

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032620\
 Data File : BF119553.D
 Acq On : 27 Mar 2020 7:48
 Operator : CG/JU
 Sample : L2043-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2A

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-BF031820.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	2.122	3	6	14	rVB	336047	425091	5.18%	0.631%
2	2.228	19	24	33	rVB	115356	150024	1.83%	0.223%
3	5.051	496	504	511	rBV	694261	928696	11.32%	1.379%
4	5.451	566	572	575	rBV	6465843	6293908	76.69%	9.348%
5	6.457	737	743	746	rBV	5678709	6363976	77.55%	9.453%
6	6.834	803	807	811	rBV	1741939	1376455	16.77%	2.044%
7	6.992	830	834	837	rBV	5703506	4975571	60.63%	7.390%
8	7.398	897	903	906	rBV	4193278	4147593	50.54%	6.160%
9	8.116	1021	1025	1029	rBV	1947678	1791825	21.83%	2.661%
10	9.198	1203	1209	1212	rBV	7591609	7622473	92.88%	11.322%
11	9.316	1221	1229	1233	rVB	316651	294450	3.59%	0.437%
12	9.592	1271	1276	1279	rBV	437692	357883	4.36%	0.532%
13	9.880	1319	1325	1328	rBV	2359703	2171842	26.46%	3.226%
14	10.669	1453	1459	1462	rBV	5449651	5249910	63.97%	7.798%
15	10.769	1473	1476	1486	rVB	257858	236418	2.88%	0.351%
16	11.369	1573	1578	1581	rBV	2524850	2346796	28.60%	3.486%
17	11.898	1663	1668	1684	rBV	1317458	1516931	18.48%	2.253%
18	12.686	1798	1802	1816	rBV	869613	939086	11.44%	1.395%
19	12.963	1842	1849	1852	rBV	8056460	8206526	100.00%	12.189%
20	13.439	1922	1930	1936	rBV	156127	141153	1.72%	0.210%
21	13.863	1994	2002	2022	rBV	721152	1350839	16.46%	2.006%
22	14.010	2022	2027	2030	rBV	2460440	2173571	26.49%	3.228%
23	14.180	2053	2056	2060	rBV	601839	505042	6.15%	0.750%
24	14.486	2105	2108	2112	rBV	934644	818583	9.97%	1.216%
25	14.792	2157	2160	2165	rVB	1045185	1067768	13.01%	1.586%
26	15.133	2214	2218	2226	rBV	991012	1135016	13.83%	1.686%
27	15.474	2270	2276	2280	rBV	1577887	1921968	23.42%	2.855%
28	15.521	2280	2284	2291	rVB	818563	1072731	13.07%	1.593%
29	15.957	2353	2358	2366	rBV	541933	819523	9.99%	1.217%
30	16.468	2438	2445	2452	rBV2	259119	457613	5.58%	0.680%
31	17.021	2528	2539	2541	rBV3	55354	171999	2.10%	0.255%
32	17.062	2543	2546	2556	rVB3	99971	210717	2.57%	0.313%
33	17.780	2662	2668	2677	rVB2	32143	83714	1.02%	0.124%

LSC Area Percent Report

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
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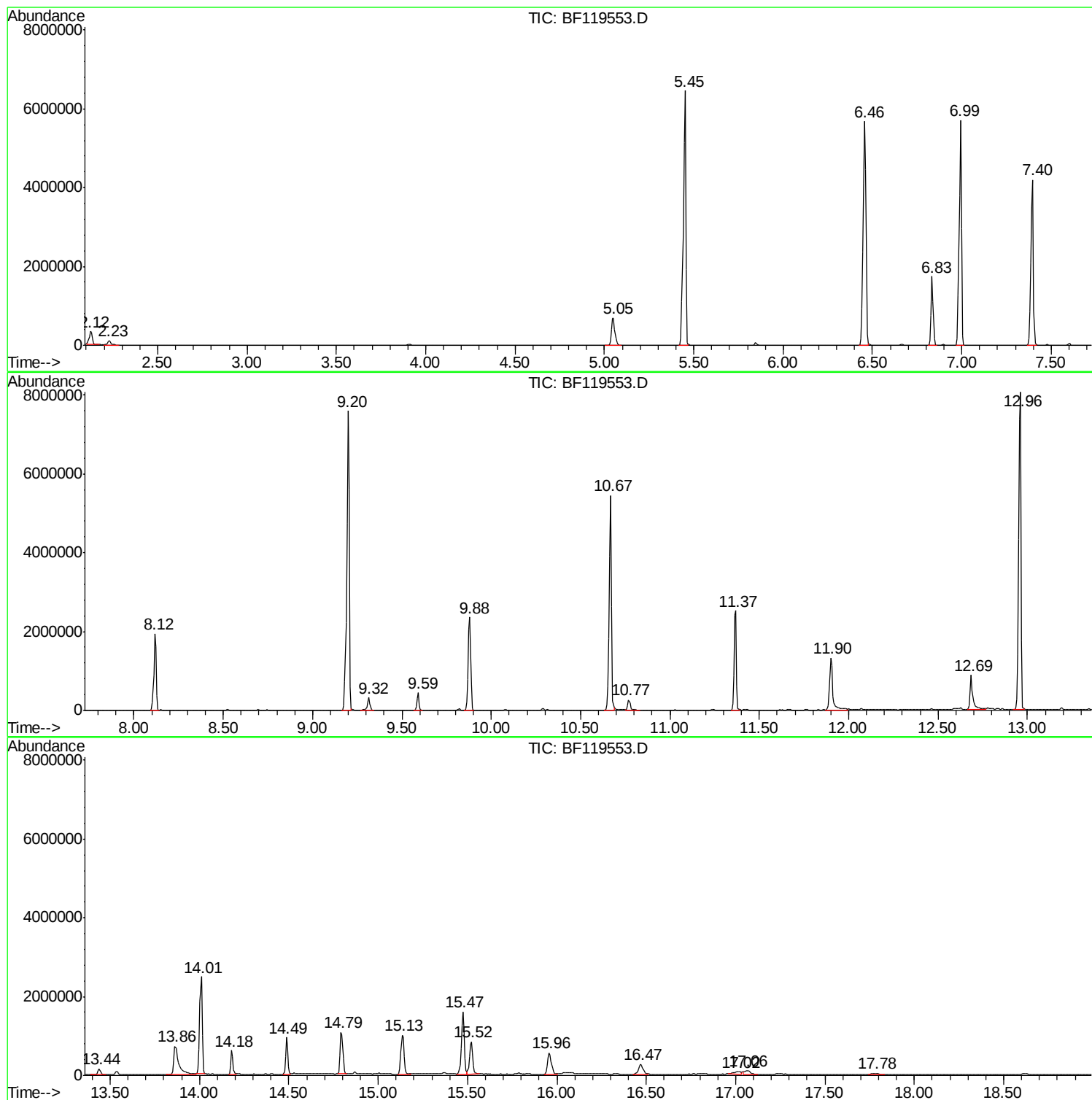
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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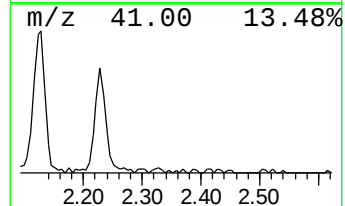
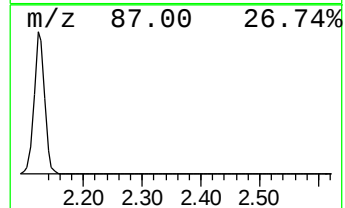
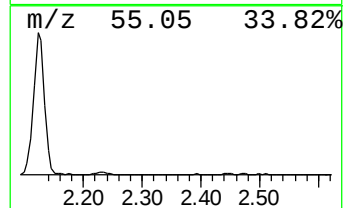
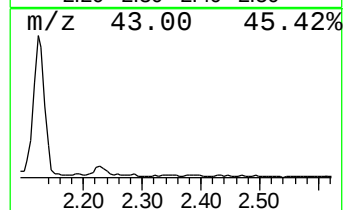
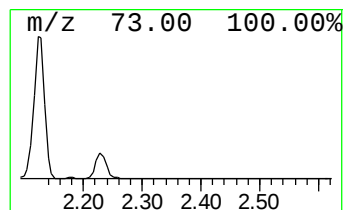
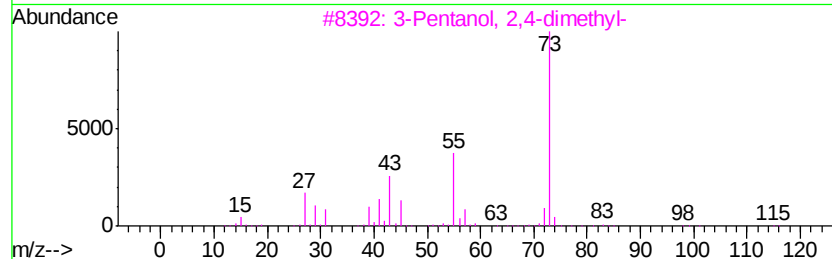
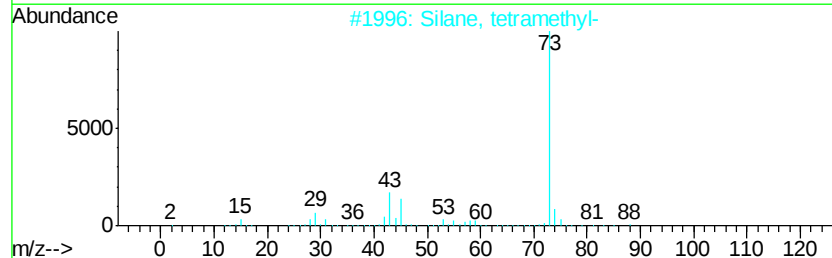
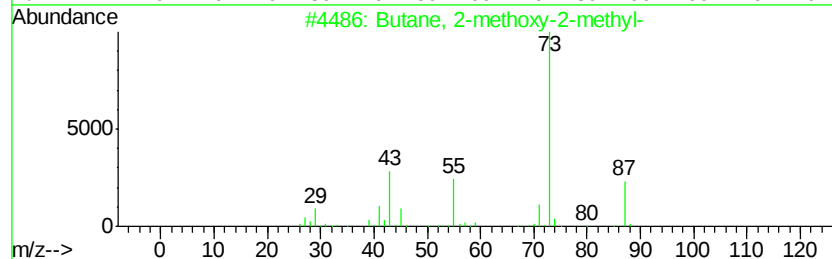
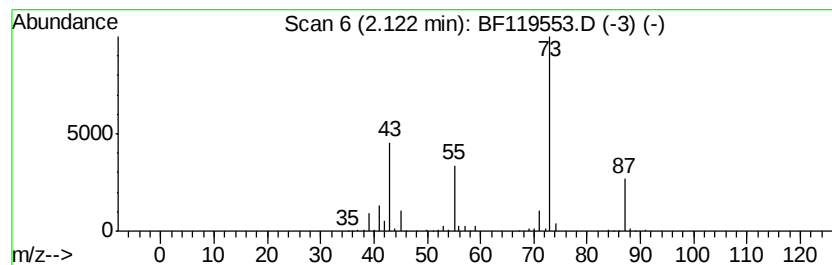
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	6.18 ng	425091	1,4-Dichlorobenzene-d4	6.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	23



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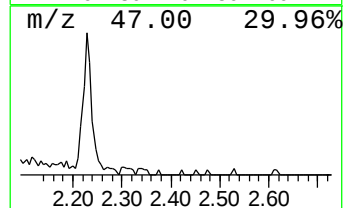
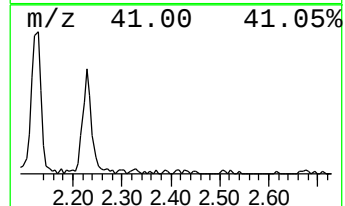
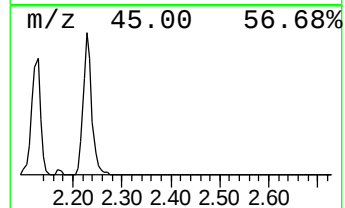
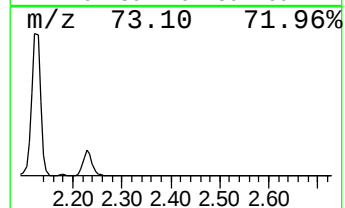
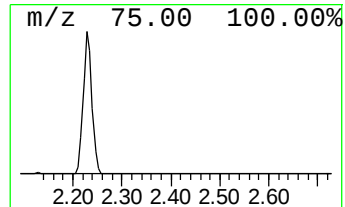
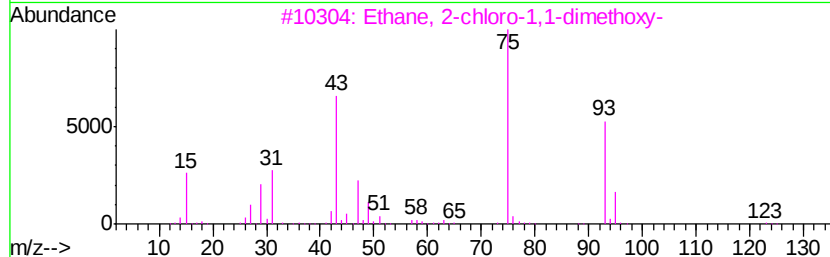
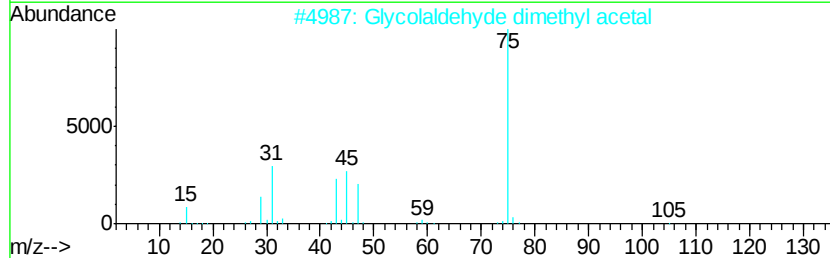
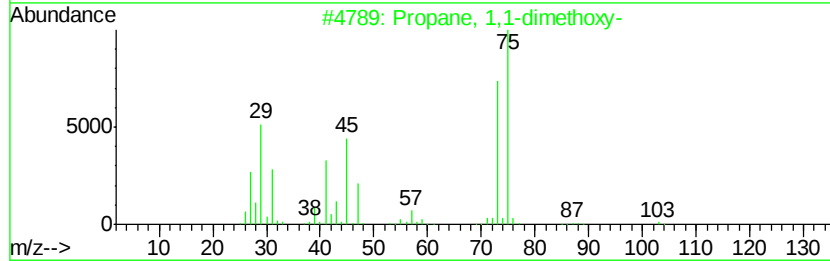
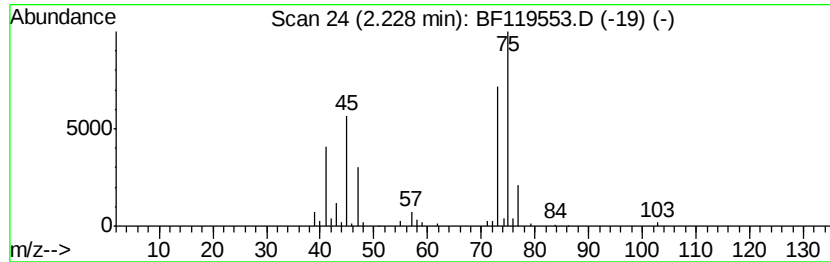
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propane, 1,1-dimethoxy- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	2.18 ng	150024	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	53
2		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	33
3		Ethane, 2-chloro-1,1-dimethoxy-	124	C4H9ClO2	000097-97-2	28
4		Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	28
5		Ethanethioamide	75	C2H5NS	000062-55-5	9



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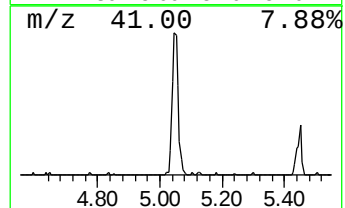
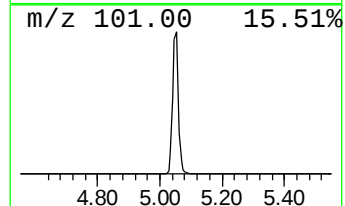
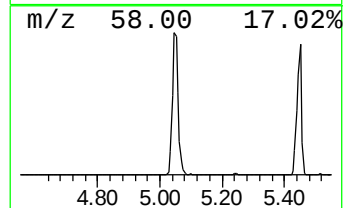
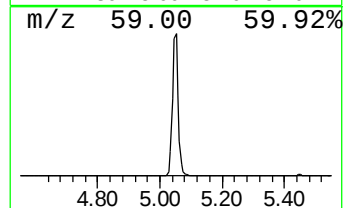
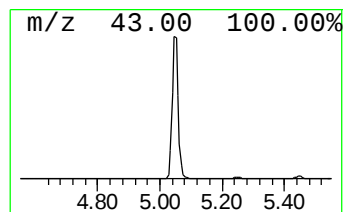
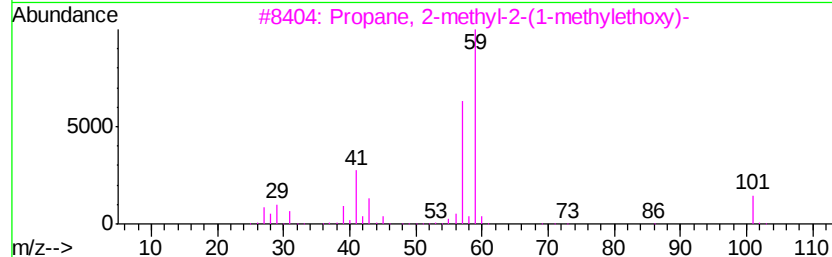
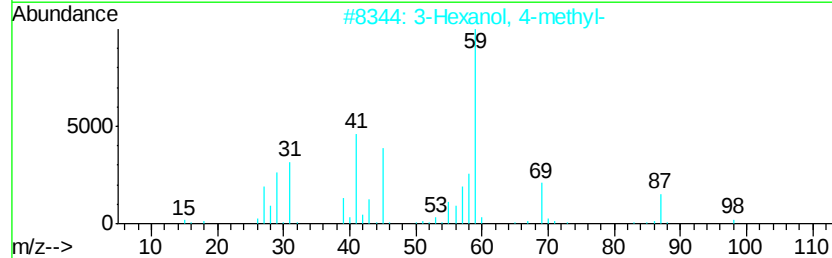
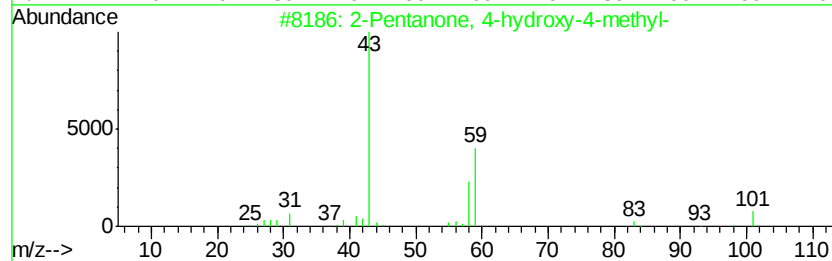
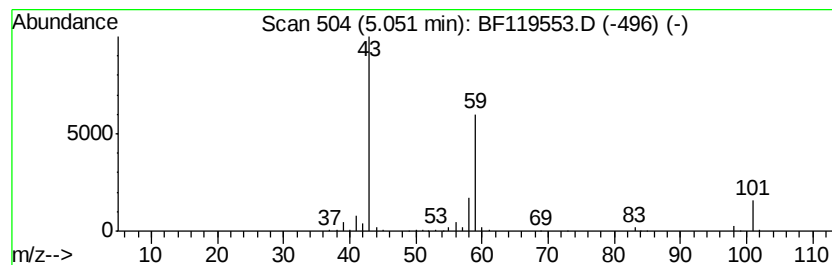
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	13.49 ng	928696	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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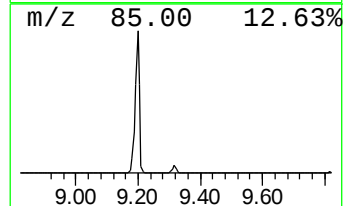
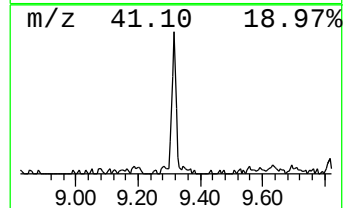
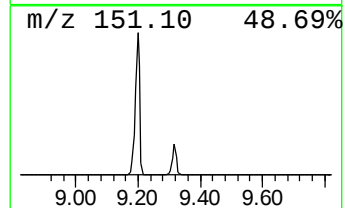
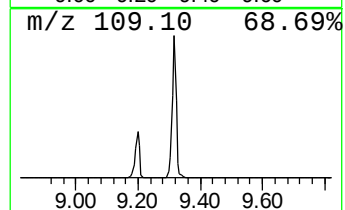
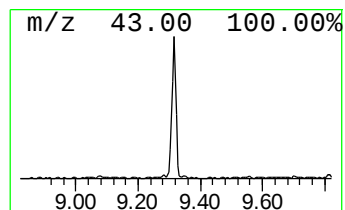
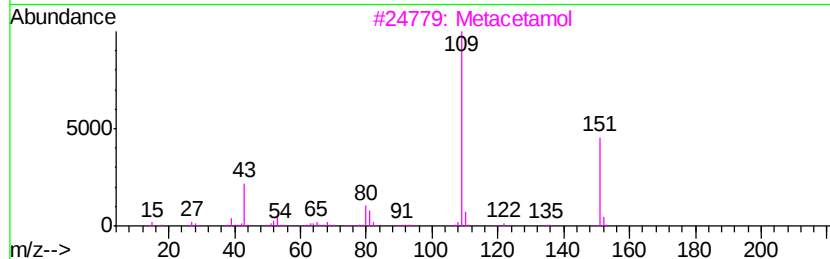
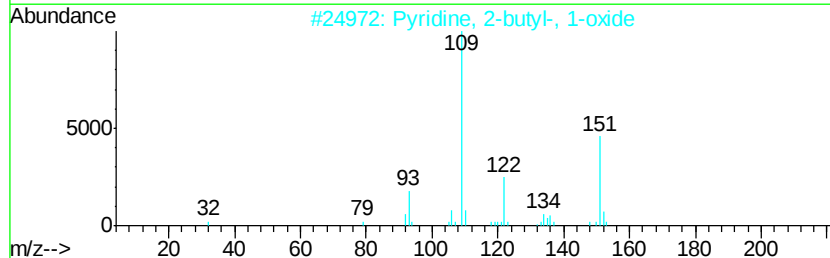
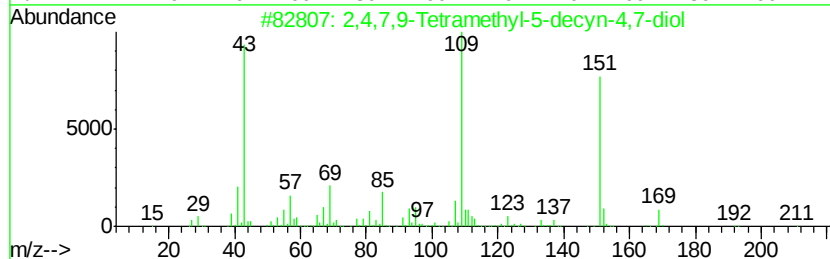
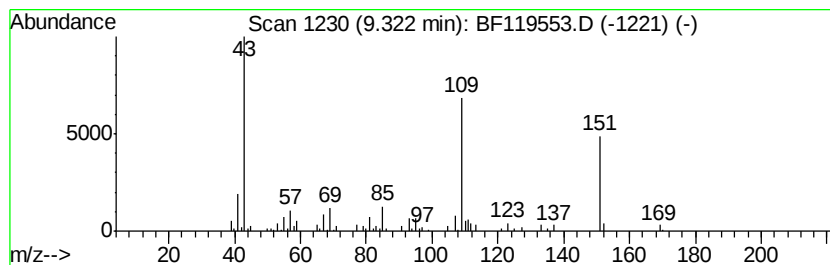
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.32	2.71 ng	294450	Acenaphthene-d10	9.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2		000126-86-3	86
2	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO		031396-32-4	53
3	Metacetamol	151	C8H9NO2		000621-42-1	49
4	Acetaminophen	151	C8H9NO2		000103-90-2	47
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2		001006-96-8	47



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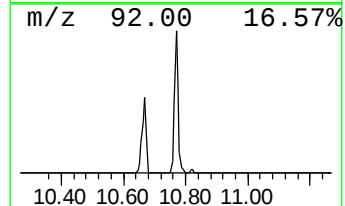
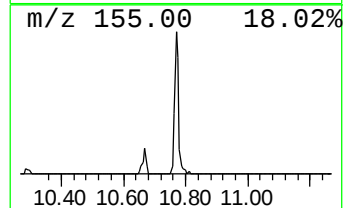
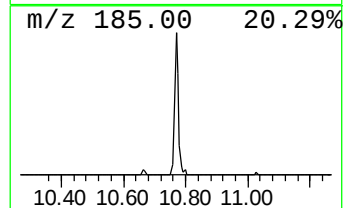
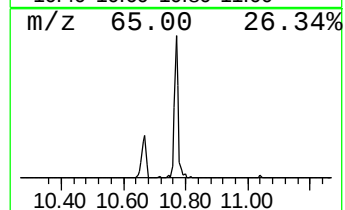
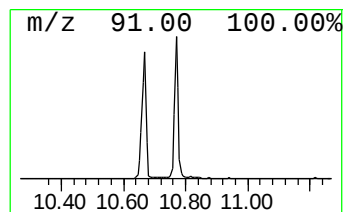
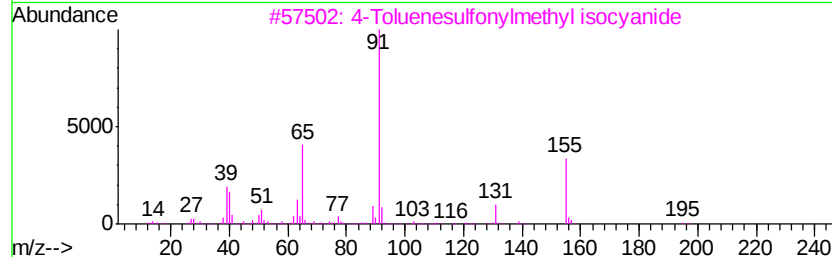
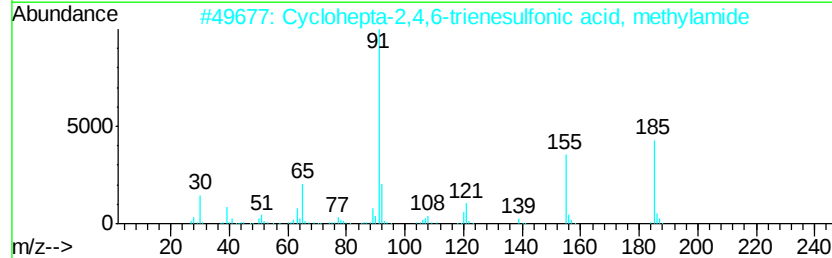
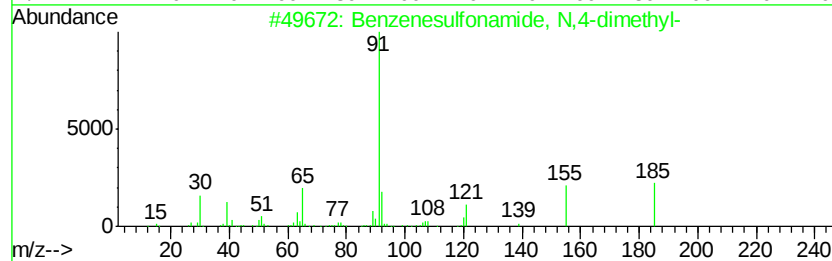
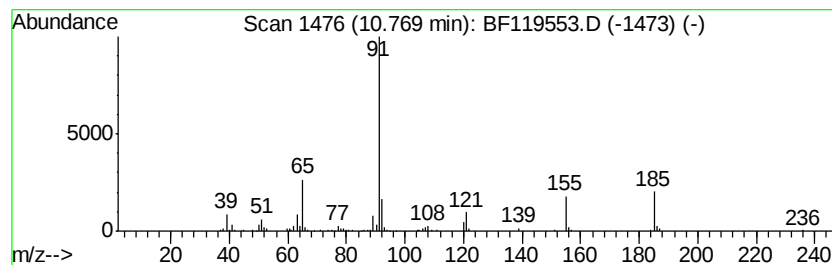
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzenesulfonamide, N,4-dim... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	2.01 ng	236418	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	93
2		Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	49
3		4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	47
4		Benzoic acid, 2-hydroxy-, phenyl...	228	C14H12O3	000118-58-1	43
5		Benzenesulfonamide, N,4-dimethyl...	214	C8H10N2O3S	000080-11-5	38



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Acq On : 27 Mar 2020 7:48
Operator : CG/JU
Sample : L2043-02
Misc :
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ClientSampleId :
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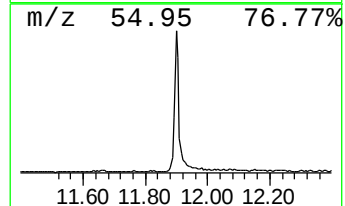
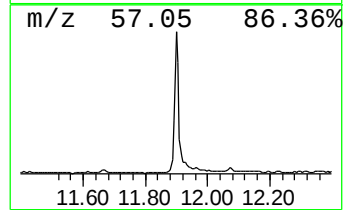
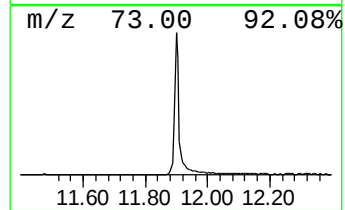
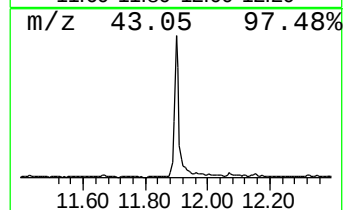
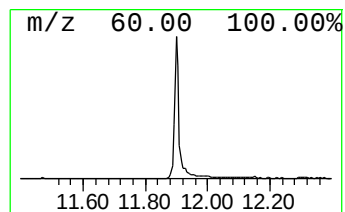
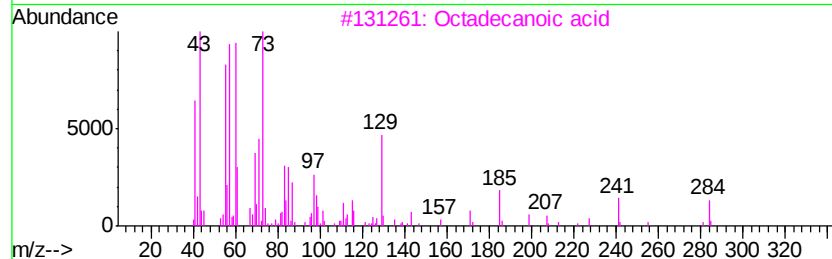
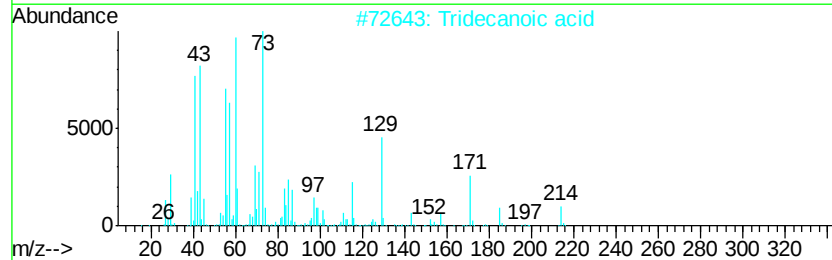
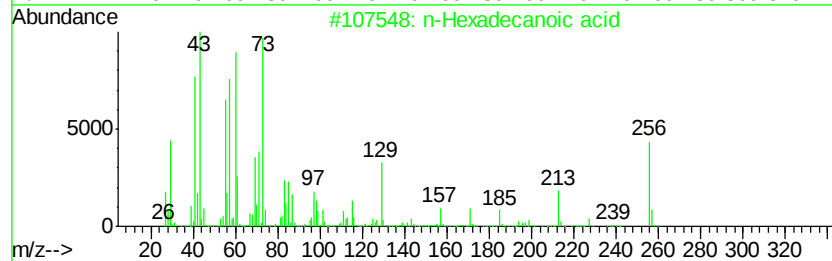
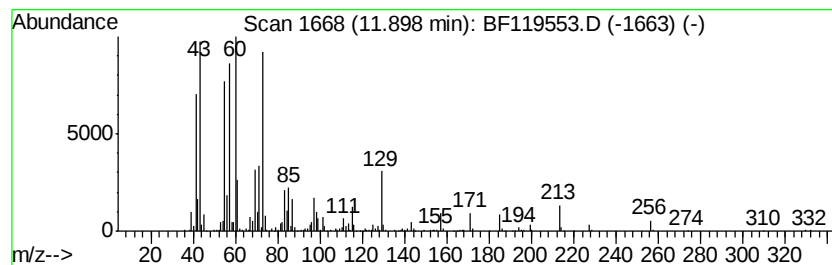
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	12.93 ng	1516930	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Octadecanoic acid	284	C18H36O2	000057-11-4	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
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Acq On : 27 Mar 2020 7:48
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ClientSampleId :
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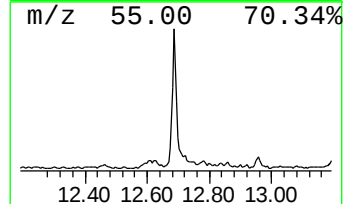
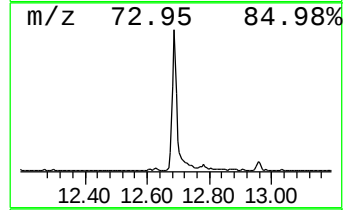
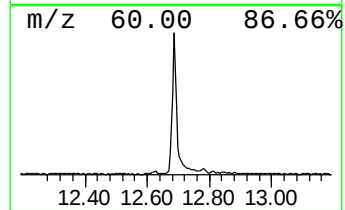
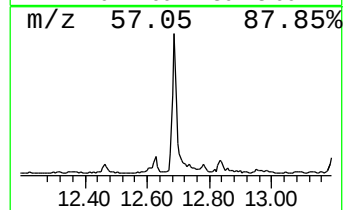
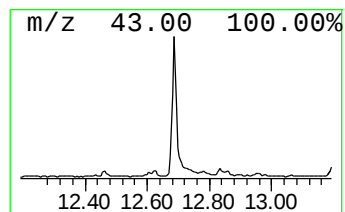
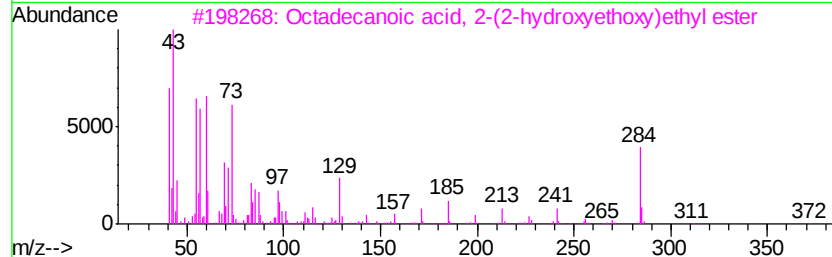
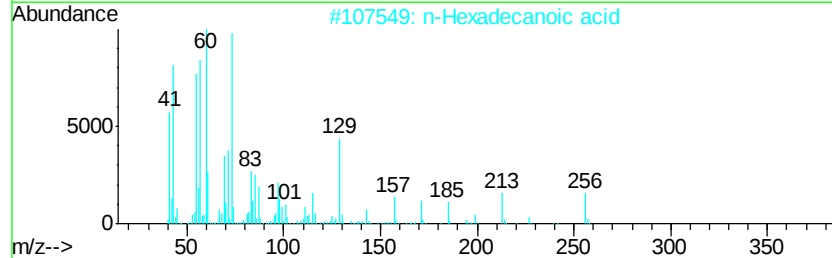
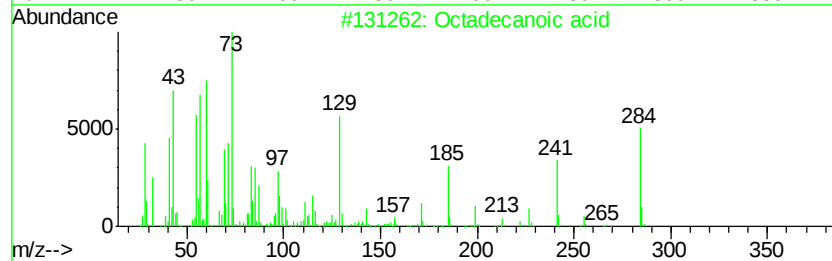
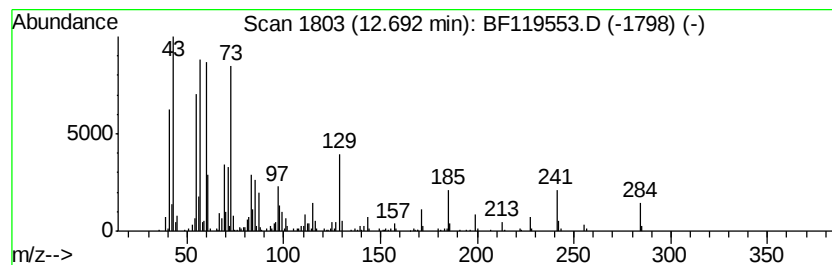
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	8.00 ng	939086	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
4		Tridecanoic acid	214	C13H26O2	000638-53-9	59
5		n-Decanoic acid	172	C10H20O2	000334-48-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
Data File : BF119553.D
Acq On : 27 Mar 2020 7:48
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ClientSampleId :
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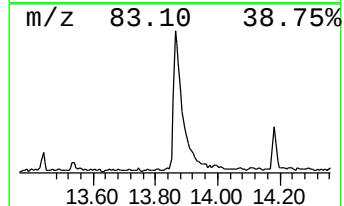
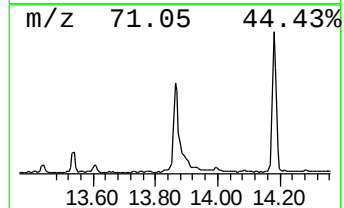
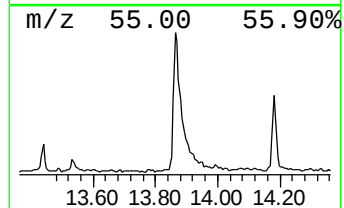
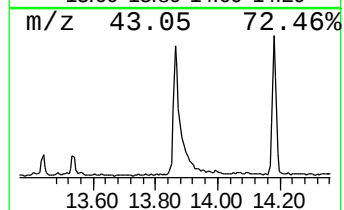
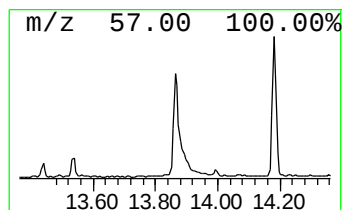
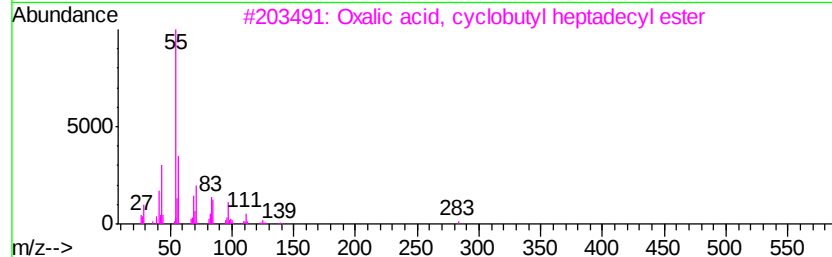
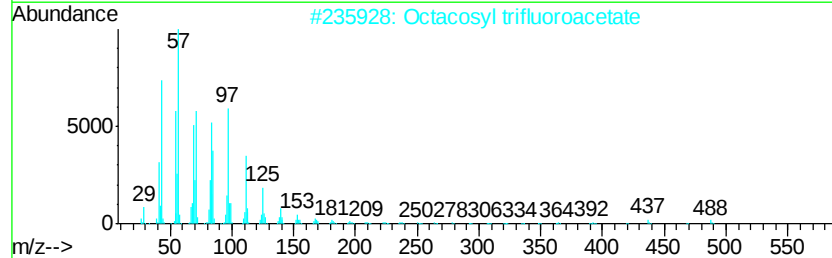
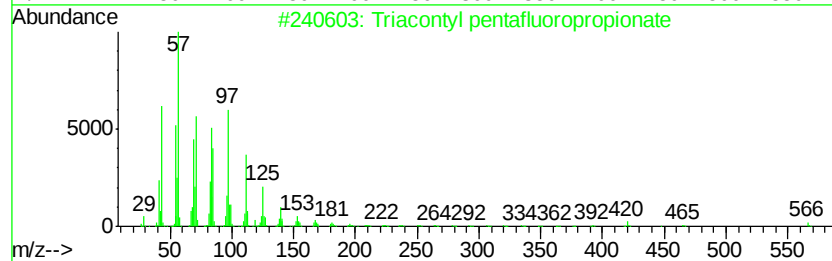
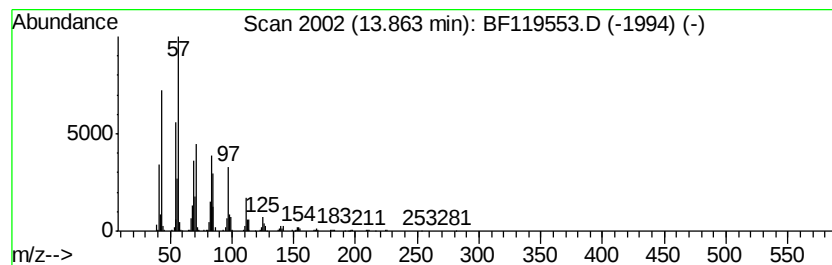
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Triacontyl pentafluoropropi... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	12.43 ng	1350840	Chrysene-d12	14.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	91
2		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
3		Oxalic acid, cyclobutyl heptadec...	382	C23H42O4	1000309-70-7	91
4		Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	91
5		Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
Data File : BF119553.D
Acq On : 27 Mar 2020 7:48
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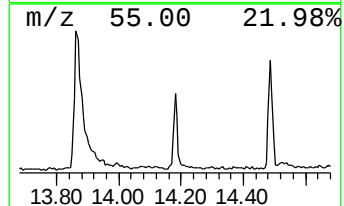
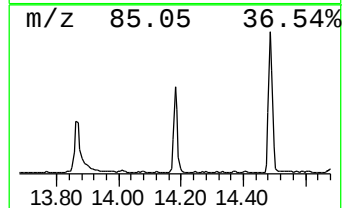
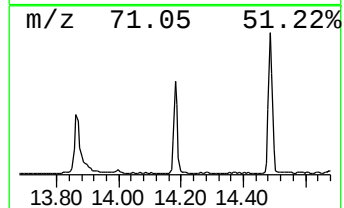
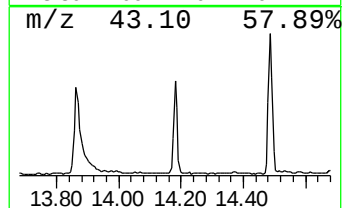
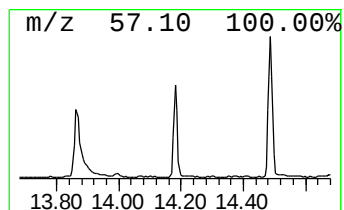
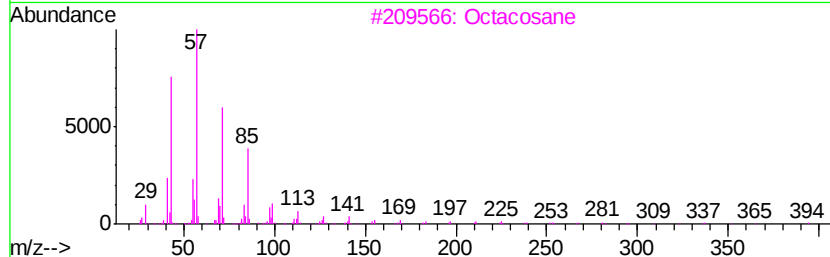
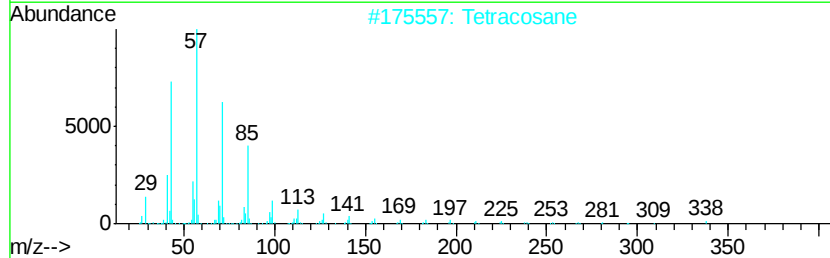
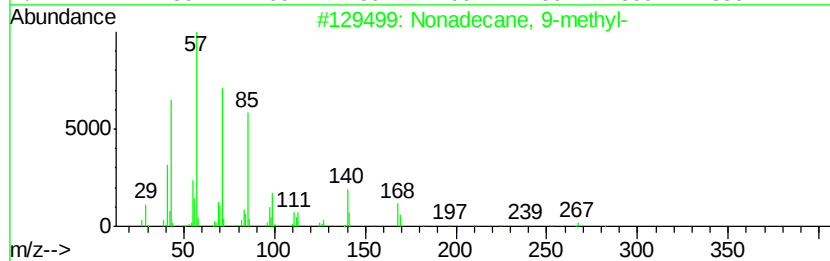
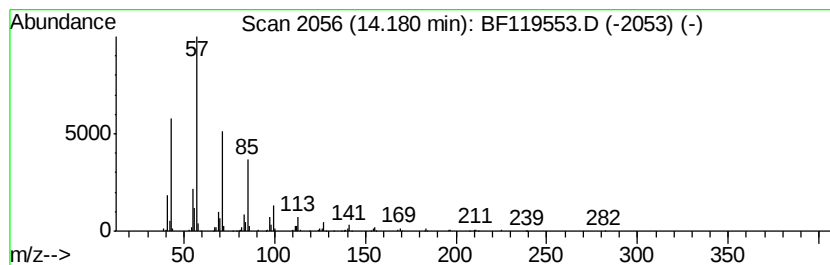
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Nonadecane, 9-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	4.65 ng	505042	Chrysene-d12	14.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Octacosane	394	C28H58	000630-02-4	90
4			Nonadecane, 4-methyl-	282	C20H42	025117-27-5	90
5			Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
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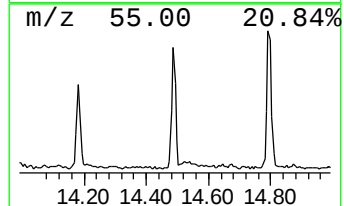
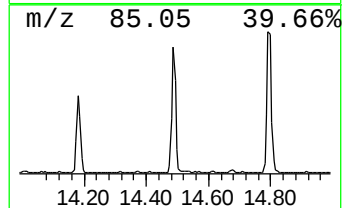
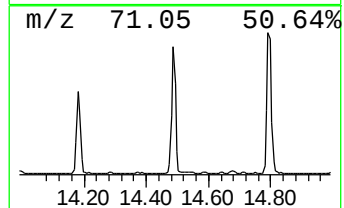
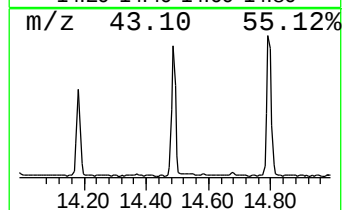
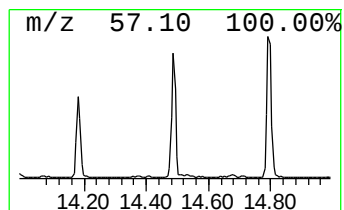
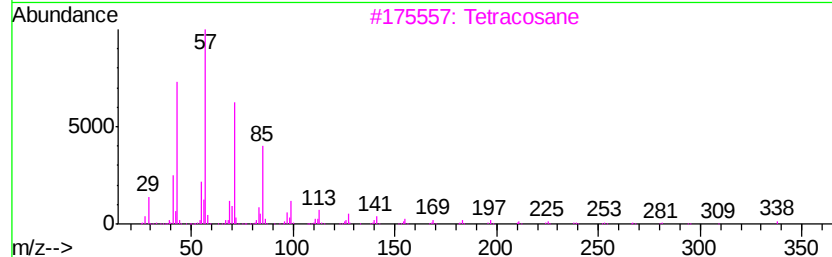
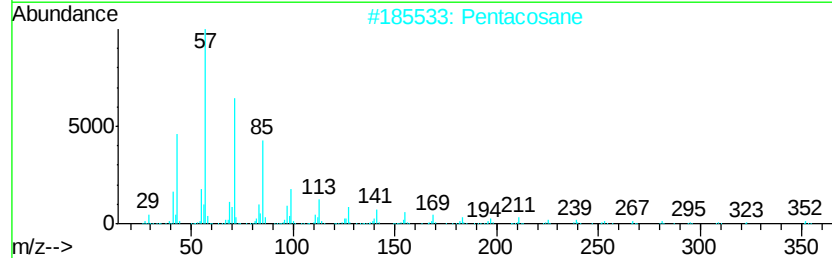
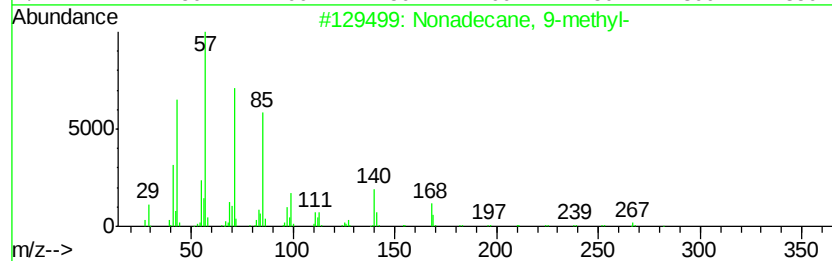
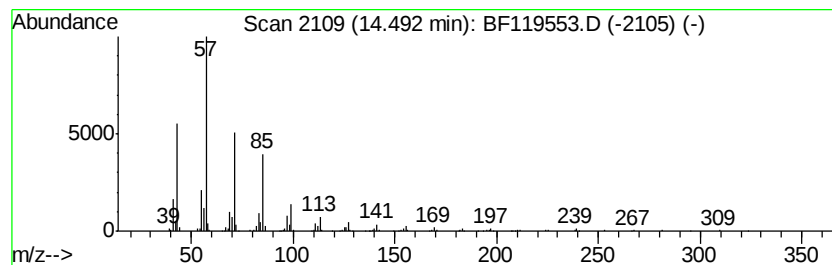
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Pentacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	7.53 ng	818583	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	95
2		Pentacosane	352	C25H52	000629-99-2	91
3		Tetracosane	338	C24H50	000646-31-1	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\BF032620\
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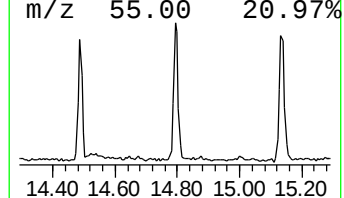
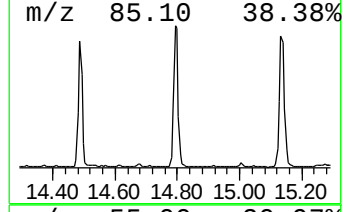
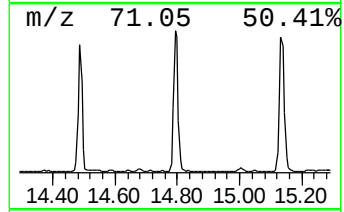
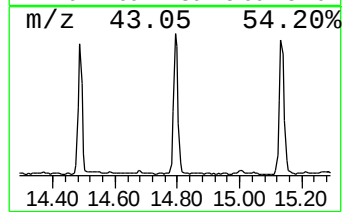
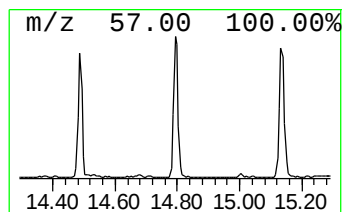
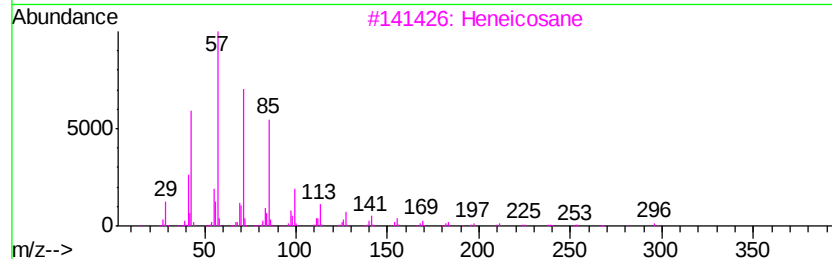
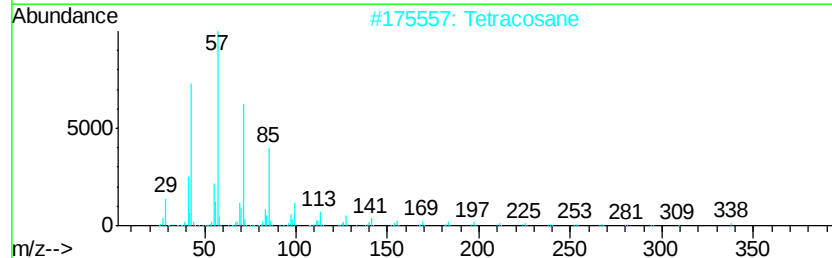
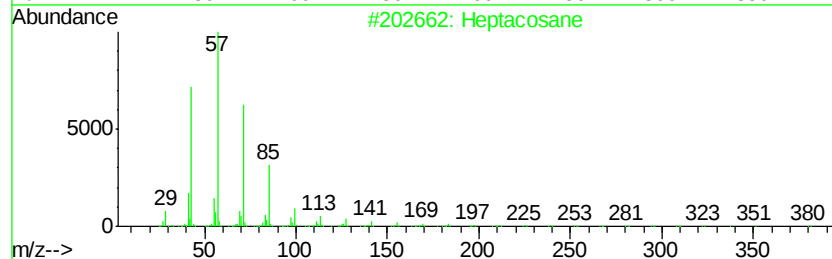
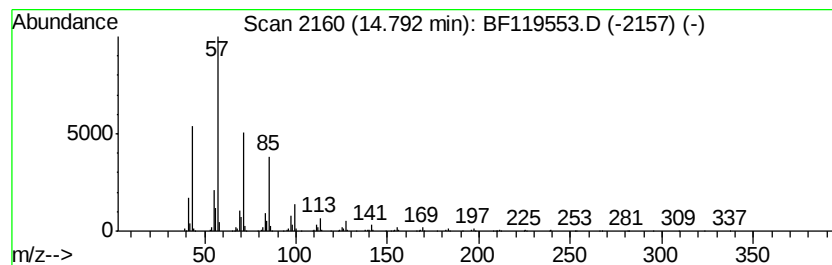
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 12 Heptacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	11.11 ng	1067770	Perylene-d12	15.47

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Tetracosane		338	C24H50	000646-31-1	91
3	Heneicosane		296	C21H44	000629-94-7	90
4	Pentacosane		352	C25H52	000629-99-2	90
5	Nonadecane		268	C19H40	000629-92-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
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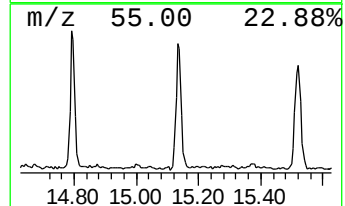
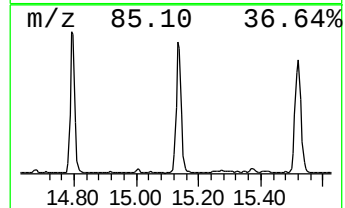
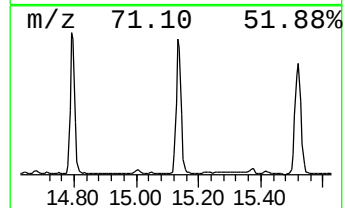
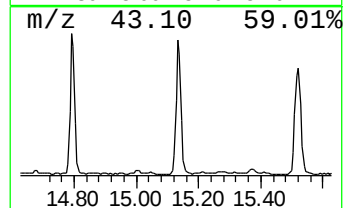
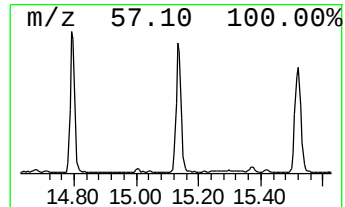
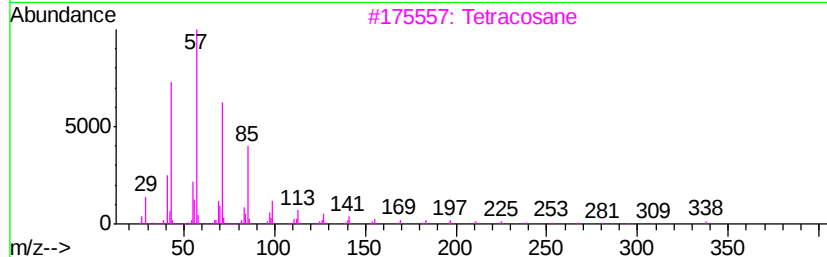
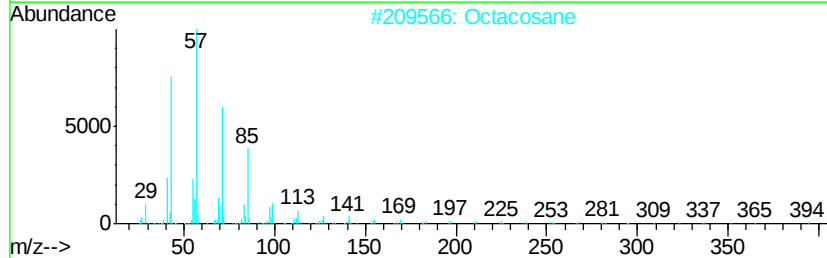
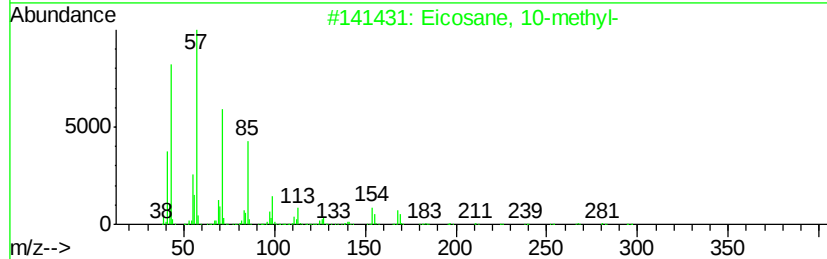
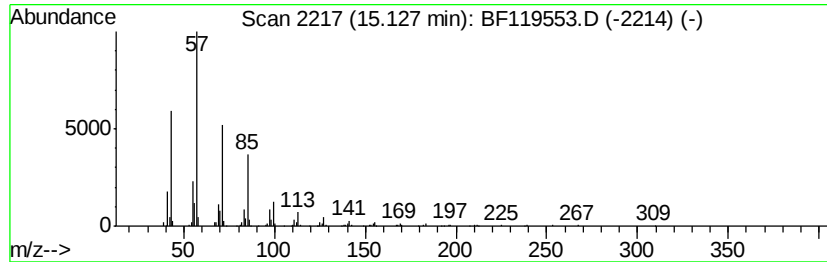
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane, 10-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	11.81 ng	1135020	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
Data File : BF119553.D
Acq On : 27 Mar 2020 7:48
Operator : CG/JU
Sample : L2043-02
Misc :
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Instrument :
BNA_F
ClientSampleId :
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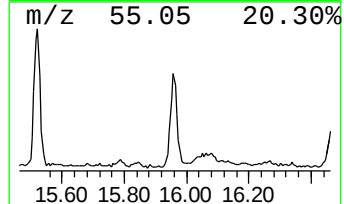
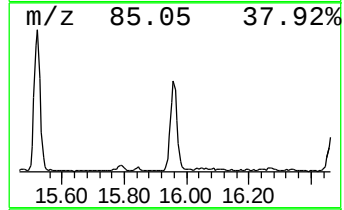
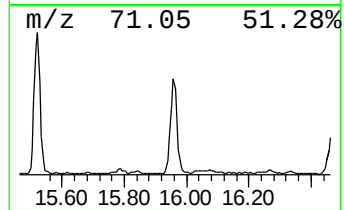
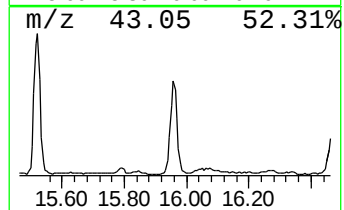
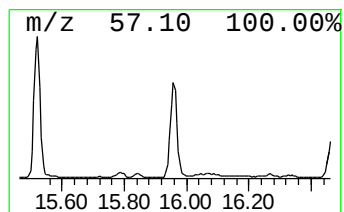
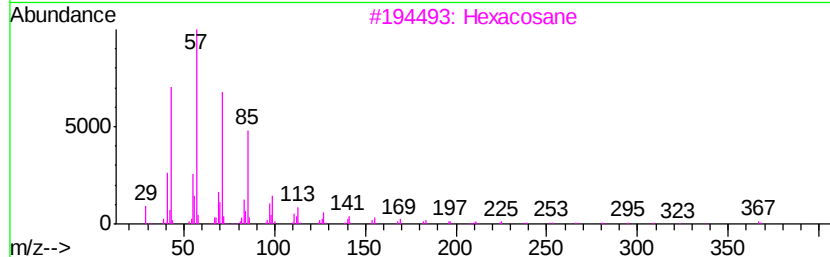
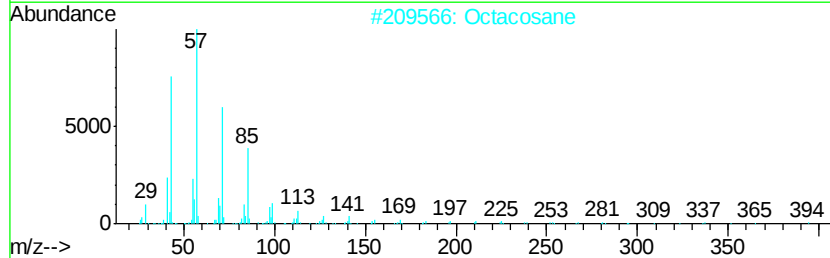
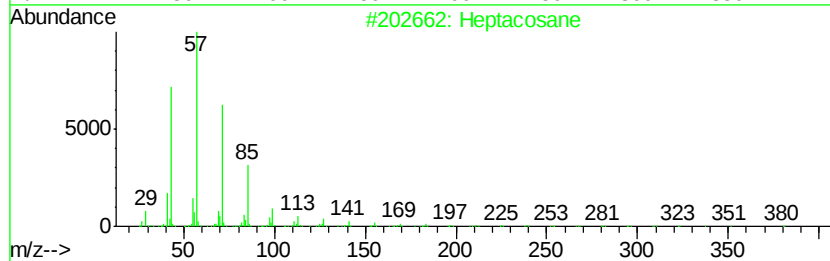
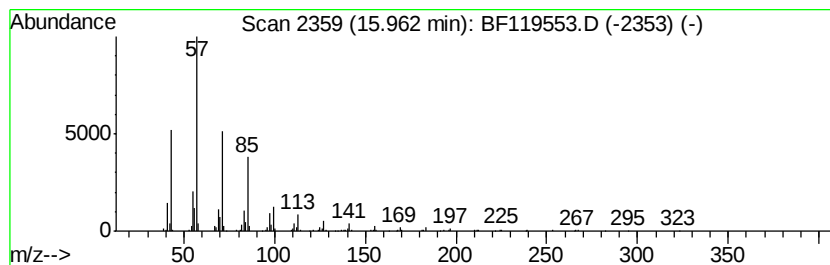
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	8.53 ng	819523	Perylene-d12	15.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	90
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
Data File : BF119553.D
Acq On : 27 Mar 2020 7:48
Operator : CG/JU
Sample : L2043-02
Misc :
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Instrument :
BNA_F
ClientSampleId :
TP-2A

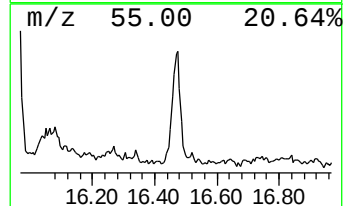
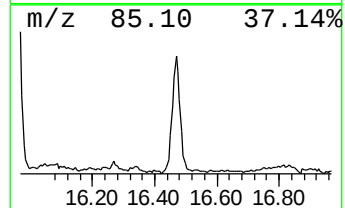
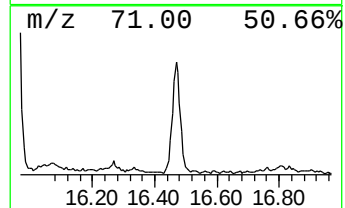
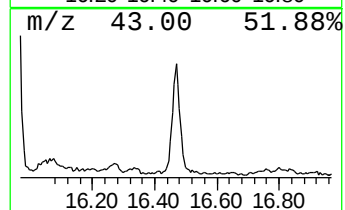
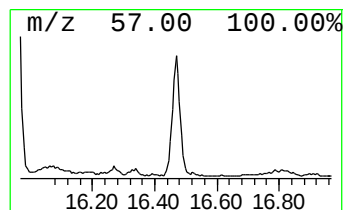
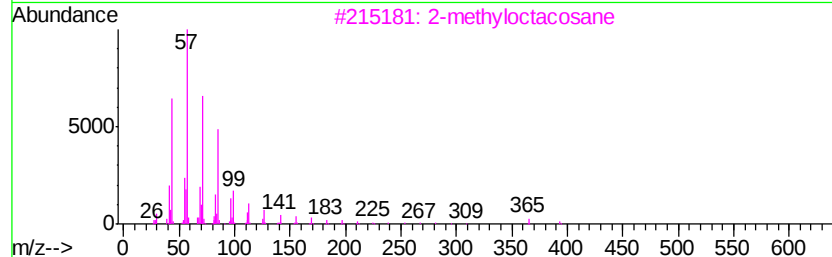
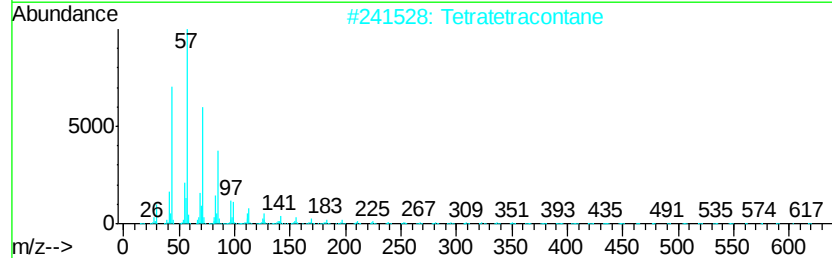
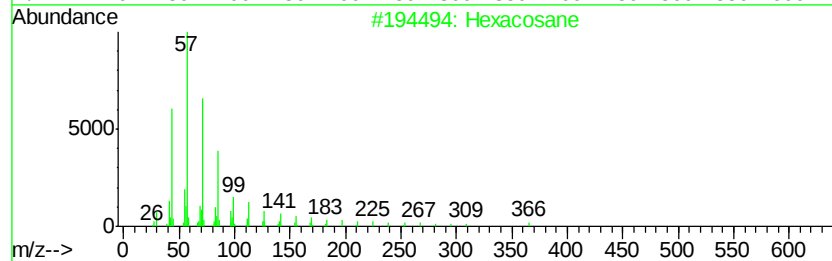
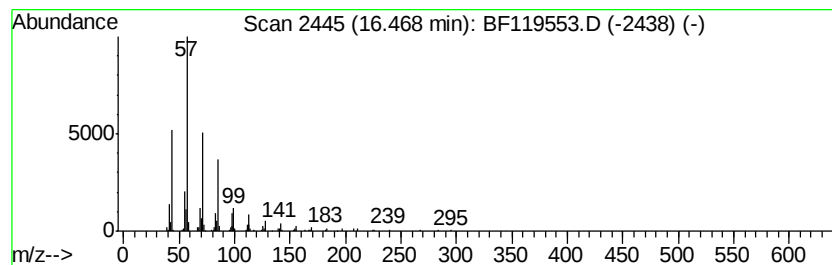
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	4.76 ng	457613	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Tetratetracontane	619	C44H90	007098-22-8	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Pentacosane	352	C25H52	000629-99-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
Data File : BF119553.D
Acq On : 27 Mar 2020 7:48
Operator : CG/JU
Sample : L2043-02
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-2A

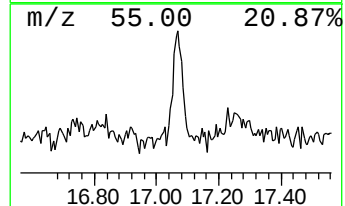
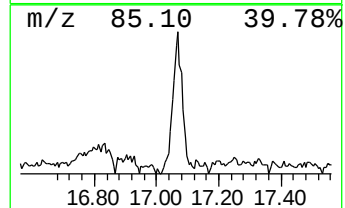
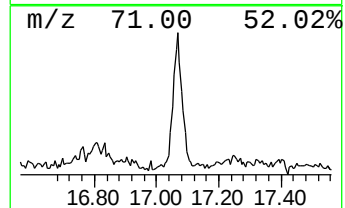
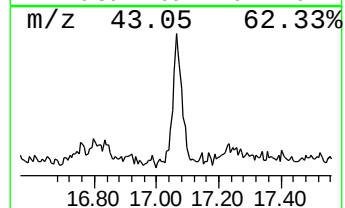
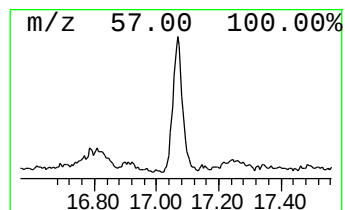
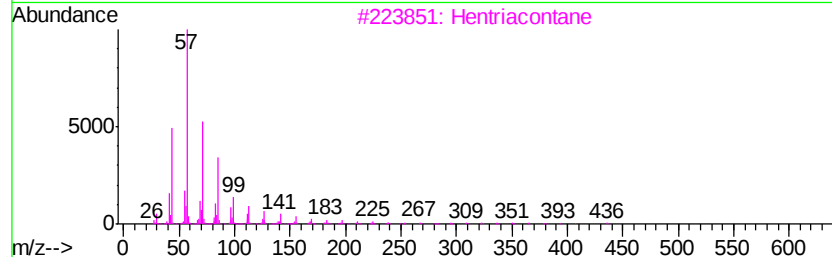
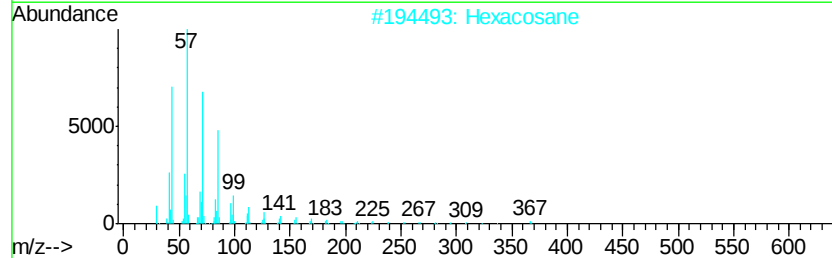
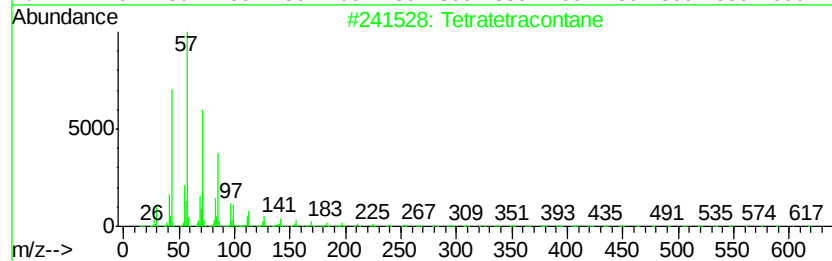
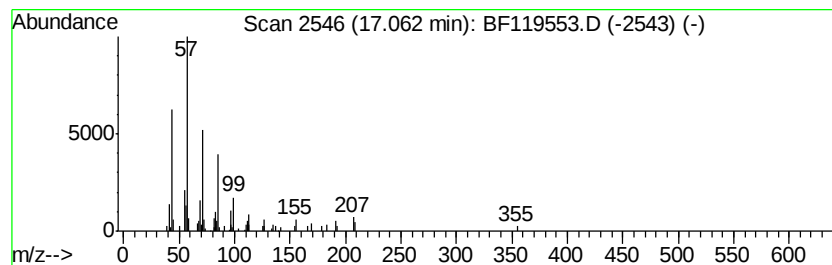
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetratetracontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.06	2.19 ng	210717	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	80
2		Hexacosane	366	C26H54	000630-01-3	80
3		Hentriacontane	437	C31H64	000630-04-6	80
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	72
5		Pentacosane	352	C25H52	000629-99-2	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032620\
 Data File : BF119553.D
 Acq On : 27 Mar 2020 7:48
 Operator : CG/JU
 Sample : L2043-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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...	2.12	6.2	ng	425091	1	6.83	1376460	20.0
Propane, 1,1-dime...	2.23	2.2	ng	150024	1	6.83	1376460	20.0
2-Pentanone, 4-hy...	5.05	13.5	ng	928696	1	6.83	1376460	20.0
2,4,7,9-Tetrameth...	9.32	2.7	ng	294450	3	9.88	2171840	20.0
Benzenesulfonamid...	10.77	2.0	ng	236418	4	11.37	2346800	20.0
n-Hexadecanoic acid	11.90	12.9	ng	1516930	4	11.37	2346800	20.0
Octadecanoic acid	12.69	8.0	ng	939086	4	11.37	2346800	20.0
Triacetyl penta...	13.86	12.4	ng	1350840	5	14.01	2173570	20.0
Nonadecane, 9-met...	14.18	4.7	ng	505042	5	14.01	2173570	20.0
Pentacosane	14.49	7.5	ng	818583	5	14.01	2173570	20.0
Heptacosane	14.79	11.1	ng	1067770	6	15.47	1921970	20.0
Eicosane, 10-methyl-	15.13	11.8	ng	1135020	6	15.47	1921970	20.0
Octacosane	15.96	8.5	ng	819523	6	15.47	1921970	20.0
Hexacosane	16.47	4.8	ng	457613	6	15.47	1921970	20.0
Tetratetracontane	17.06	2.2	ng	210717	6	15.47	1921970	20.0