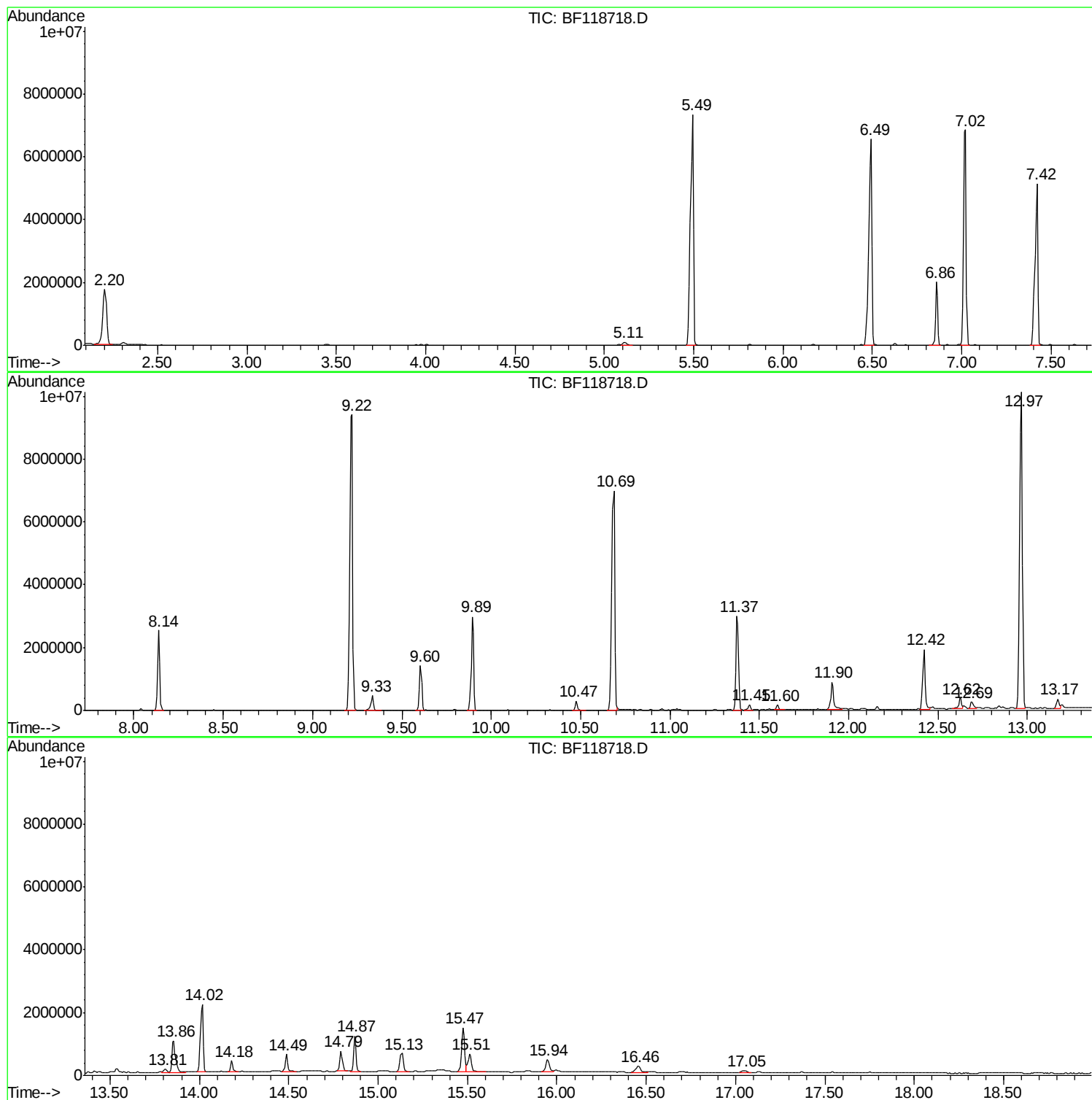
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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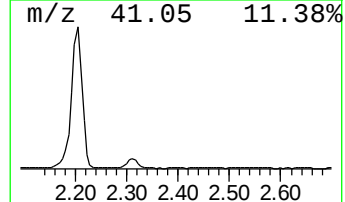
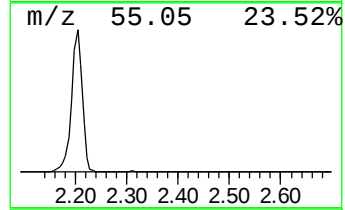
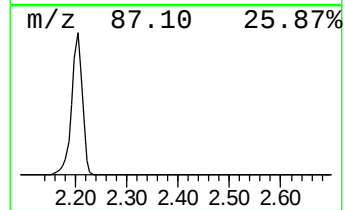
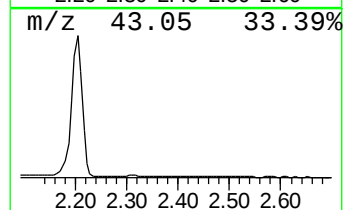
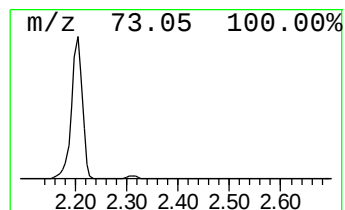
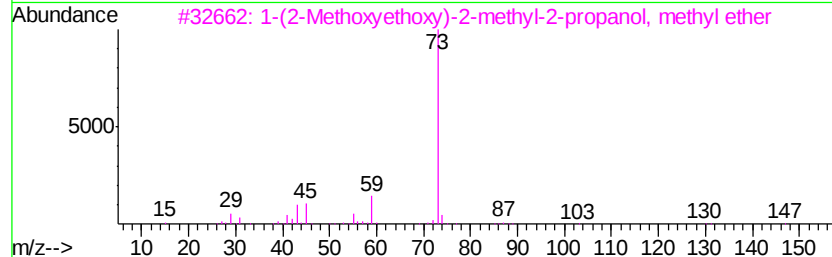
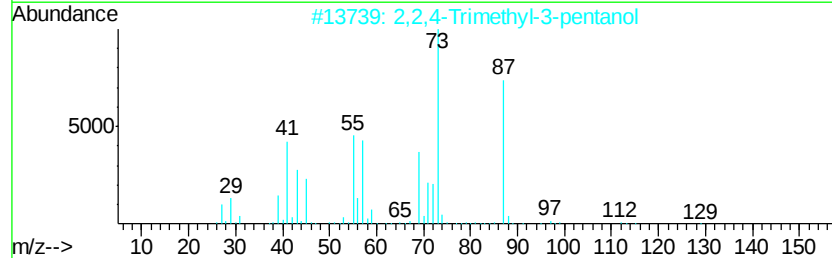
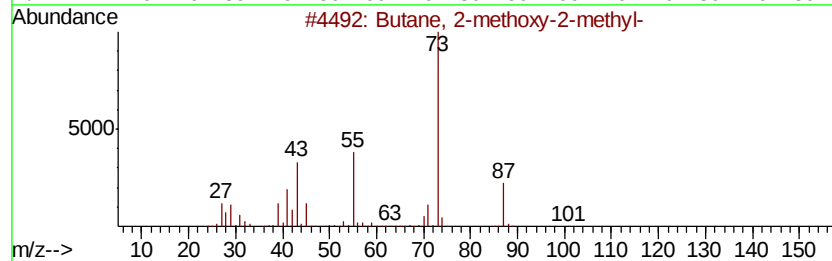
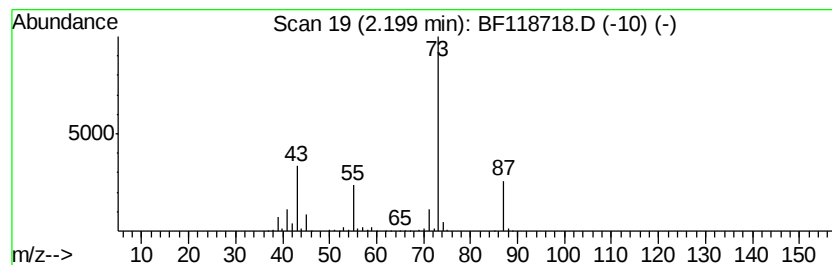
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.20	28.94 ng	2532480	1,4-Dichlorobenzene-d4	6.86	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	1-(2-Methoxyethoxy)-2-methyl-2-propanol	162	C8H18O3	1000364-24-4	25
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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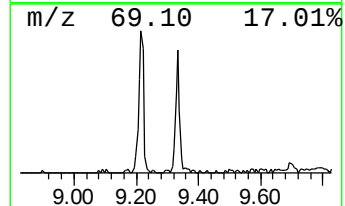
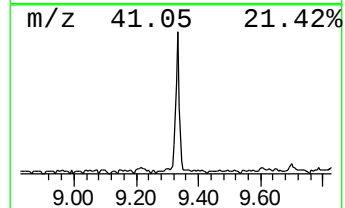
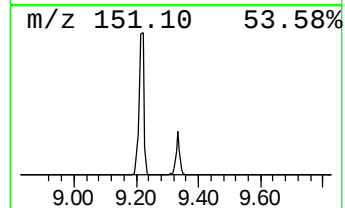
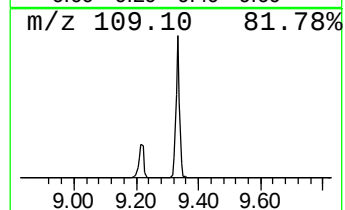
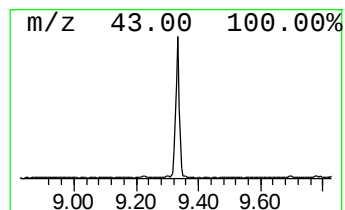
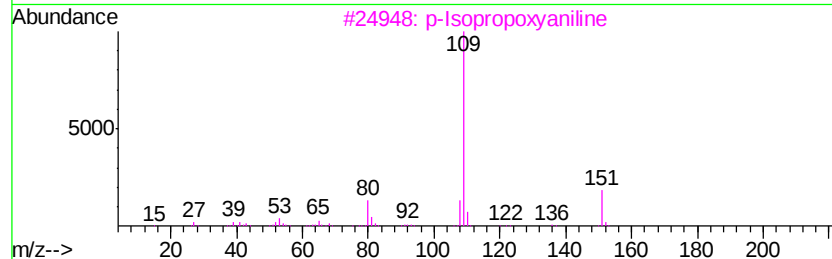
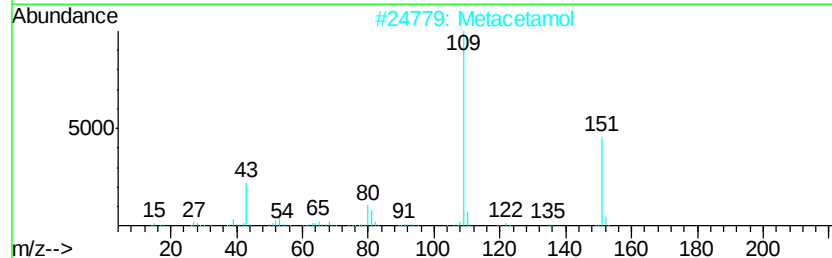
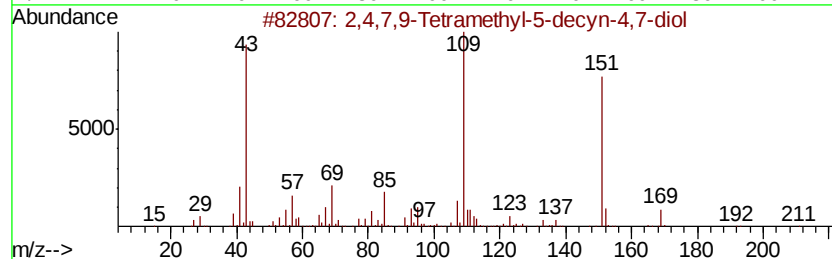
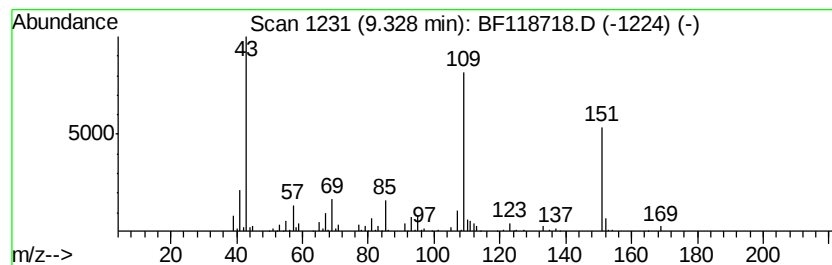
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.33	3.26 ng	445865	Acenaphthene-d10	9.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	64
2		Metacetamol	151	C8H9NO2	000621-42-1	58
3		p-Isopropoxvaniline	151	C9H13NO	007664-66-6	58
4		3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	53
5		m-Isopropoxyaniline	151	C9H13NO	041406-00-2	53



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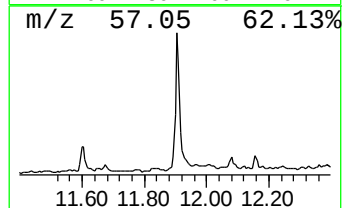
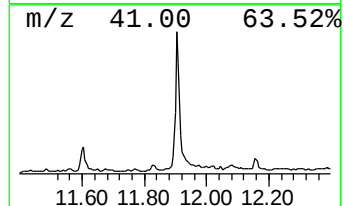
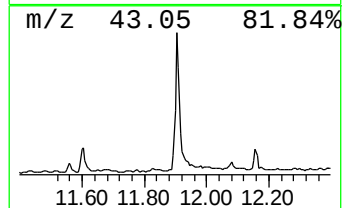
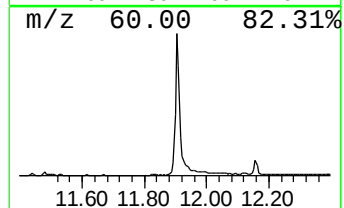
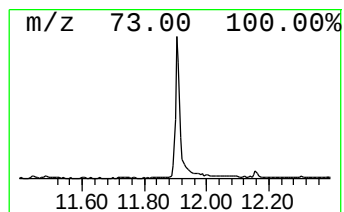
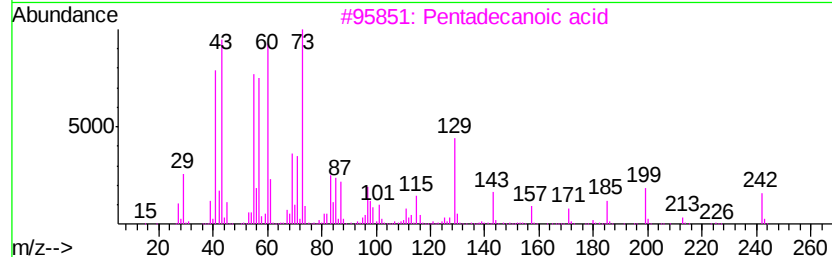
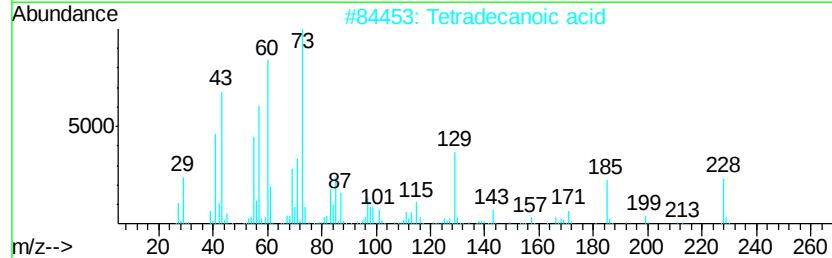
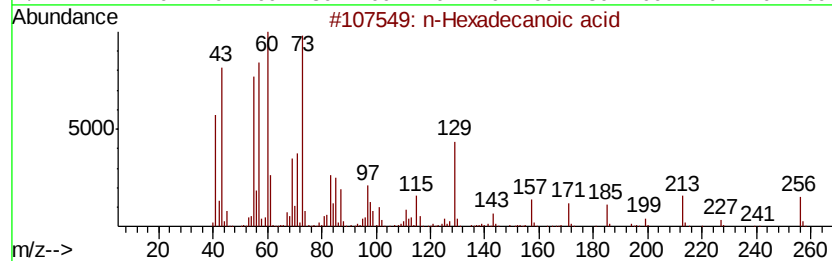
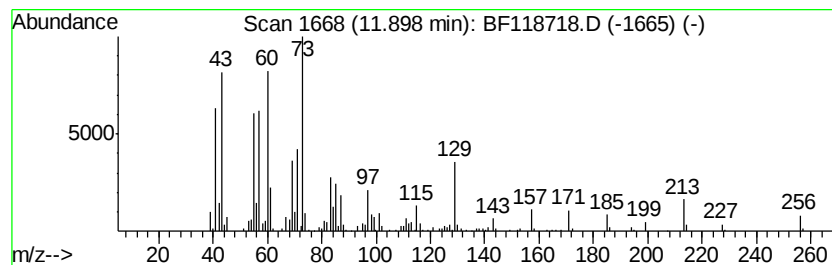
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	6.99 ng	957634	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Tridecanoic acid	214	C13H26O2	000638-53-9	83
5		Octadecanoic acid	284	C18H36O2	000057-11-4	81



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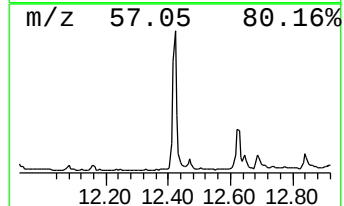
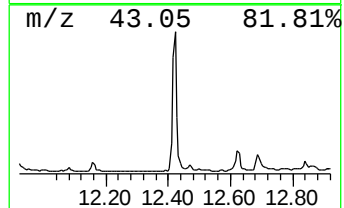
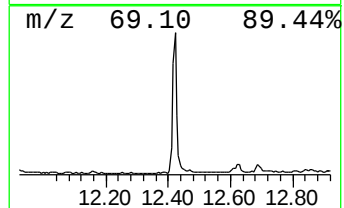
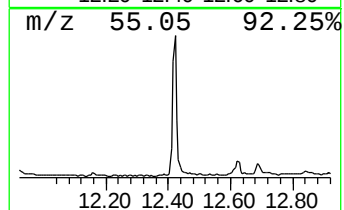
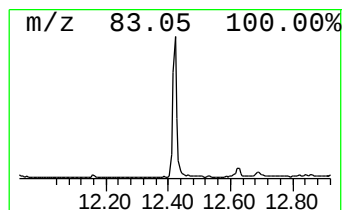
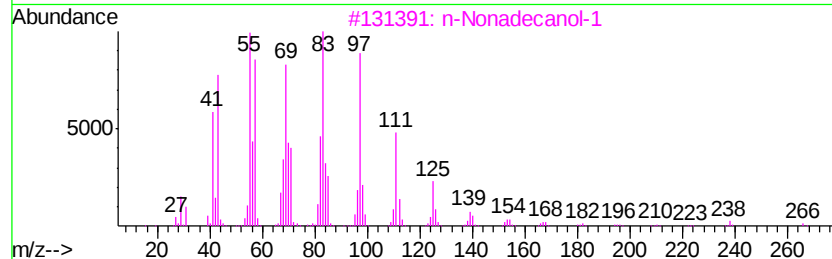
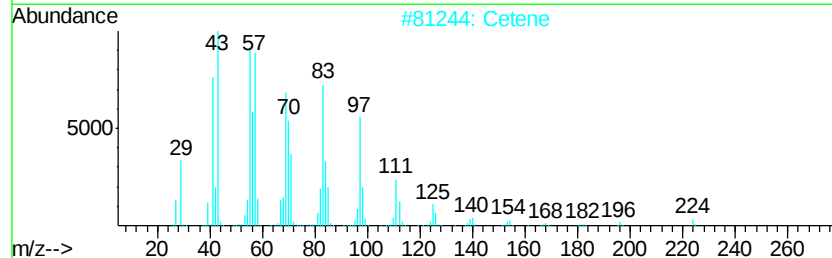
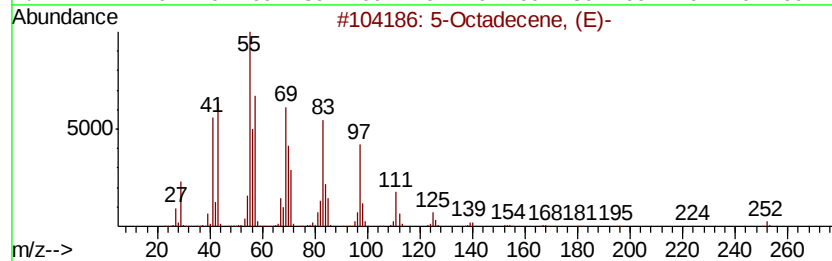
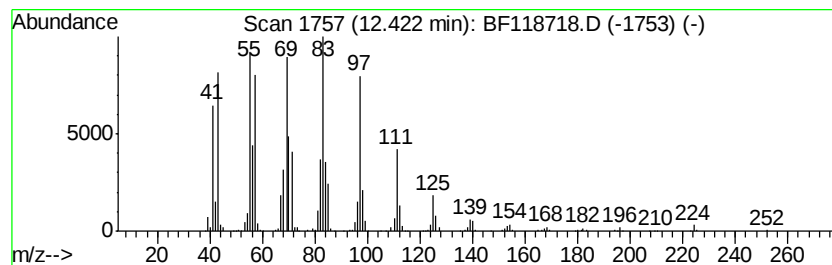
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 5-Octadecene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.42	12.82 ng	1757470	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Octadecene, (E)-	252	C18H36	007206-21-5	97
2		Cetene	224	C16H32	000629-73-2	95
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4		n-Heptadecanol-1	256	C17H36O	001454-85-9	95
5		Cyclohexadecane	224	C16H32	000295-65-8	94



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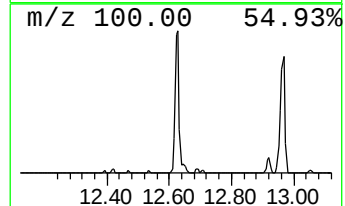
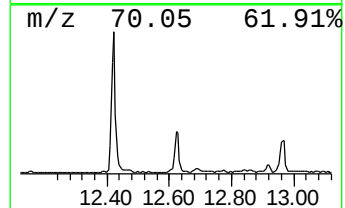
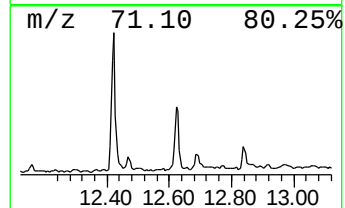
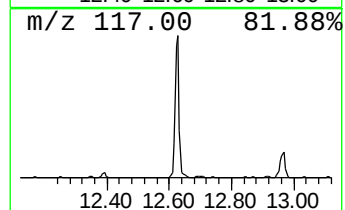
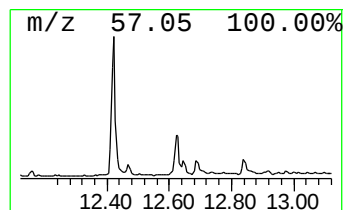
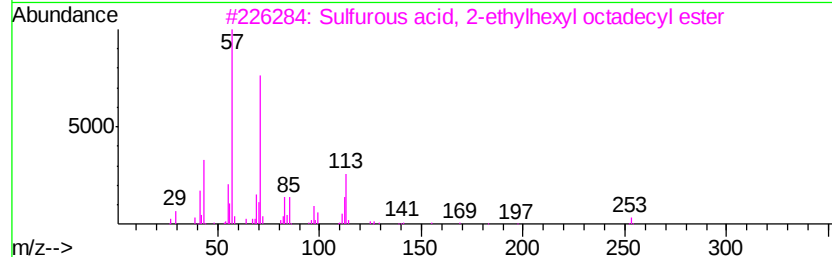
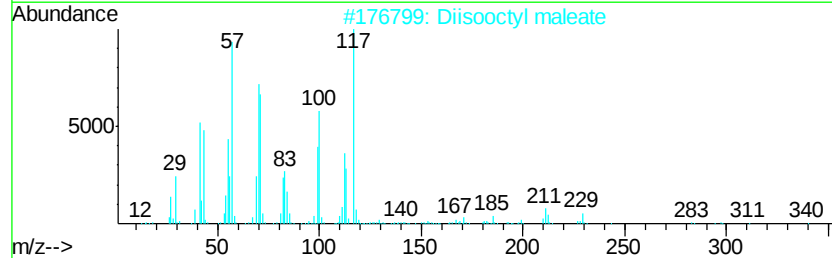
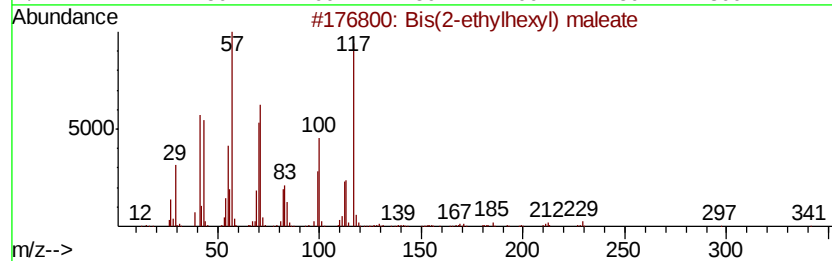
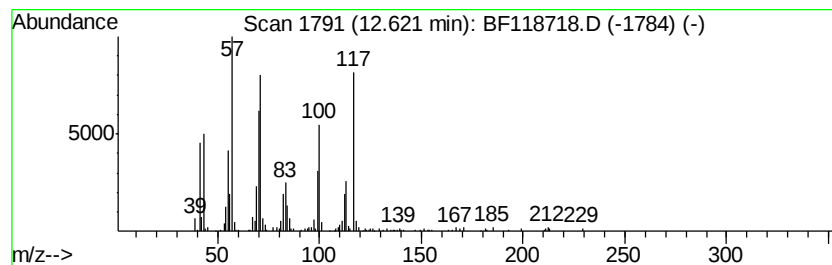
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Bis(2-ethylhexyl) maleate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.62	3.02 ng	414334	Phenanthrene-d10	11.37		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	91
2		Diisooctyl maleate	340	C20H36O4	001330-76-3	52
3		Sulfurous acid, 2-ethylhexyl oct...	446	C26H54O3S	1000309-20-1	35
4		Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	27
5		Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
Data File : BF118718.D
Acq On : 27 Jan 2020 14:09
Operator : CG/JU
Sample : L1241-10
Misc :
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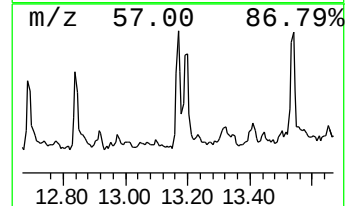
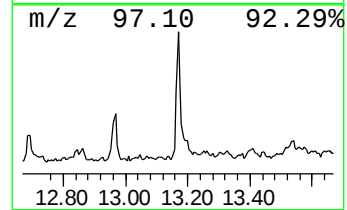
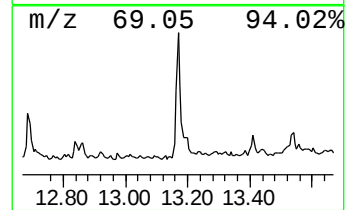
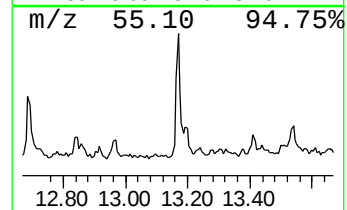
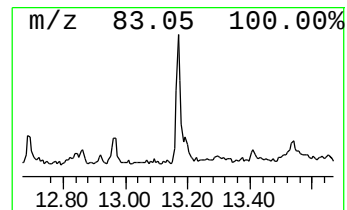
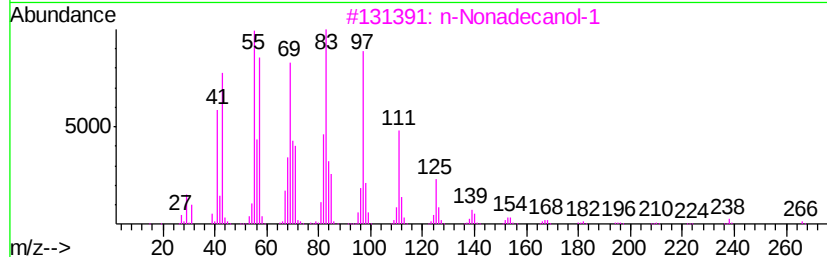
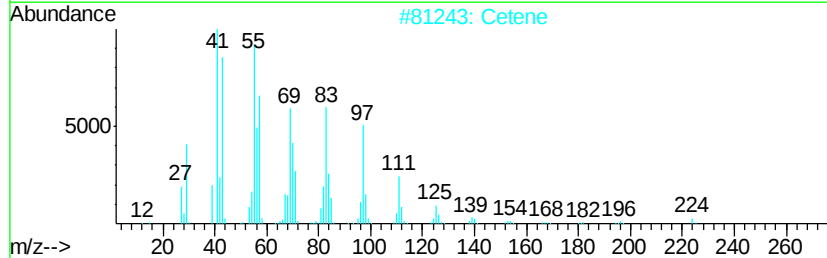
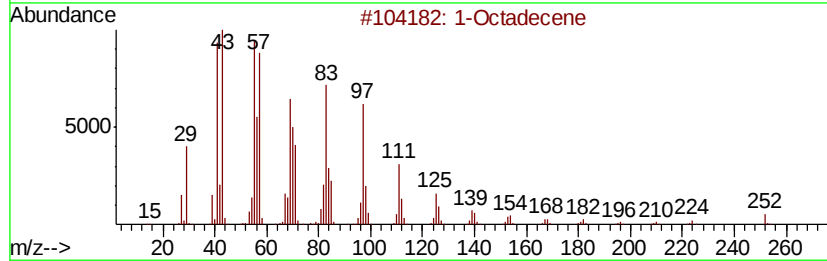
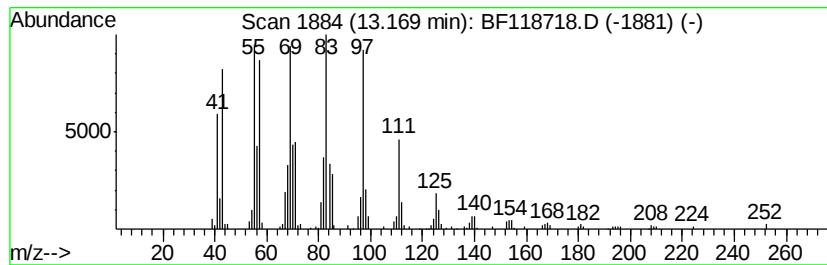
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Octadecene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.17	2.85 ng	304420	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	98
2	Cetene	224	C16H32	000629-73-2	96
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4	5-Octadecene, (E)-	252	C18H36	007206-21-5	95
5	3-Octadecene, (E)-	252	C18H36	007206-19-1	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
Data File : BF118718.D
Acq On : 27 Jan 2020 14:09
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Sample : L1241-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

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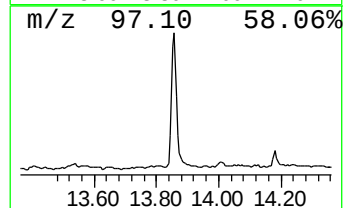
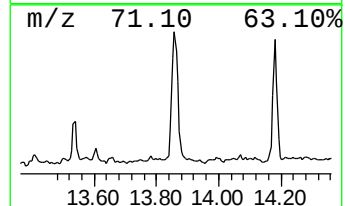
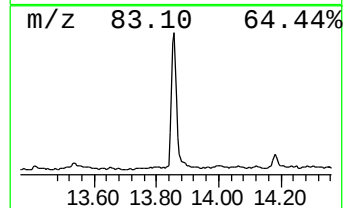
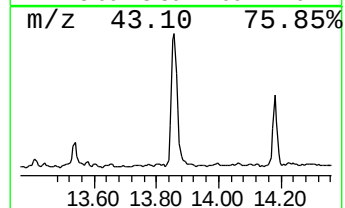
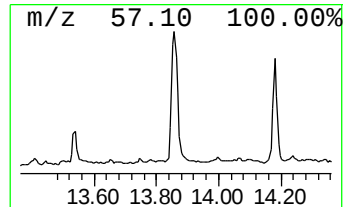
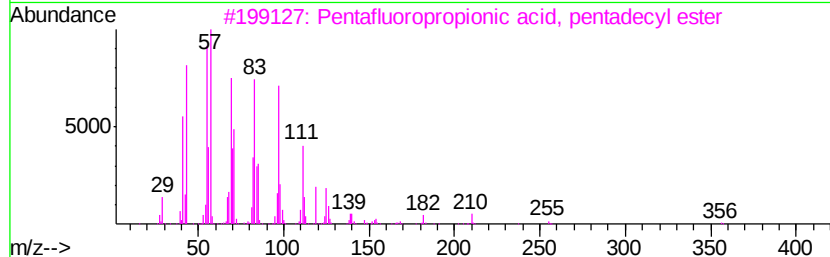
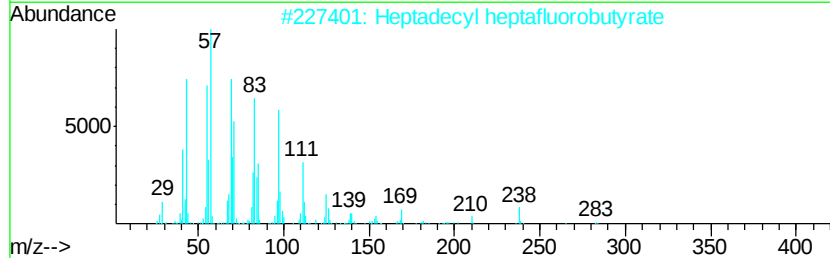
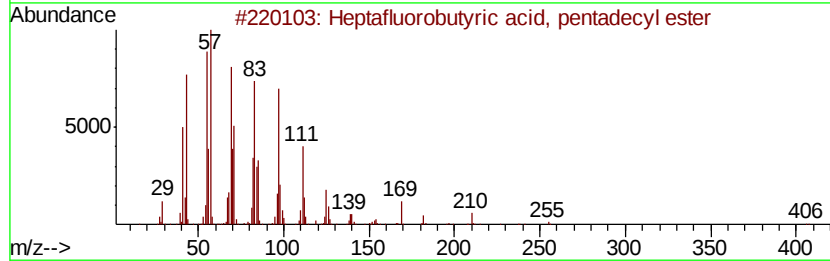
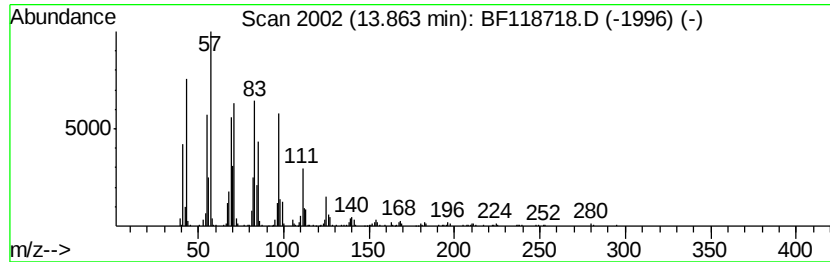
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptafluorobutyric acid, pe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	12.73 ng	1358410	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4		Behenic alcohol	326	C22H46O	000661-19-8	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118718.D
 Acq On : 27 Jan 2020 14:09
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 Sample : L1241-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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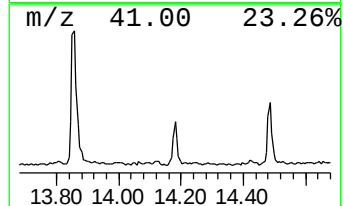
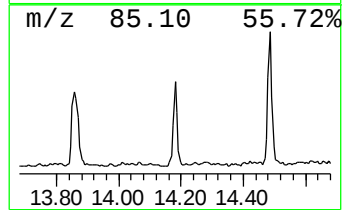
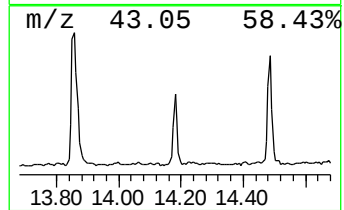
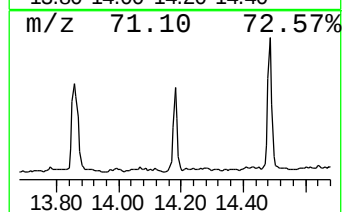
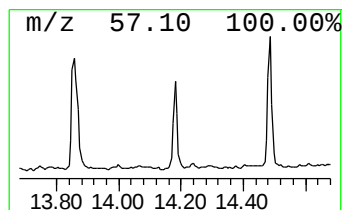
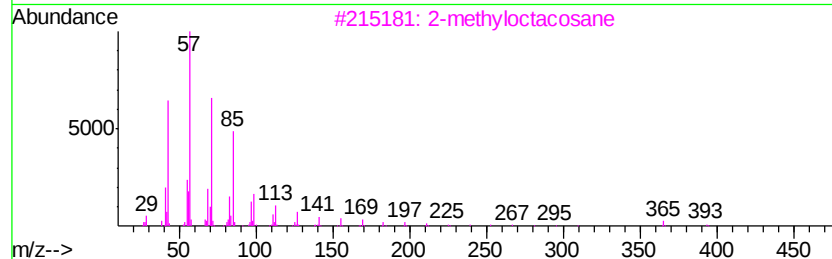
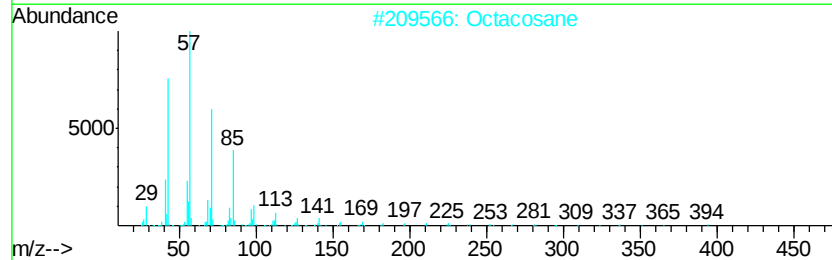
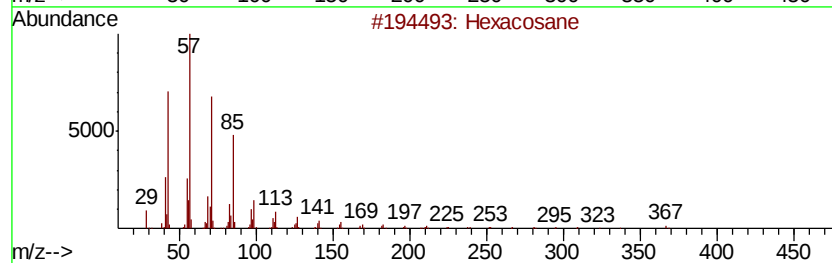
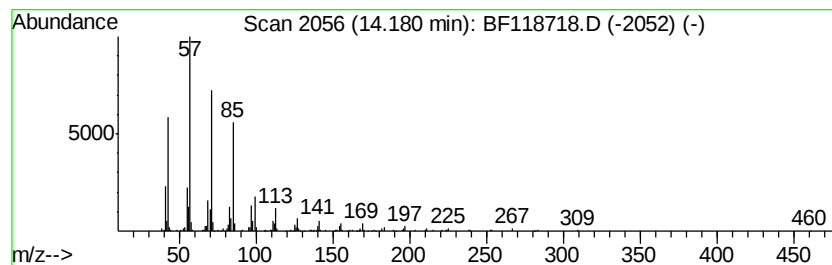
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Hexacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	3.21 ng	343104	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	93
2	Octacosane		394	C28H58	000630-02-4	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Heneicosane		296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
Data File : BF118718.D
Acq On : 27 Jan 2020 14:09
Operator : CG/JU
Sample : L1241-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

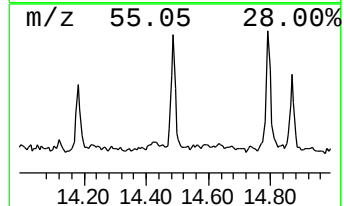
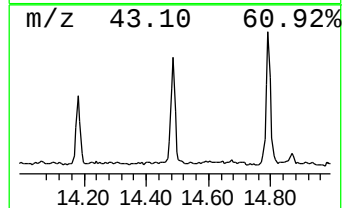
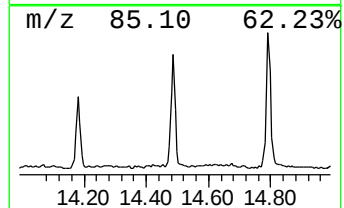
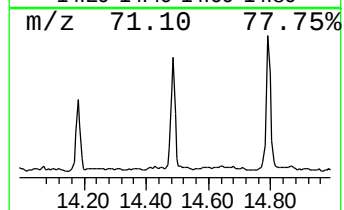
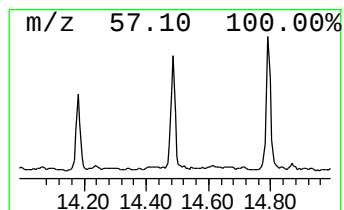
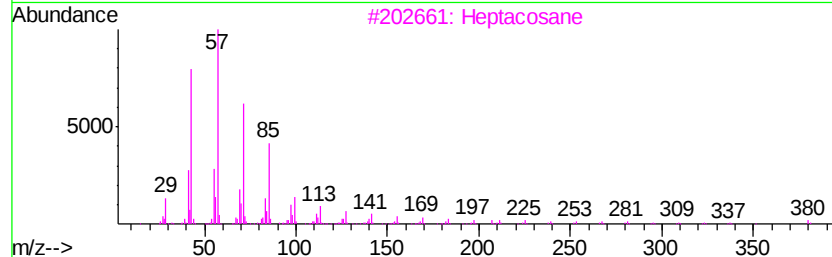
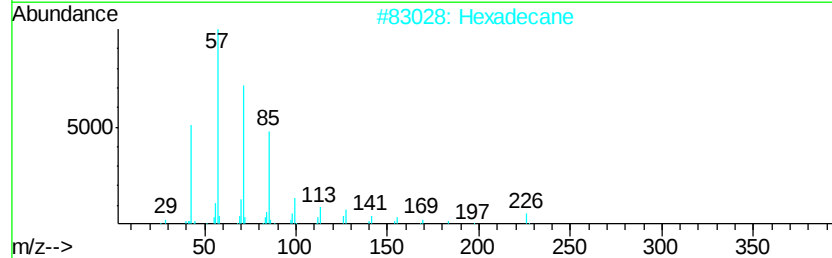
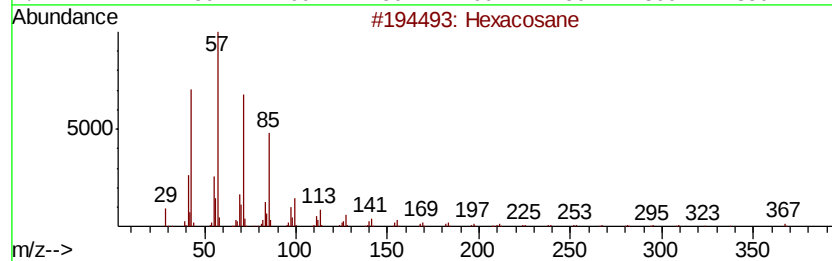
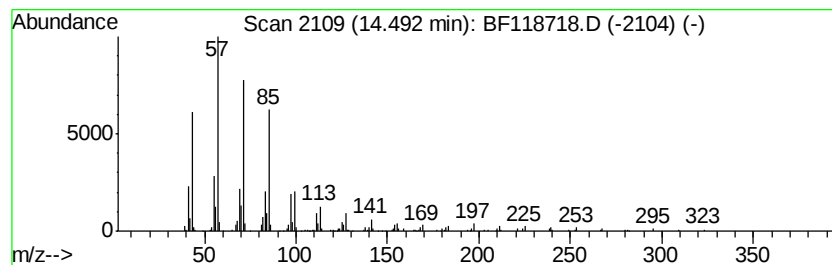
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.49	4.88 ng	521014	Chrysene-d12		14.02
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	94
2	Hexadecane	226	C16H34	000544-76-3	93
3	Heptacosane	380	C27H56	000593-49-7	91
4	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
5	Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
Data File : BF118718.D
Acq On : 27 Jan 2020 14:09
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Sample : L1241-10
Misc :
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Instrument :
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ClientSampleId :
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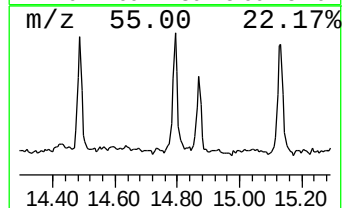
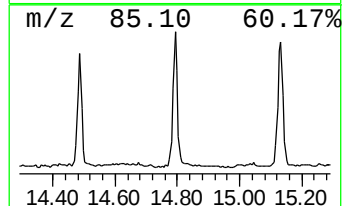
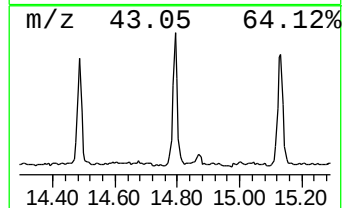
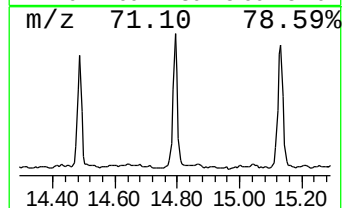
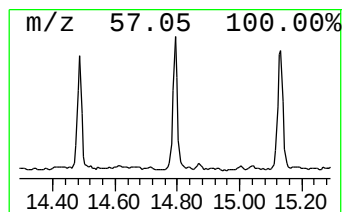
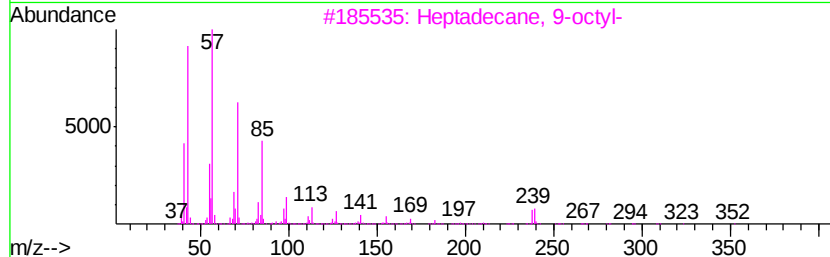
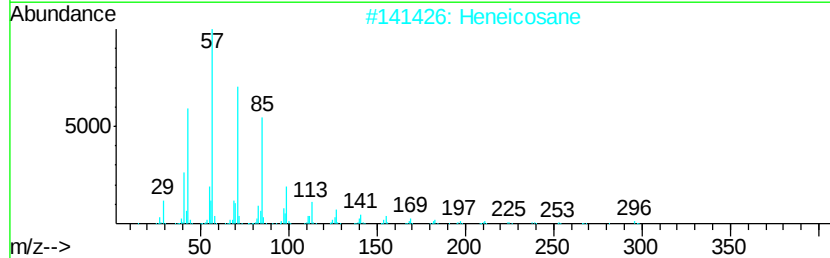
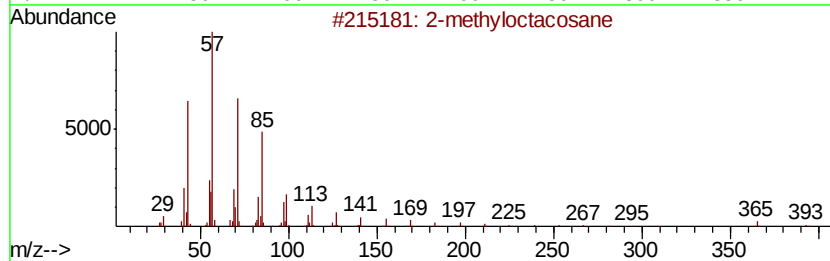
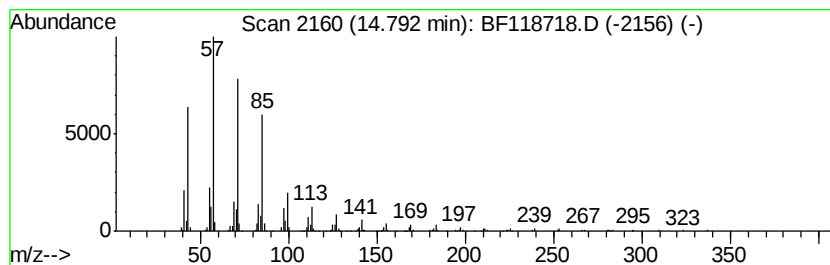
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-methyloctacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	7.09 ng	633092	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
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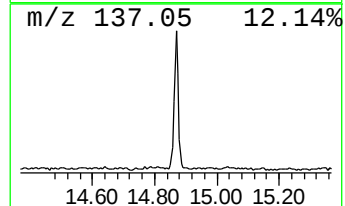
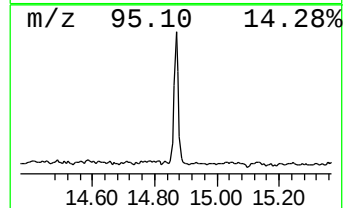
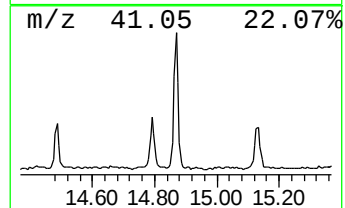
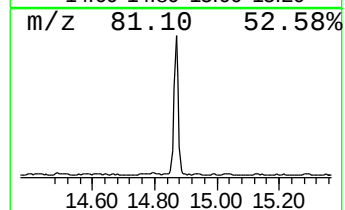
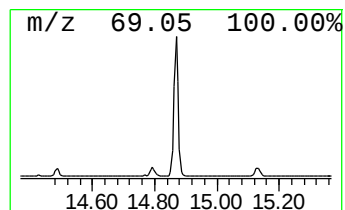
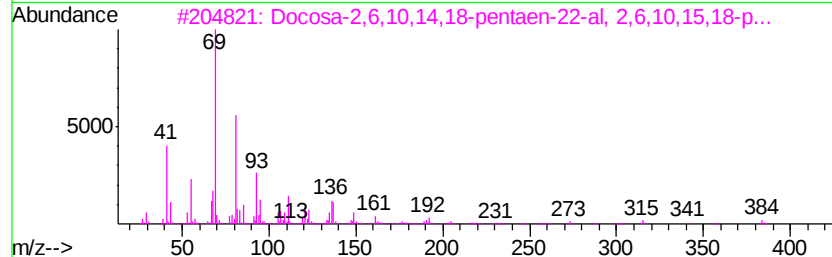
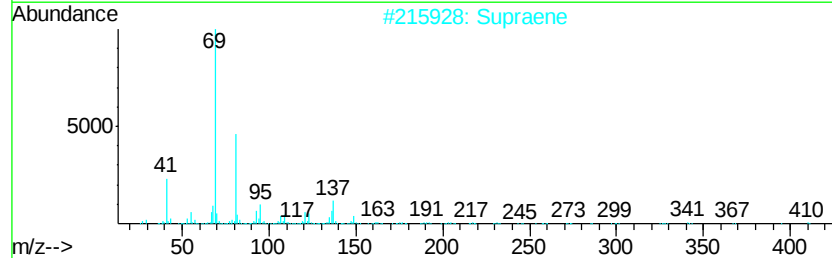
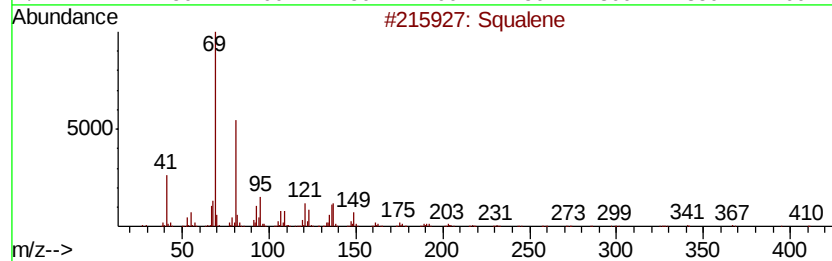
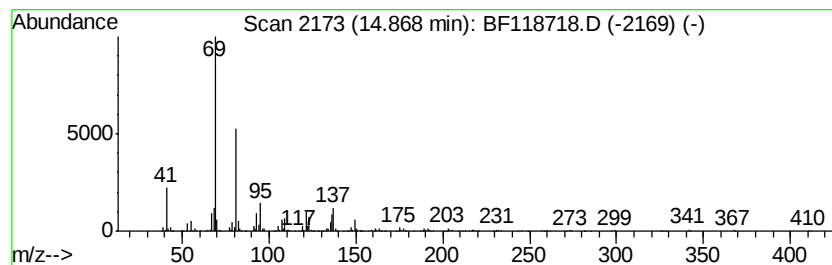
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	12.25 ng	1094370	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	96
2		Supraene	410	C30H50	007683-64-9	91
3		Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	90
4		2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	64
5		1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
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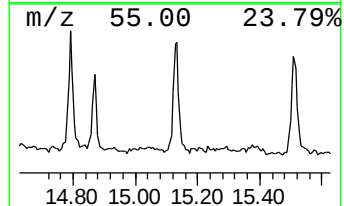
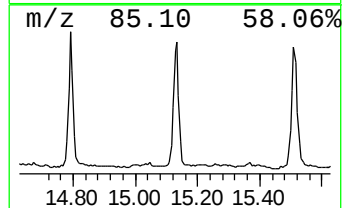
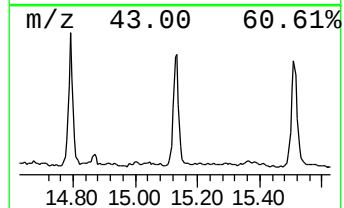
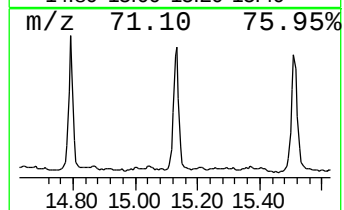
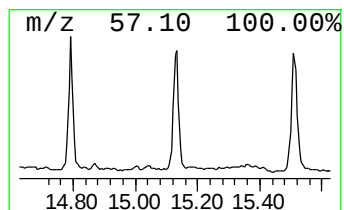
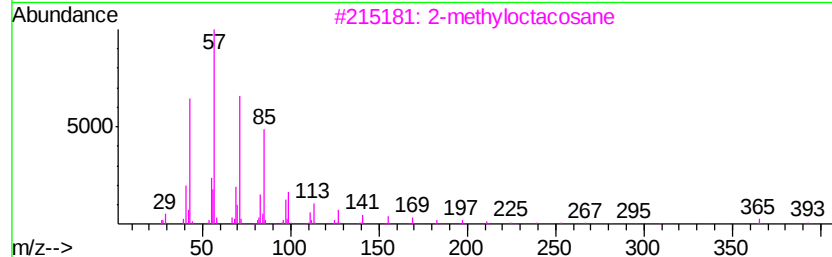
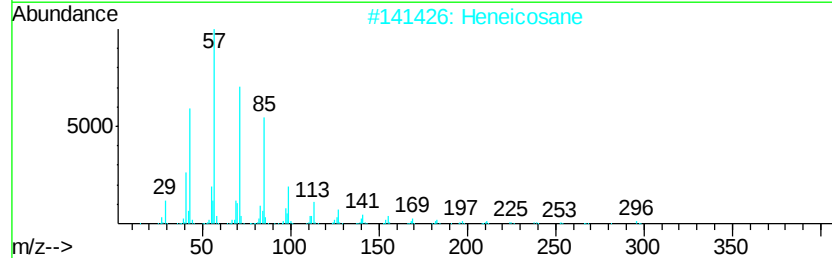
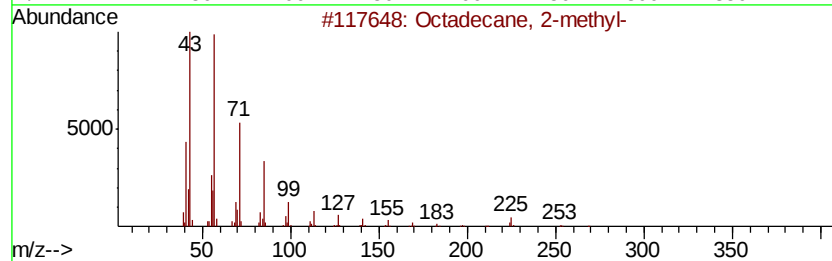
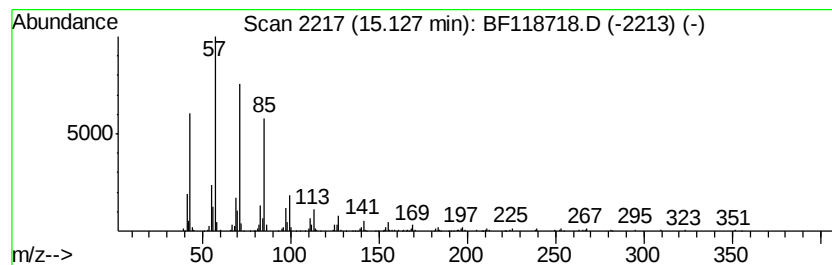
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Octadecane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	7.84 ng	700622	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		triacontane	422	C30H62	000638-68-6	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
Data File : BF118718.D
Acq On : 27 Jan 2020 14:09
Operator : CG/JU
Sample : L1241-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-16

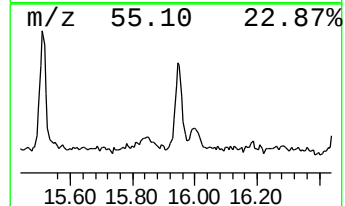
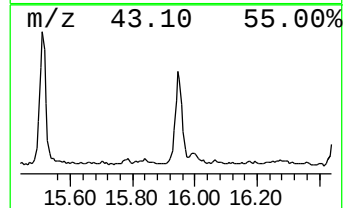
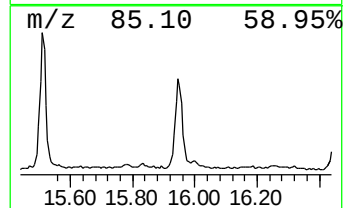
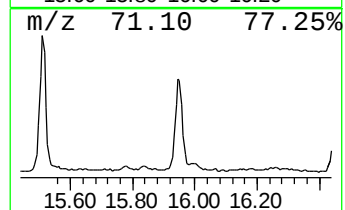
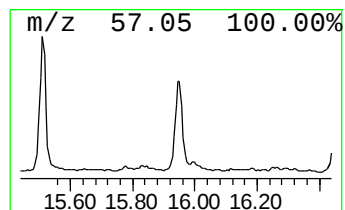
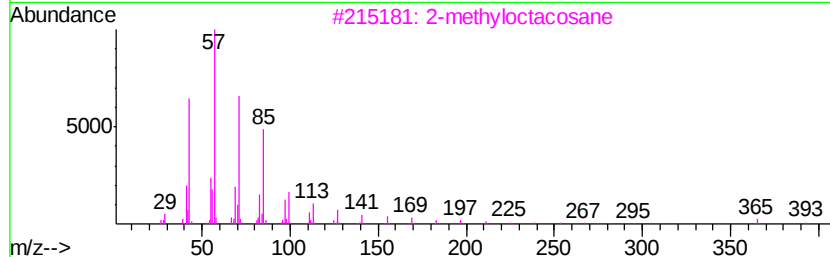
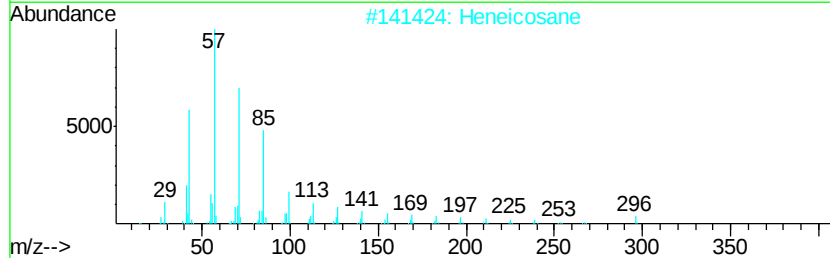
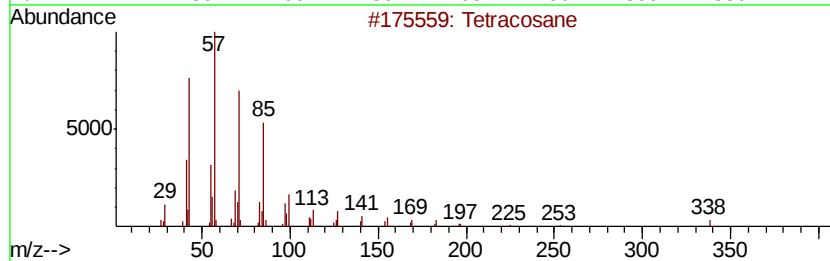
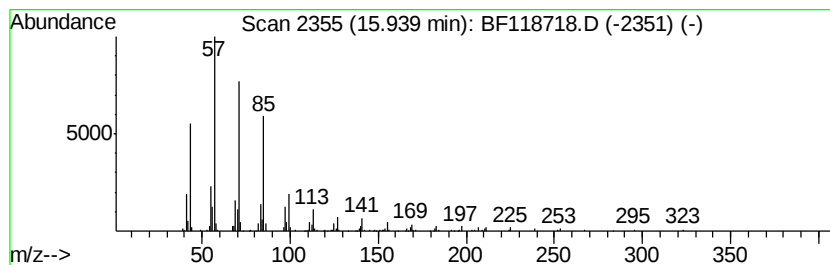
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	6.75 ng	602651	Perylene-d12	15.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			2-methyloctacosane	408	C29H60	1000376-72-8	91
4			Triacontane	422	C30H62	000638-68-6	91
5			Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118718.D
 Acq On : 27 Jan 2020 14:09
 Operator : CG/JU
 Sample : L1241-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-16

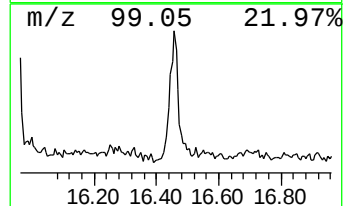
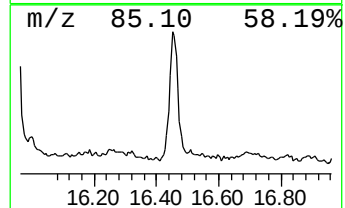
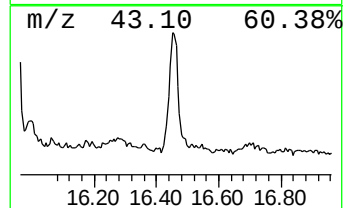
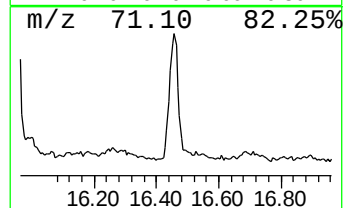
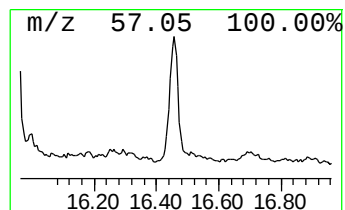
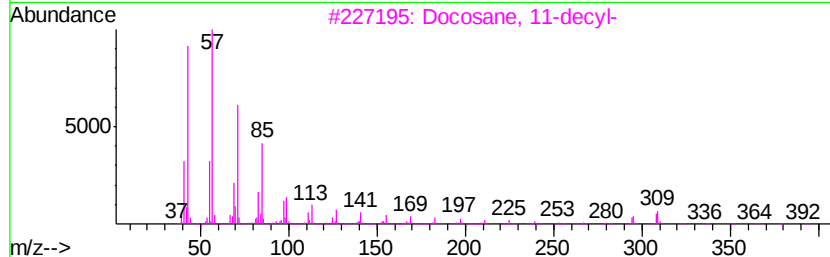
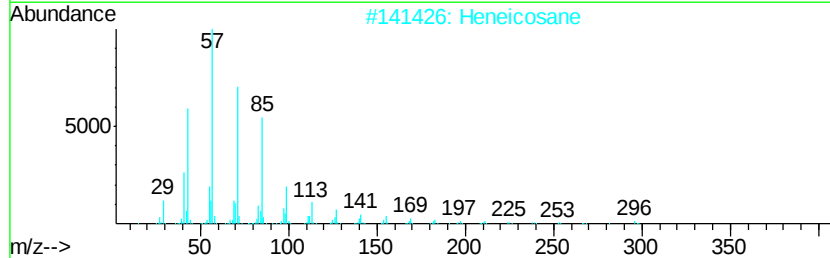
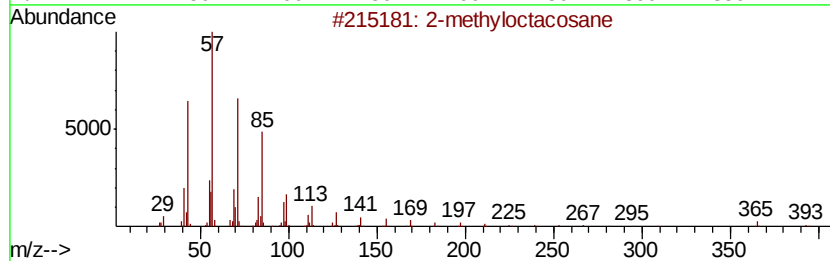
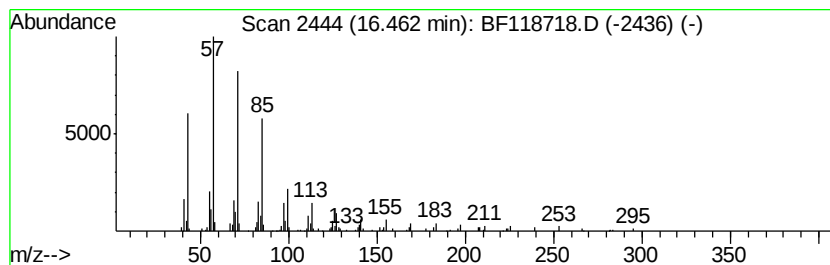
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	4.66 ng	415923	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Docosane, 11-decyl-	451	C32H66	055401-55-3	87
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012720\
 Data File : BF118718.D
 Acq On : 27 Jan 2020 14:09
 Operator : CG/JU
 Sample : L1241-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-16

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.20	28.9	ng	2532480	1	6.86	1750220	20.0
2,4,7,9-Tetrameth...	9.33	3.3	ng	445865	3	9.89	2731430	20.0
n-Hexadecanoic acid	11.90	7.0	ng	957634	4	11.37	2741890	20.0
5-Octadecene, (E)-	12.42	12.8	ng	1757470	4	11.37	2741890	20.0
Bis(2-ethylhexyl)...	12.62	3.0	ng	414334	4	11.37	2741890	20.0
1-Octadecene	13.17	2.9	ng	304420	5	14.02	2134540	20.0
Heptafluorobutyri...	13.86	12.7	ng	1358410	5	14.02	2134540	20.0
Hexacosane	14.18	3.2	ng	343104	5	14.02	2134540	20.0
Hexadecane	14.49	4.9	ng	521014	5	14.02	2134540	20.0
2-methyloctacosane	14.79	7.1	ng	633092	6	15.47	1786710	20.0
Squalene	14.87	12.3	ng	1094370	6	15.47	1786710	20.0
Octadecane, 2-met...	15.13	7.8	ng	700622	6	15.47	1786710	20.0
Tetracosane	15.94	6.8	ng	602651	6	15.47	1786710	20.0
Heneicosane	16.46	4.7	ng	415923	6	15.47	1786710	20.0