

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
 Data File : BF130288.D
 Acq On : 16 Sep 2022 13:12
 Operator : CG\JU
 Sample : N4582-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRUSHED-MASONRY-PILE-7

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-BF091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130288.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.198	354	359	366	rBV	312611	378688	15.53%	2.295%
2	4.869	467	473	476	rBV	821407	1034678	42.43%	6.270%
3	5.275	538	542	555	rBV	1002550	897180	36.79%	5.437%
4	6.310	713	718	730	rBV	1508203	1633625	66.99%	9.899%
5	6.675	776	780	783	rBV	744277	561630	23.03%	3.403%
6	7.239	871	876	879	rBV	1412164	1263727	51.82%	7.658%
7	7.957	994	998	1001	rVB	928745	756468	31.02%	4.584%
8	9.033	1176	1181	1184	rBV	2934206	2438776	100.00%	14.778%
9	9.704	1291	1295	1299	rBV	980272	827866	33.95%	5.017%
10	10.486	1425	1428	1434	rVB	114424	110731	4.54%	0.671%
11	11.186	1543	1547	1550	rBV	970111	769037	31.53%	4.660%
12	11.586	1611	1615	1619	rBV2	36632	43252	1.77%	0.262%
13	11.721	1625	1638	1657	rBV	619282	974090	39.94%	5.903%
14	11.986	1680	1683	1688	rVB	58906	62392	2.56%	0.378%
15	12.374	1745	1749	1753	rBV2	58995	94146	3.86%	0.571%
16	12.504	1767	1771	1785	rVB	469008	638969	26.20%	3.872%
17	12.774	1812	1817	1820	rBV	2317519	1963815	80.52%	11.900%
18	13.486	1929	1938	1945	rBV4	57199	177902	7.29%	1.078%
19	13.680	1967	1971	1974	rBV	35158	34963	1.43%	0.212%
20	13.815	1990	1994	1997	rBV	529060	464873	19.06%	2.817%
21	13.998	2021	2025	2029	rBV	38756	42402	1.74%	0.257%
22	14.257	2063	2069	2074	rBV3	58675	127237	5.22%	0.771%
23	14.298	2074	2076	2079	rVB	46541	42323	1.74%	0.256%
24	14.592	2123	2126	2131	rVB	41008	37407	1.53%	0.227%
25	14.874	2169	2174	2177	rBV	50325	89610	3.67%	0.543%
26	15.209	2226	2231	2235	rVB	429346	463943	19.02%	2.811%
27	15.551	2284	2289	2296	rVB2	35273	75861	3.11%	0.460%
28	15.651	2302	2306	2310	rBV6	23475	48638	1.99%	0.295%
29	15.862	2340	2342	2359	rVB7	50369	125800	5.16%	0.762%
30	17.274	2578	2582	2591	rVB4	23308	55195	2.26%	0.334%
31	17.950	2689	2697	2706	rVB3	43814	133316	5.47%	0.808%
32	18.121	2719	2726	2739	rVB3	37192	133711	5.48%	0.810%

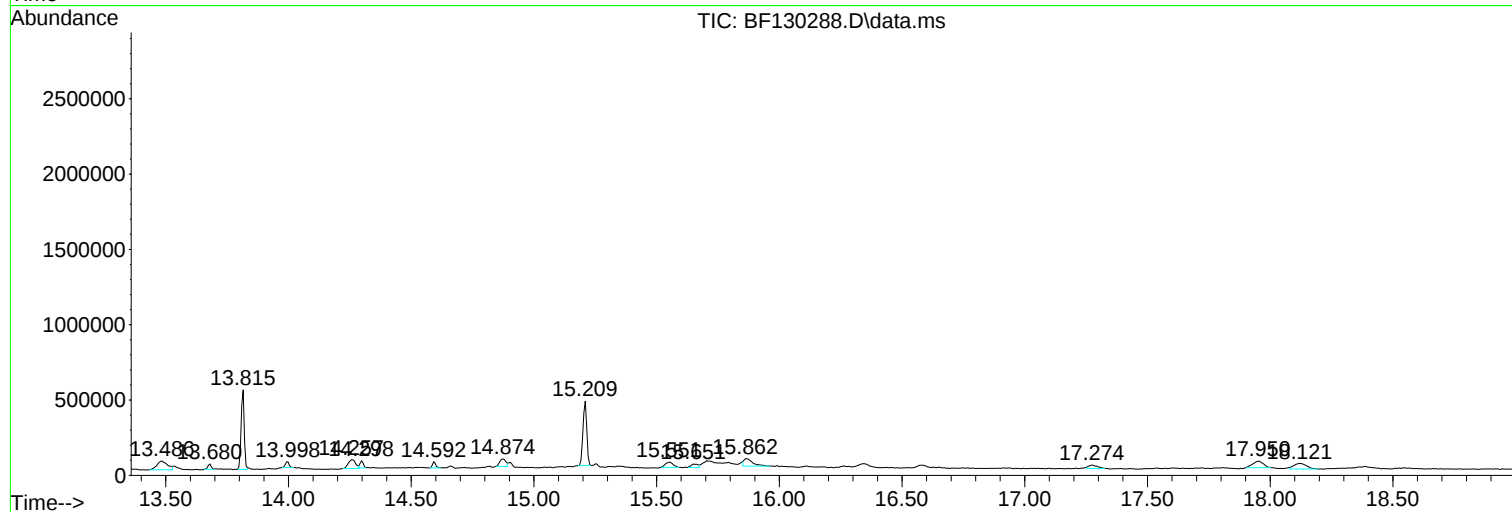
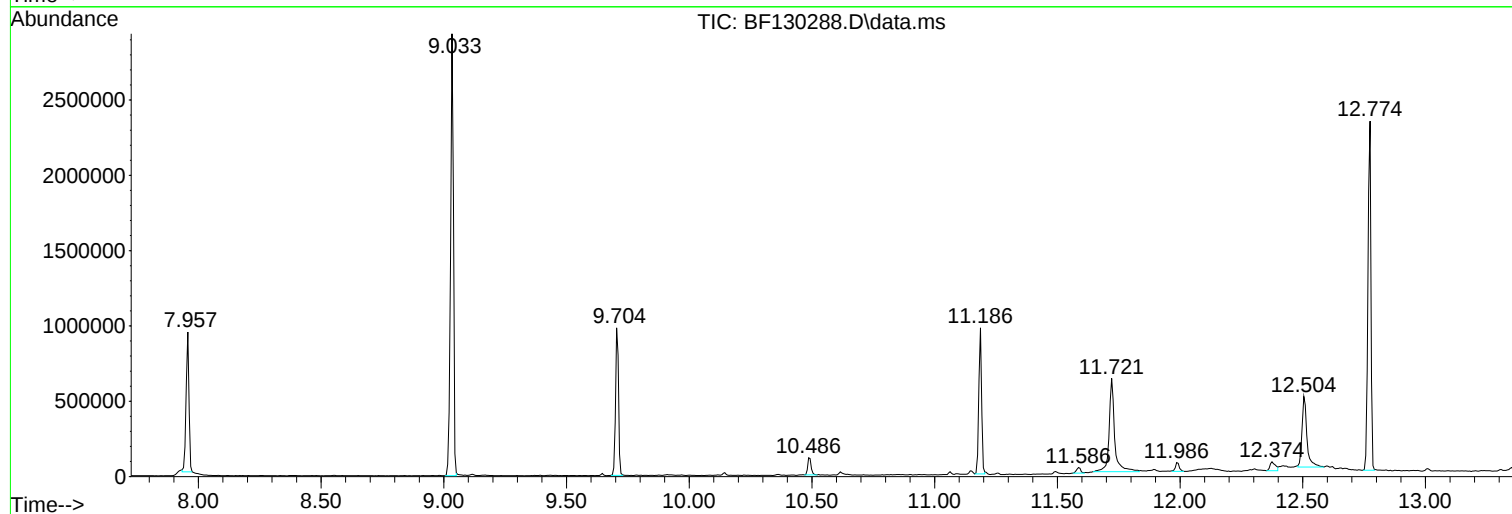
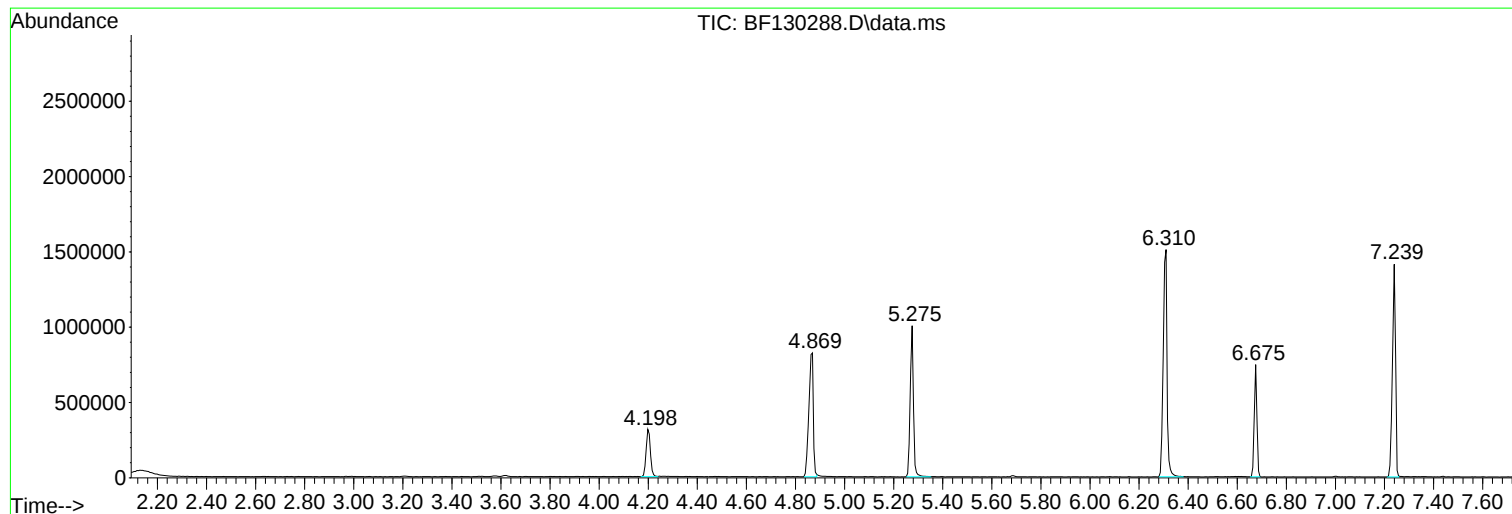
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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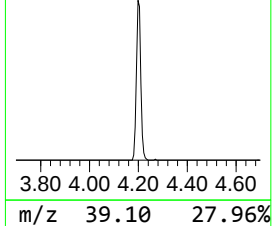
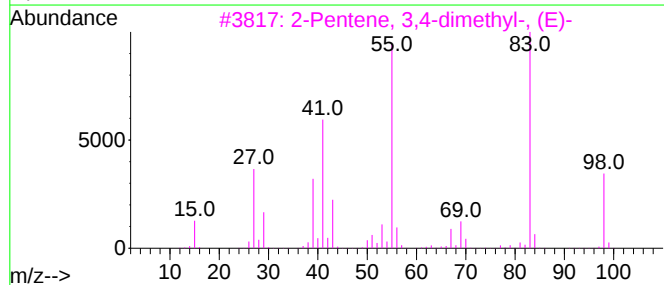
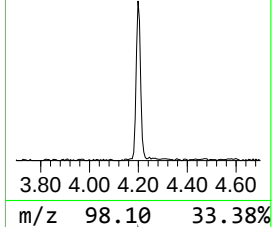
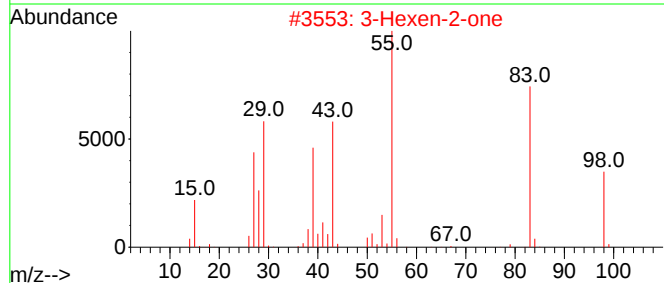
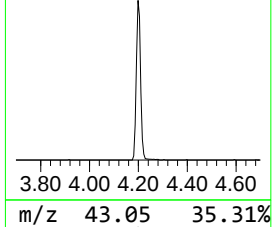
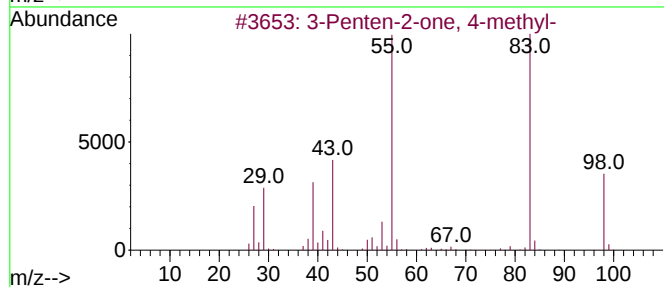
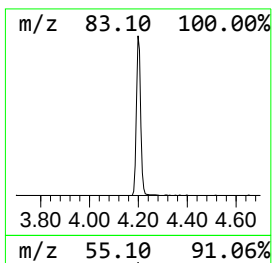
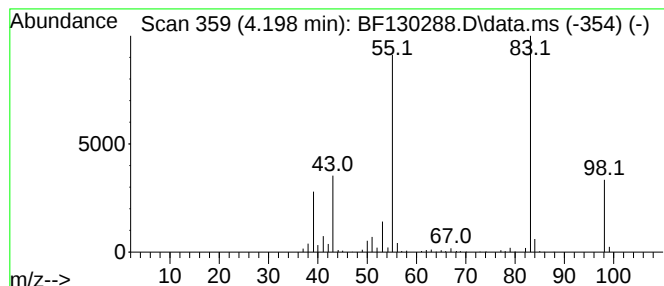
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.198	13.49 ng	378688	1,4-Dichlorobenzene-d4	6.675

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	87
5		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86



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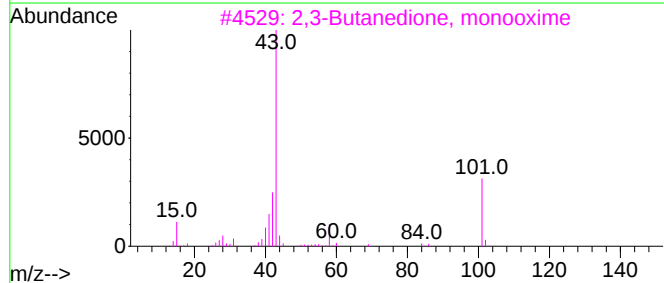
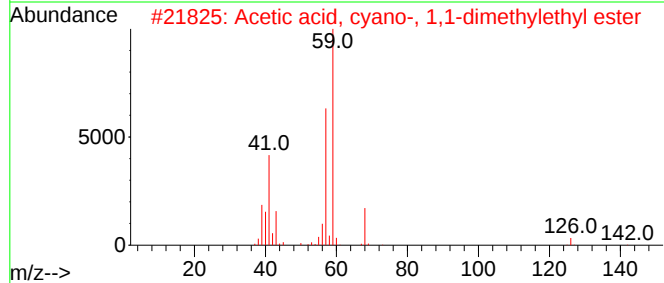
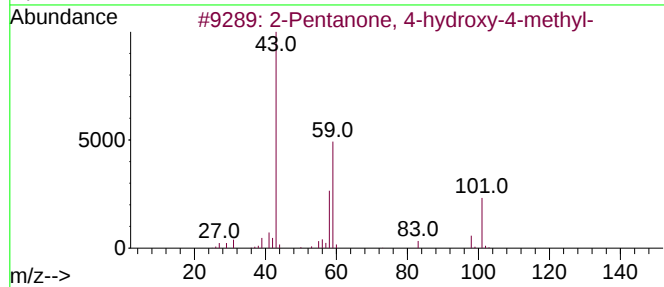
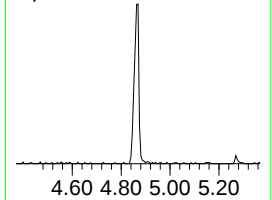
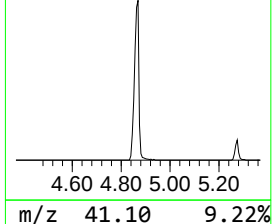
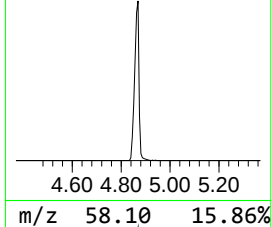
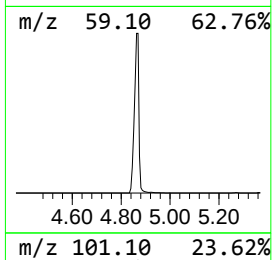
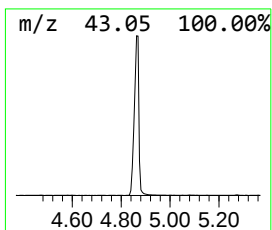
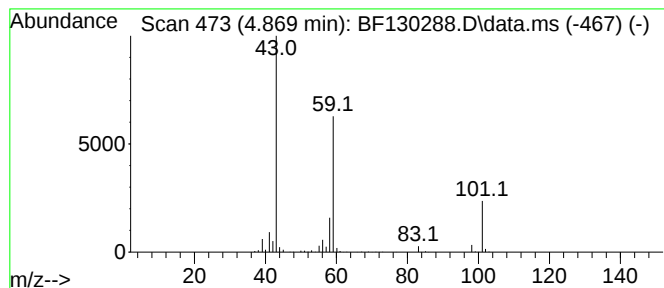
Instrument :
BNA_F
ClientSampleId :
CRUSHED-MASONRY-PILE-7

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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.869	36.85 ng	1034680	1,4-Dichlorobenzene-d4	6.675	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4	Acetone	58	C3H6O	000067-64-1	10
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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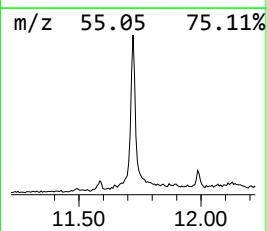
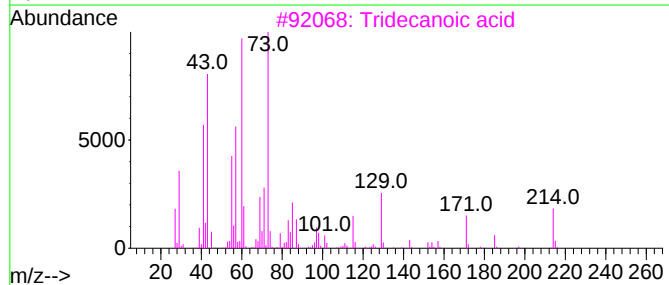
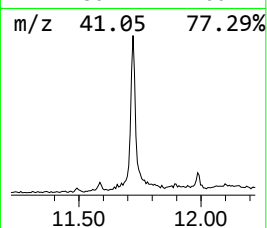
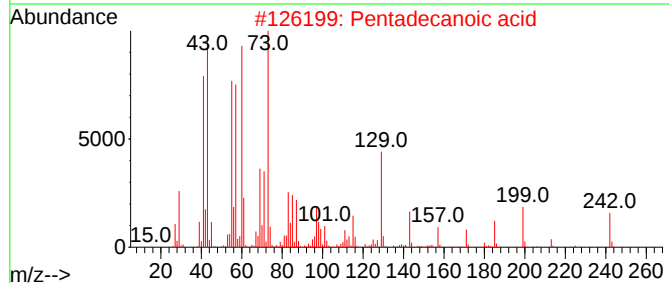
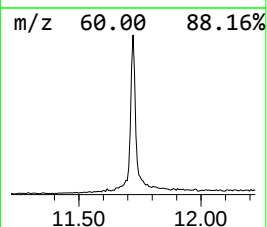
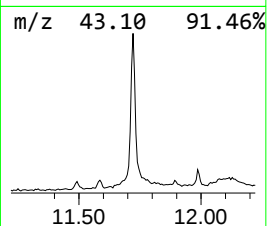
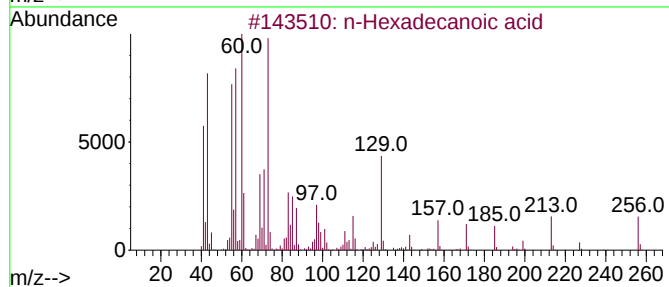
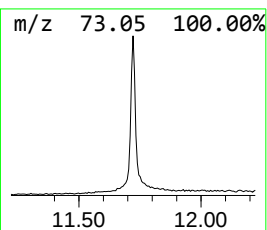
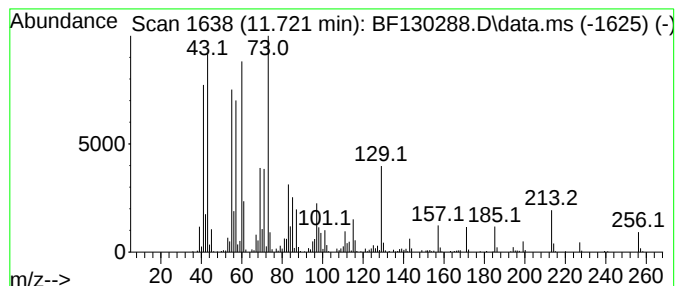
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.722	25.33 ng	974090	Phenanthrene-d10	11.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99		
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93		
3	Tridecanoic acid	214	C13H26O2	000638-53-9	83		
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80		
5	n-Decanoic acid	172	C10H20O2	000334-48-5	76		



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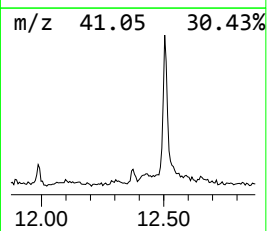
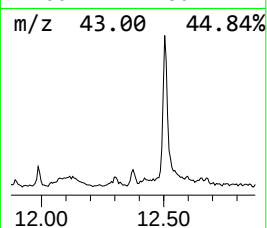
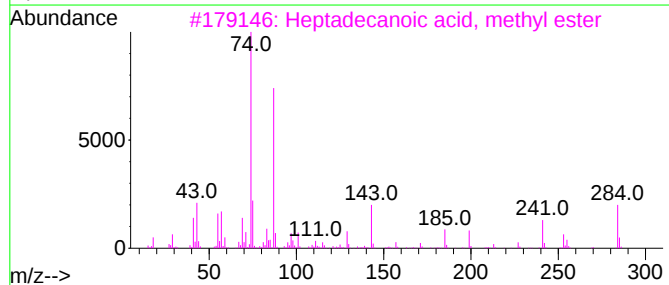
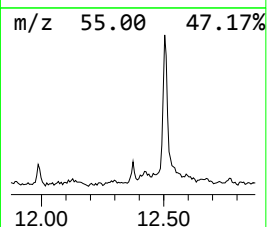
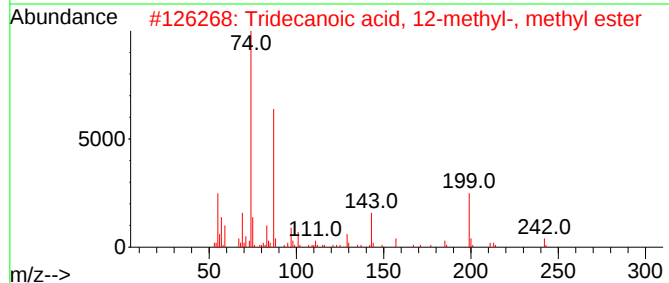
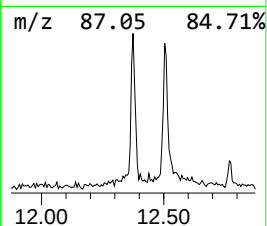
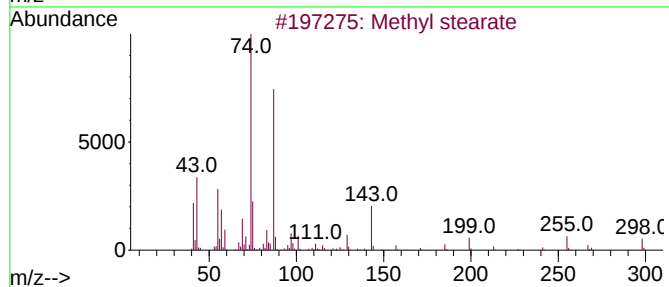
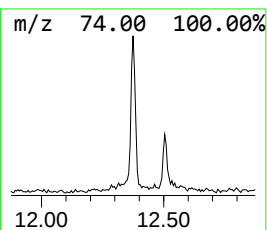
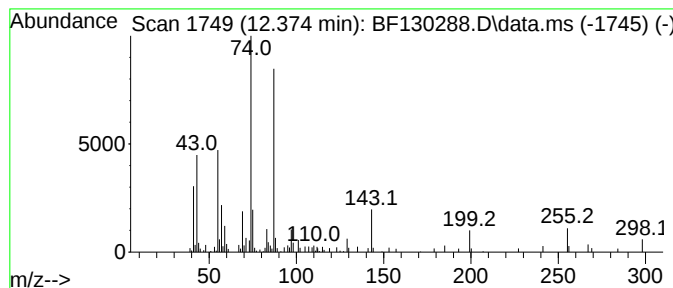
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Peak Number 4 Methyl stearate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.374	2.45 ng	94146	Phenanthrene-d10	11.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	86
3			Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	86
4			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	80
5			Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	80



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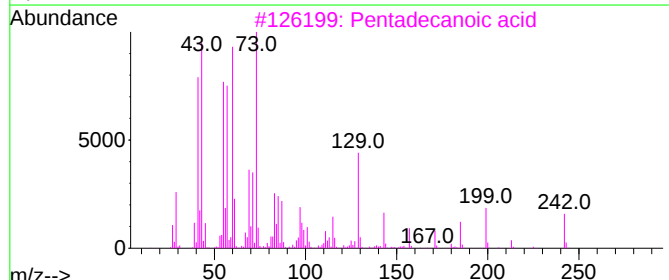
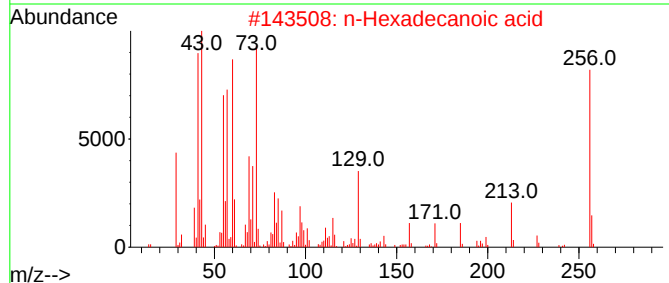
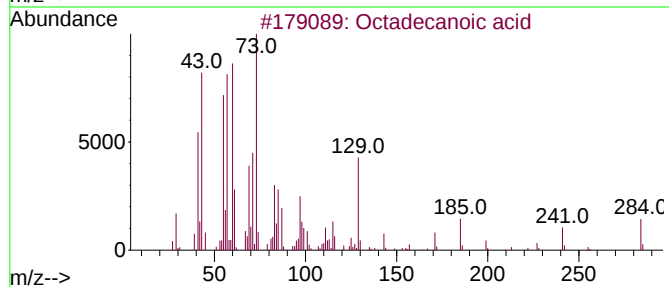
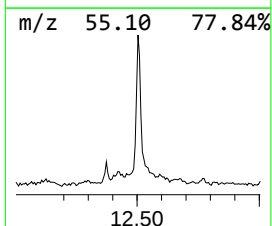
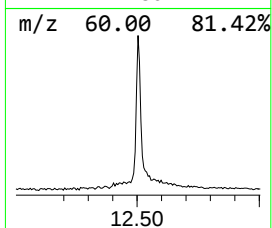
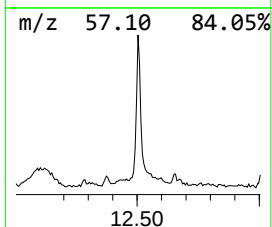
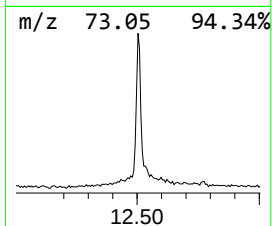
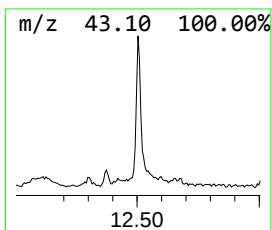
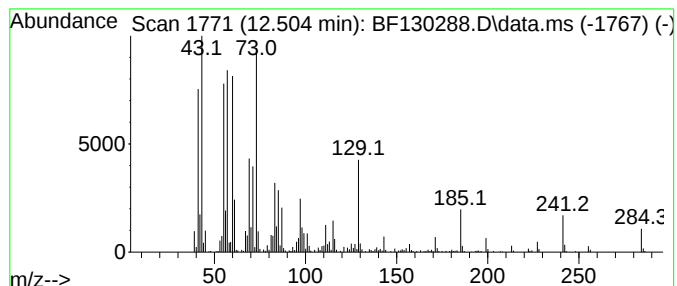
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Peak Number 5 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.504	27.49 ng	638969	Chrysene-d12	13.815

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
5			Tridecanoic acid	214	C13H26O2	000638-53-9	83



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CRUSHED-MASONRY-PILE-7

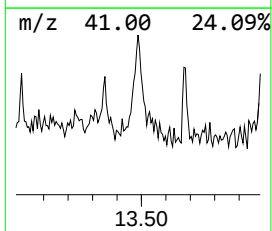
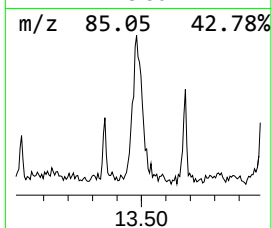
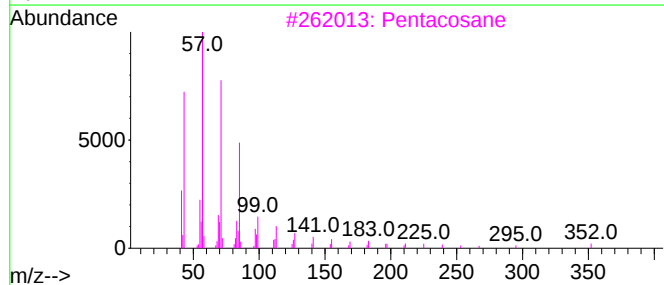
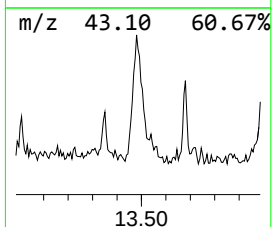
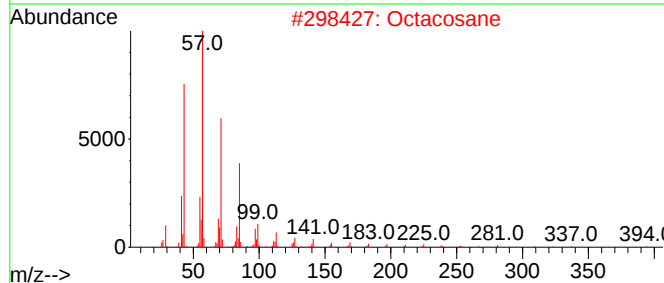
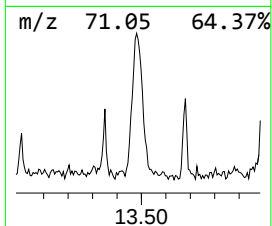
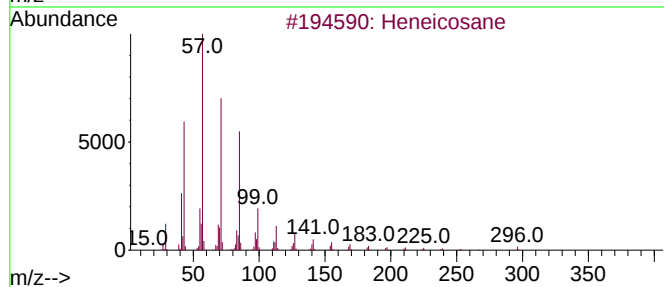
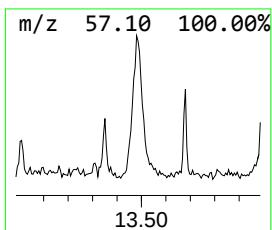
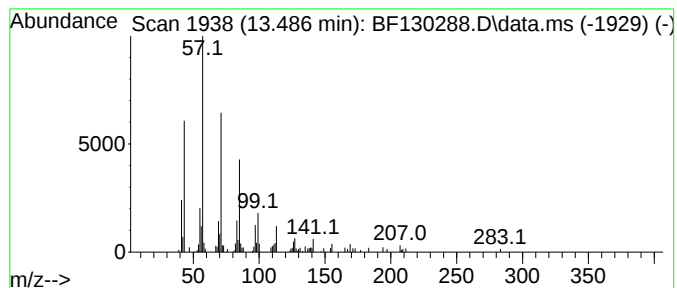
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.486	7.65 ng	177902	Chrysene-d12	13.815

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Pentacosane	352	C25H52	000629-99-2	91
4			Hentriacontane	437	C31H64	000630-04-6	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CRUSHED-MASONRY-PILE-7

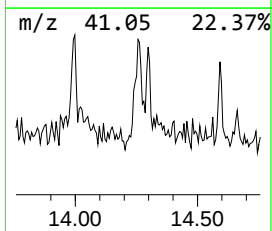
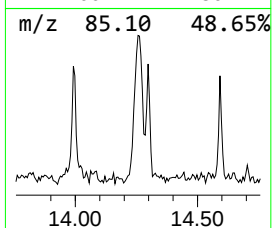
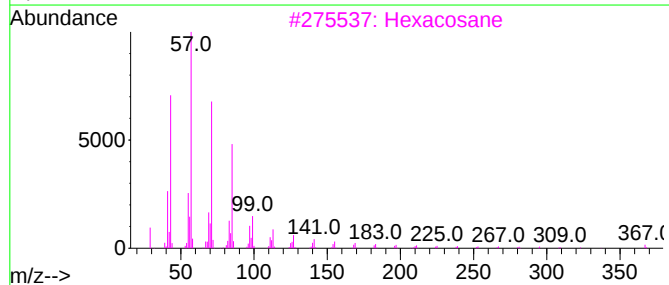
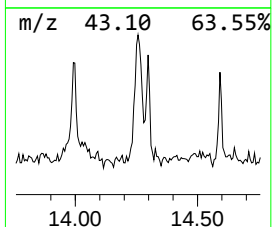
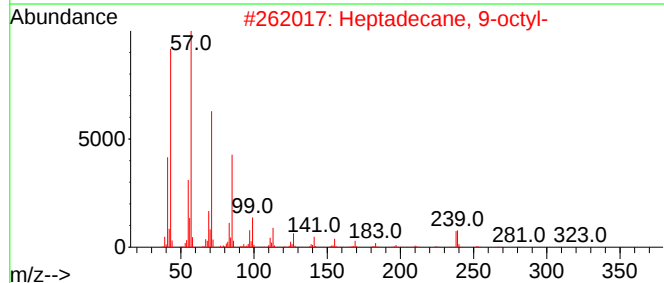
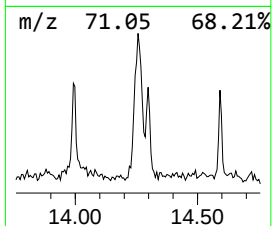
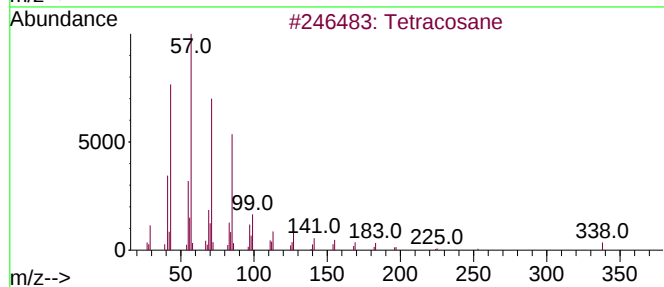
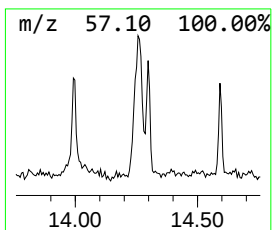
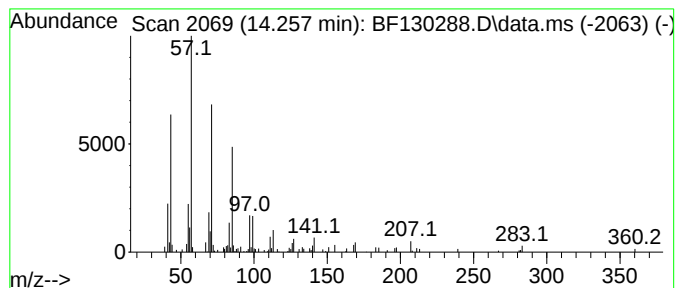
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Tetracosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.257	5.47 ng	127237	Chrysene-d12	13.815

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
3			Hexacosane	366	C26H54	000630-01-3	87
4			Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	86
5			Nonacosane	408	C29H60	000630-03-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CRUSHED-MASONRY-PILE-7

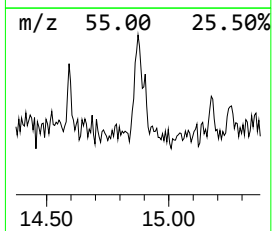
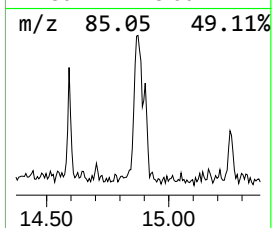
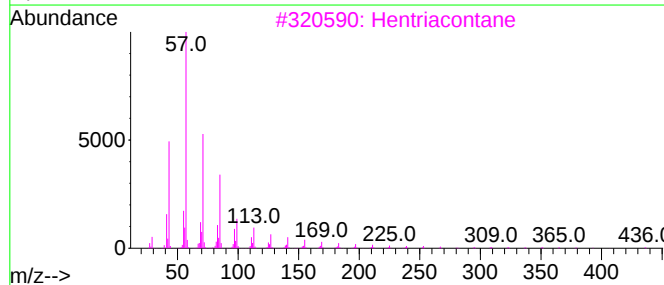
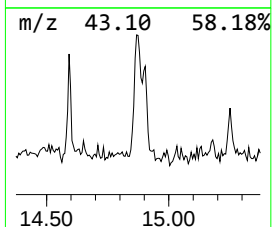
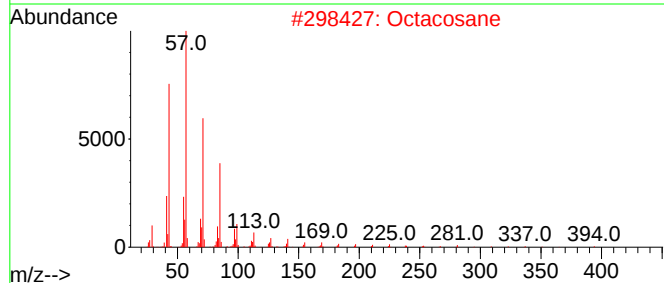
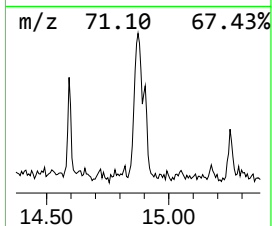
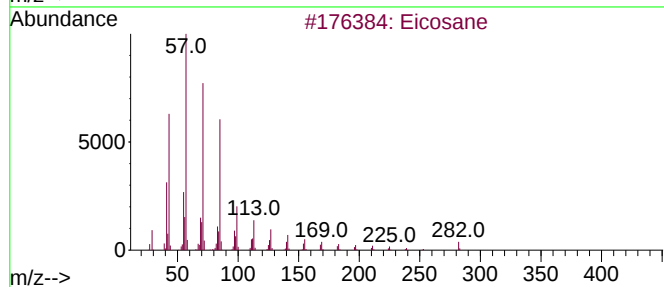
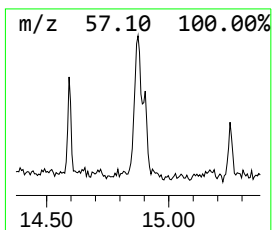
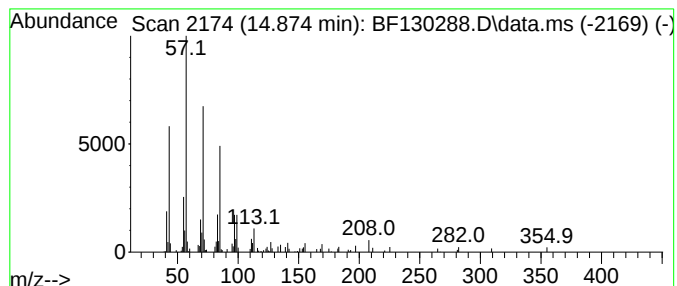
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.874	3.86 ng	89610	Perylene-d12	15.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	90
2			Octacosane	394	C28H58	000630-02-4	87
3			Hentriacontane	437	C31H64	000630-04-6	87
4			Tetratetracontane	619	C44H90	007098-22-8	87
5			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CRUSHED-MASONRY-PILE-7

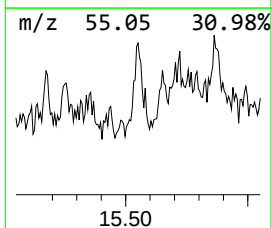
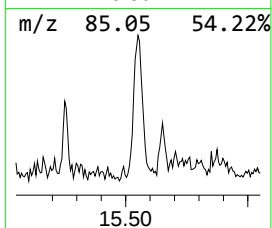
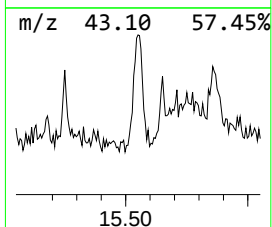
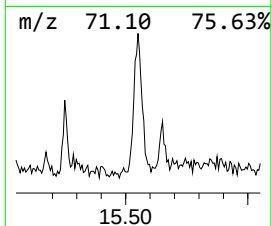
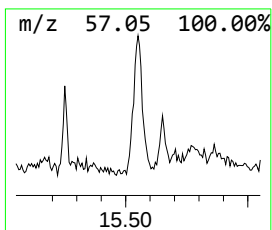
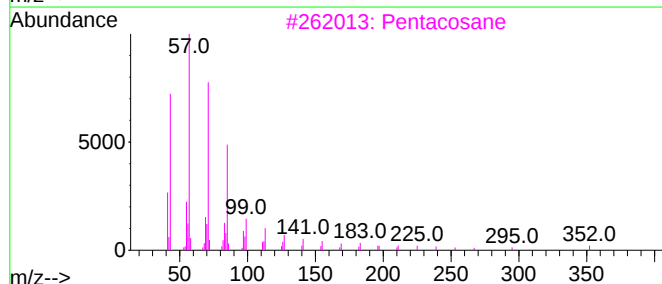
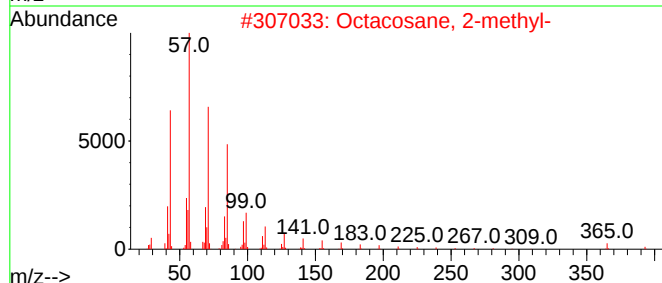
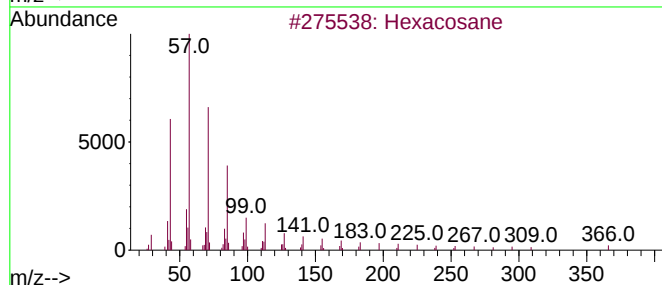
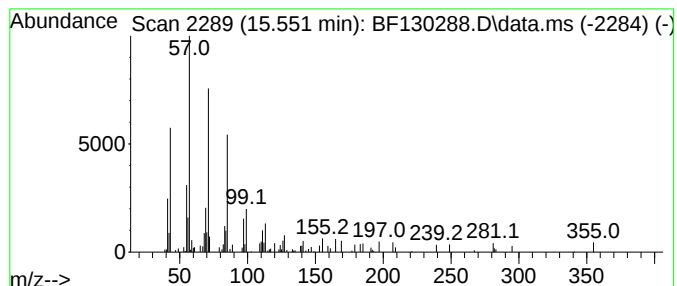
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.551	3.27 ng	75861	Perylene-d12	15.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	87
2			Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
3			Pentacosane	352	C25H52	000629-99-2	87
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	87
5			Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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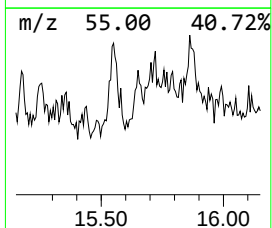
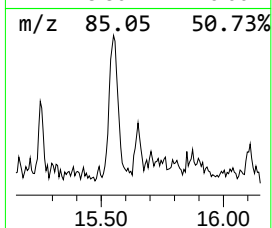
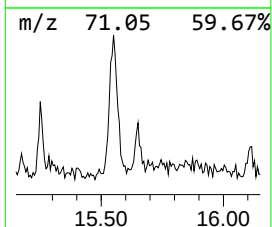
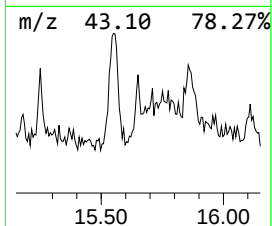
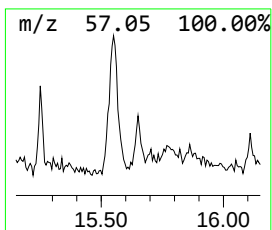
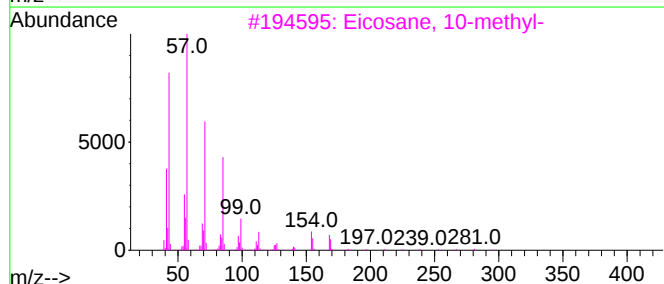
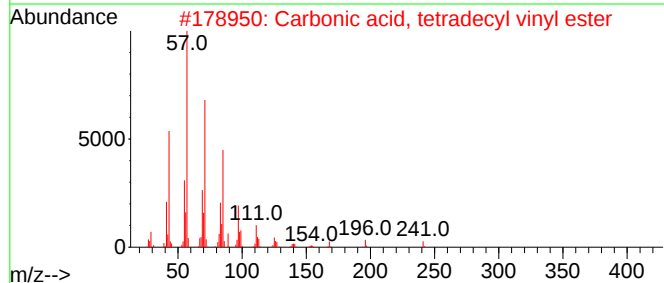
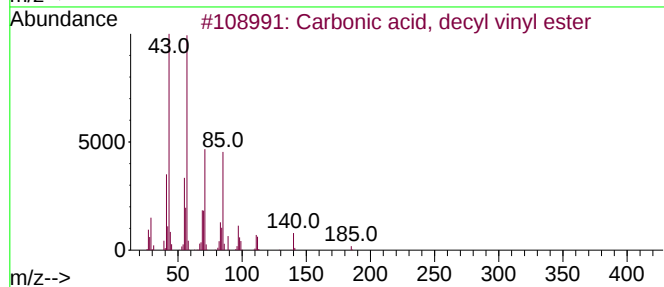
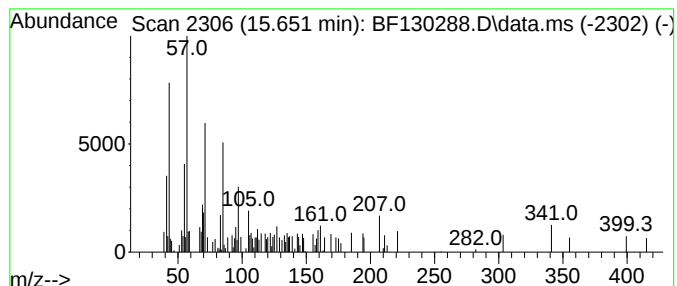
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown15.651 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	2.10 ng	48638	Perylene-d12	15.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, decyl vinyl ester	228	C13H24O3	1000383-25-7	47
2			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	47
3			Eicosane, 10-methyl-	296	C21H44	054833-23-7	43
4			Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	43
5			Succinic acid, tridec-2-yn-1-yl ...	380	C22H36O5	1000390-72-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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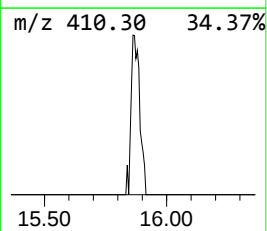
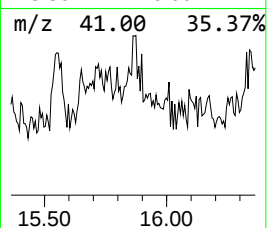
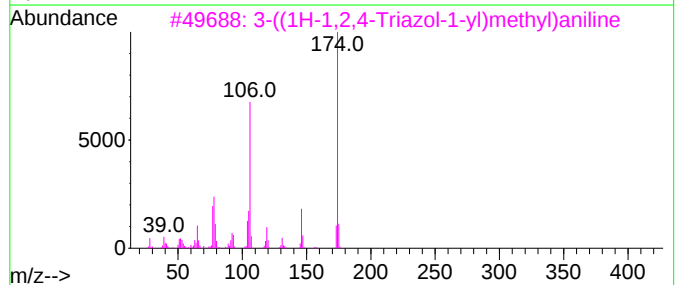
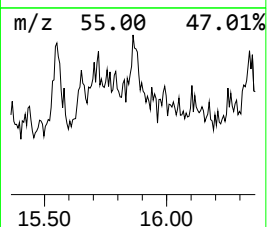
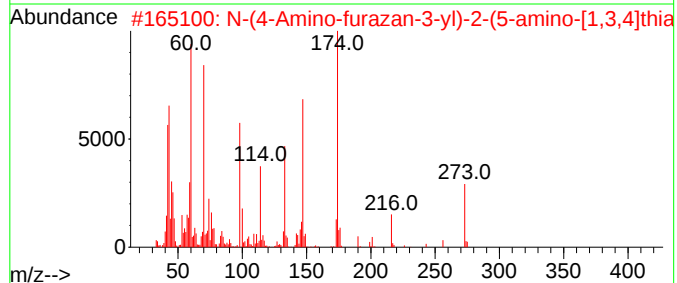
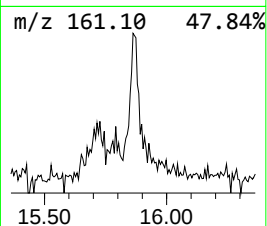
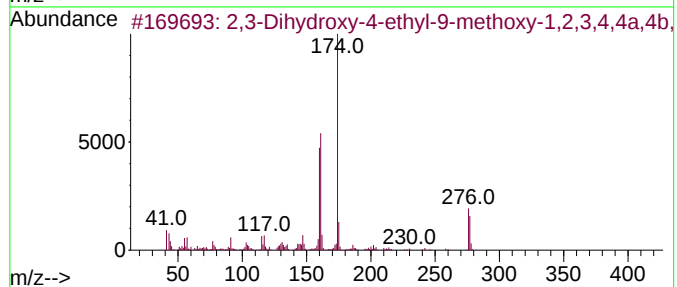
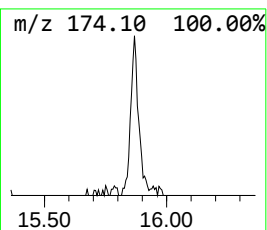
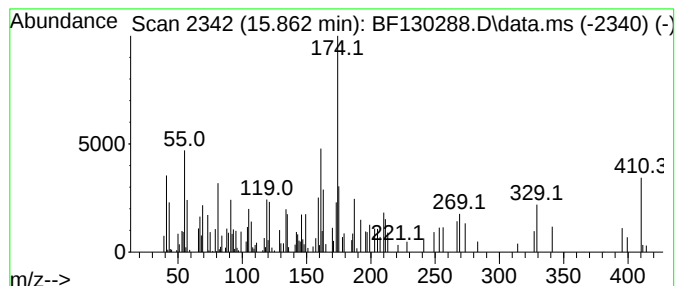
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown15.862 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.862	5.42 ng	125800	Perylene-d12	15.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dihydroxy-4-ethyl-9-methoxy-...	277	C16H23NO3	1000111-32-1	37		
2	N-(4-Amino-furazan-3-yl)-2-(5-am...	273	C6H7N7O2S2	330656-14-9	30		
3	3-((1H-1,2,4-Triazol-1-yl)methyl...	174	C9H10N4	127988-22-1	25		
4	Benzo [f]-1,5-diazabicyclo[3.2.2...	174	C11H14N2	007092-76-4	25		
5	Methyl p-(2-cyanoethoxy)benzoate	205	C11H11NO3	024126-61-2	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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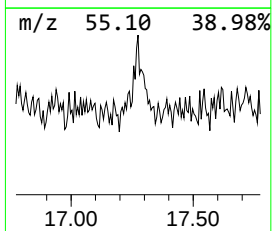
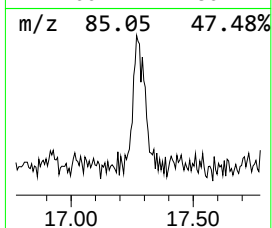
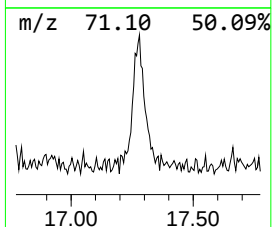
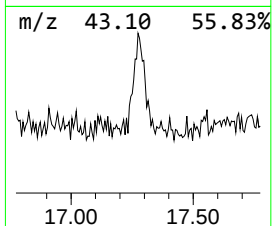
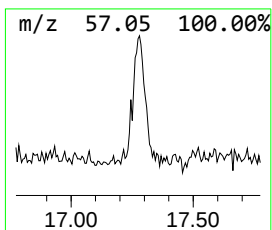
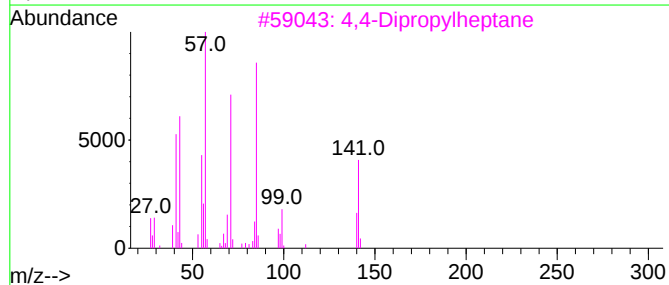
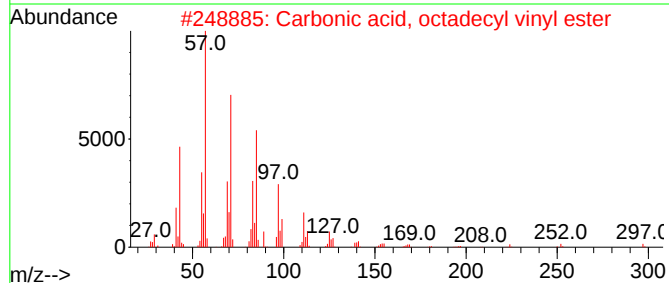
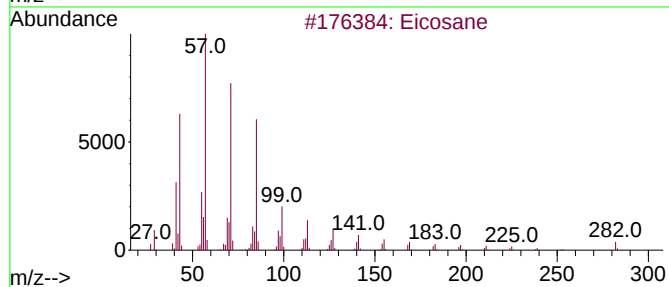
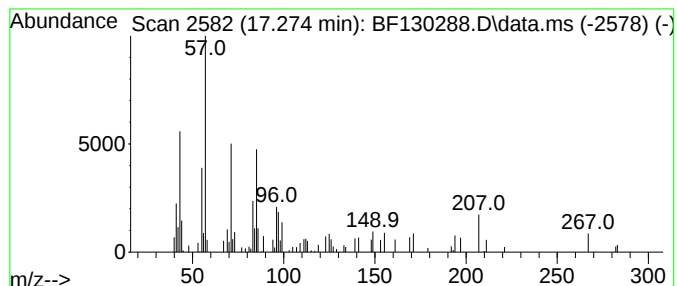
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Carbonic acid, octadecyl vi... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.274	2.38 ng	55195	Perylene-d12	15.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	64
2			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	52
3			4,4-Dipropylheptane	184	C13H28	017312-72-0	50
4			Dodecane, 3-methyl-	184	C13H28	017312-57-1	47
5			Pentacosane	352	C25H52	000629-99-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
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Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

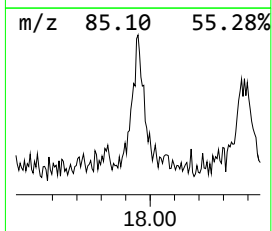
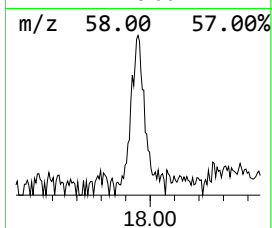
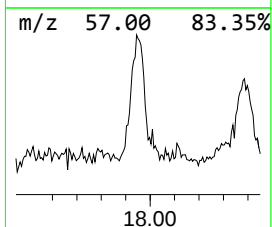
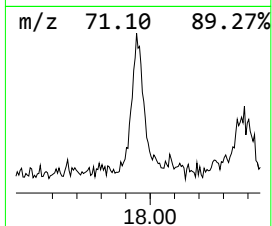
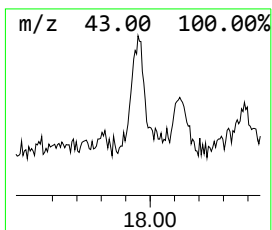
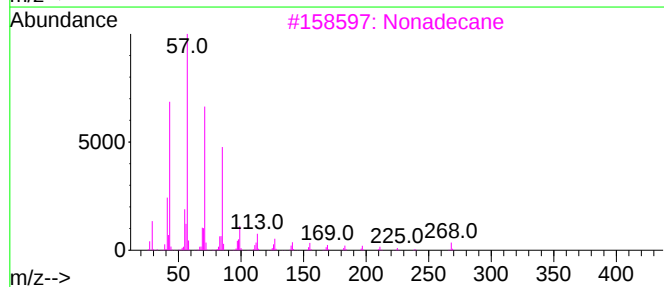
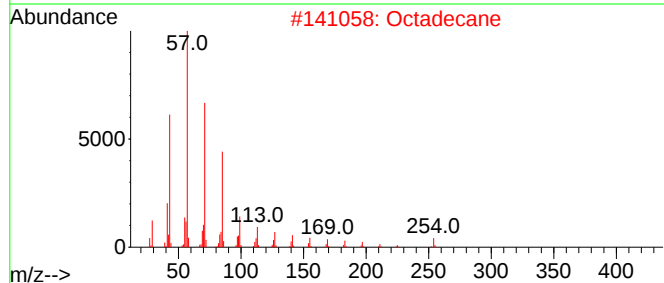
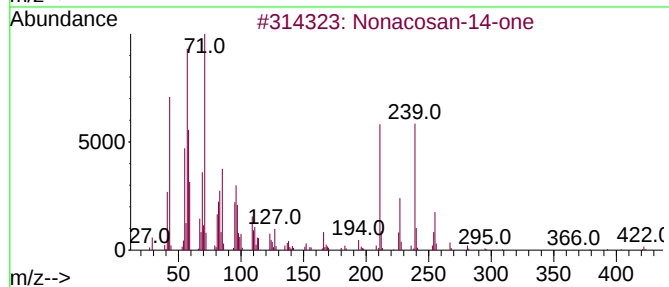
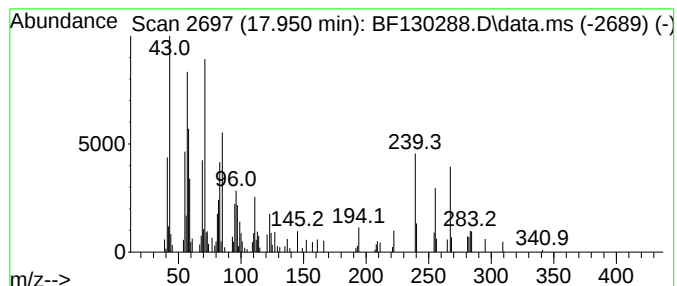
Instrument :
BNA_F
ClientSampleId :
CRUSHED-MASONRY-PILE-7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Nonacosan-14-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.951	5.75 ng	133316	Perylene-d12		15.210	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonacosan-14-one		422	C29H58O	034394-11-1	53
2	Octadecane		254	C18H38	000593-45-3	42
3	Nonadecane		268	C19H40	000629-92-5	42
4	Eicosane		282	C20H42	000112-95-8	42
5	1-Iodo-2-methylnonane		268	C10H21I	1000101-47-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CRUSHED-MASONRY-PILE-7

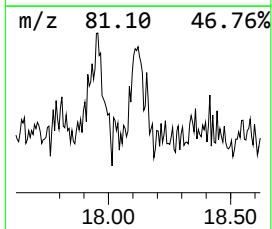
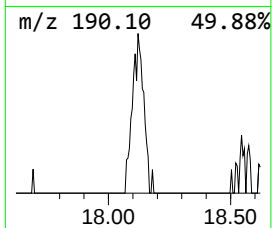
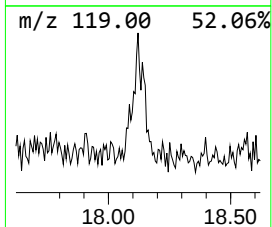
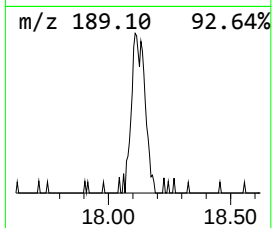
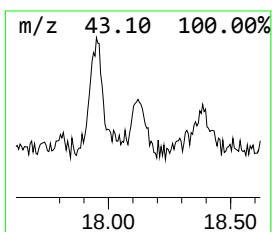
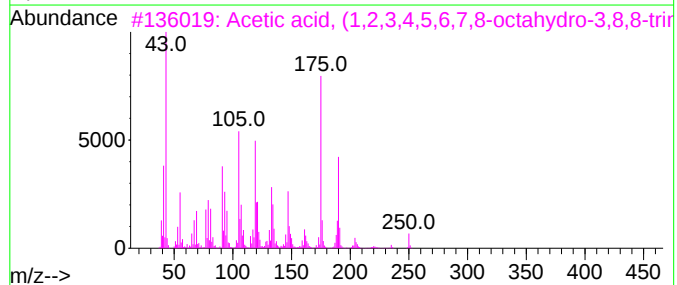
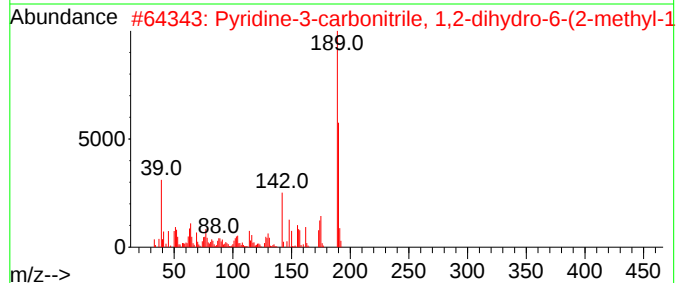
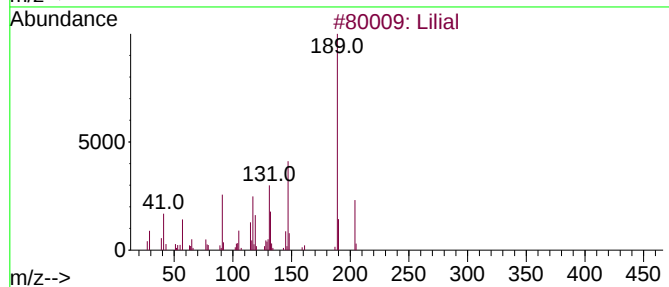
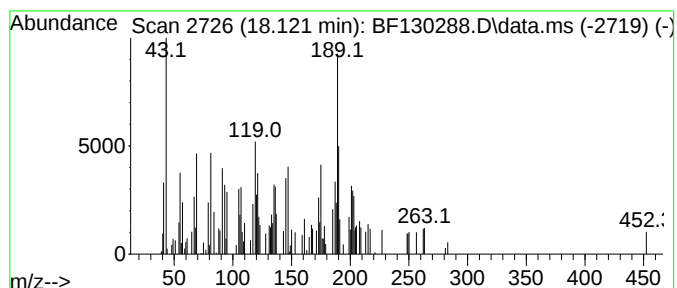
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown18.121 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.121	5.76 ng	133711	Perylene-d12	15.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lilial	204	C14H20O	000080-54-6	35
2			Pyridine-3-carbonitrile, 1,2-dih...	190	C10H10N2S	202270-50-6	27
3			Acetic acid, (1,2,3,4,5,6,7,8-oc...	250	C16H26O2	314773-27-8	25
4			2-Isopropenyl-4a,8-dimethyl-1,2...	204	C15H24	207297-57-2	25
5			.alpha.-Cyano-4-hydroxycinnamic ...	189	C10H7NO3	028166-41-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
Data File : BF130288.D
Acq On : 16 Sep 2022 13:12
Operator : CG\JU
Sample : N4582-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CRUSHED-MASONRY-PILE-7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one,...	4.198	13.5	ng	378688	1	6.675	561630	20.0
2-Pentanone, 4-...	4.869	36.9	ng	1034680	1	6.675	561630	20.0
n-Hexadecanoic ...	11.722	25.3	ng	974090	4	11.186	769037	20.0
Methyl stearate	12.374	2.5	ng	94146	4	11.186	769037	20.0
Octadecanoic acid	12.504	27.5	ng	638969	5	13.815	464873	20.0
Heneicosane	13.486	7.7	ng	177902	5	13.815	464873	20.0
Tetracosane	14.257	5.5	ng	127237	5	13.815	464873	20.0
Eicosane	14.874	3.9	ng	89610	6	15.210	463943	20.0
Hexacosane	15.551	3.3	ng	75861	6	15.210	463943	20.0
unknown15.651	15.651	2.1	ng	48638	6	15.210	463943	20.0
unknown15.862	15.862	5.4	ng	125800	6	15.210	463943	20.0
Carbonic acid, ...	17.274	2.4	ng	55195	6	15.210	463943	20.0
Nonacosan-14-one	17.951	5.8	ng	133316	6	15.210	463943	20.0
unknown18.121	18.121	5.8	ng	133711	6	15.210	463943	20.0