

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
 Data File : BF133420.D
 Acq On : 17 May 2023 18:49
 Operator : CG\JU
 Sample : 02812-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WG-0516-B5-0-10

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-BF042123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF133420.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.899	474	478	494	rVB	923543	1268193	22.53%	2.745%
2	5.328	547	551	562	rBV	4888237	5119032	90.95%	11.079%
3	5.722	614	618	624	rVB	92006	90303	1.60%	0.195%
4	6.363	722	727	730	rBV	5174403	5163961	91.75%	11.176%
5	6.710	783	786	794	rBV	1596499	1302873	23.15%	2.820%
6	7.281	878	883	886	rBV	3952911	3386407	60.17%	7.329%
7	7.998	1001	1005	1008	rBV	2088664	1754512	31.17%	3.797%
8	9.075	1184	1188	1191	rBV	6490252	5628404	100.00%	12.182%
9	9.351	1231	1235	1243	rBV	149463	143308	2.55%	0.310%
10	9.751	1299	1303	1306	rBV	2258308	1987698	35.32%	4.302%
11	10.463	1420	1424	1427	rBV	243437	192149	3.41%	0.416%
12	10.539	1433	1437	1440	rBV	3772335	3454139	61.37%	7.476%
13	11.233	1551	1555	1558	rBV	2298852	1993756	35.42%	4.315%
14	11.769	1643	1646	1656	rBV3	281158	355323	6.31%	0.769%
15	12.445	1757	1761	1766	rBV	107716	112403	2.00%	0.243%
16	12.669	1793	1799	1802	rBV	93051	114843	2.04%	0.249%
17	12.822	1820	1825	1828	rBV	5360736	4817505	85.59%	10.427%
18	13.057	1858	1865	1870	rBV	169184	218997	3.89%	0.474%
19	13.398	1918	1923	1926	rBV2	96752	98161	1.74%	0.212%
20	13.727	1975	1979	1996	rVB	862151	1386888	24.64%	3.002%
21	13.869	1998	2003	2007	rBV	1261968	1194309	21.22%	2.585%
22	14.045	2030	2033	2040	rBV	101791	141450	2.51%	0.306%
23	14.357	2081	2086	2105	rBV	1384727	2489425	44.23%	5.388%
24	14.645	2132	2135	2138	rBV	93227	94000	1.67%	0.203%
25	14.680	2138	2141	2144	rBV4	49704	75129	1.33%	0.163%
26	14.968	2186	2190	2192	rBV	194621	205965	3.66%	0.446%
27	14.998	2192	2195	2200	rBV	239334	410780	7.30%	0.889%
28	15.292	2240	2245	2249	rBV	1251081	1428751	25.38%	3.092%
29	15.321	2249	2250	2256	rVB3	86525	82662	1.47%	0.179%
30	15.521	2279	2284	2289	rBV5	77737	127843	2.27%	0.277%
31	15.733	2316	2320	2325	rVB	140128	189310	3.36%	0.410%
32	15.798	2326	2331	2335	rBV4	90350	207472	3.69%	0.449%
33	15.845	2336	2339	2347	rVB2	116299	225234	4.00%	0.487%
34	16.198	2396	2399	2414	rVB8	58038	155106	2.76%	0.336%
35	16.857	2508	2511	2517	rBV6	45065	75363	1.34%	0.163%
36	16.921	2518	2522	2524	rBV2	50829	76472	1.36%	0.166%

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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-BF042123.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	17.274	2578	2582	2590	rVB8	126276	253151	4.50%	0.548%
38	17.527	2618	2625	2629	rBV9	64678	182801	3.25%	0.396%

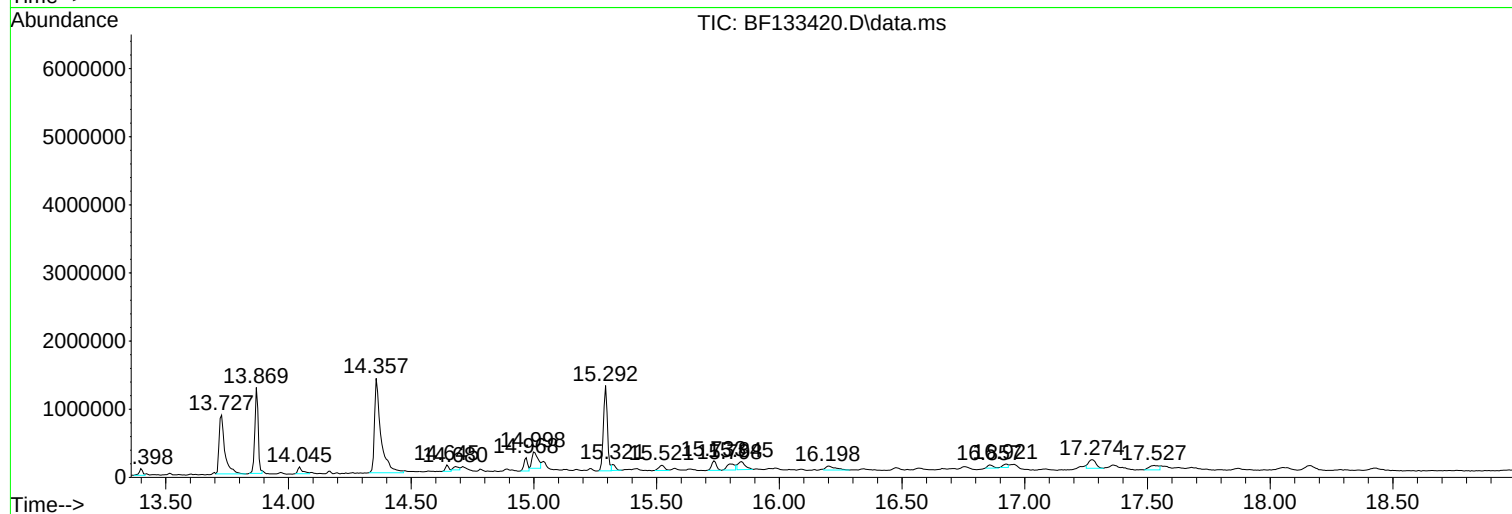
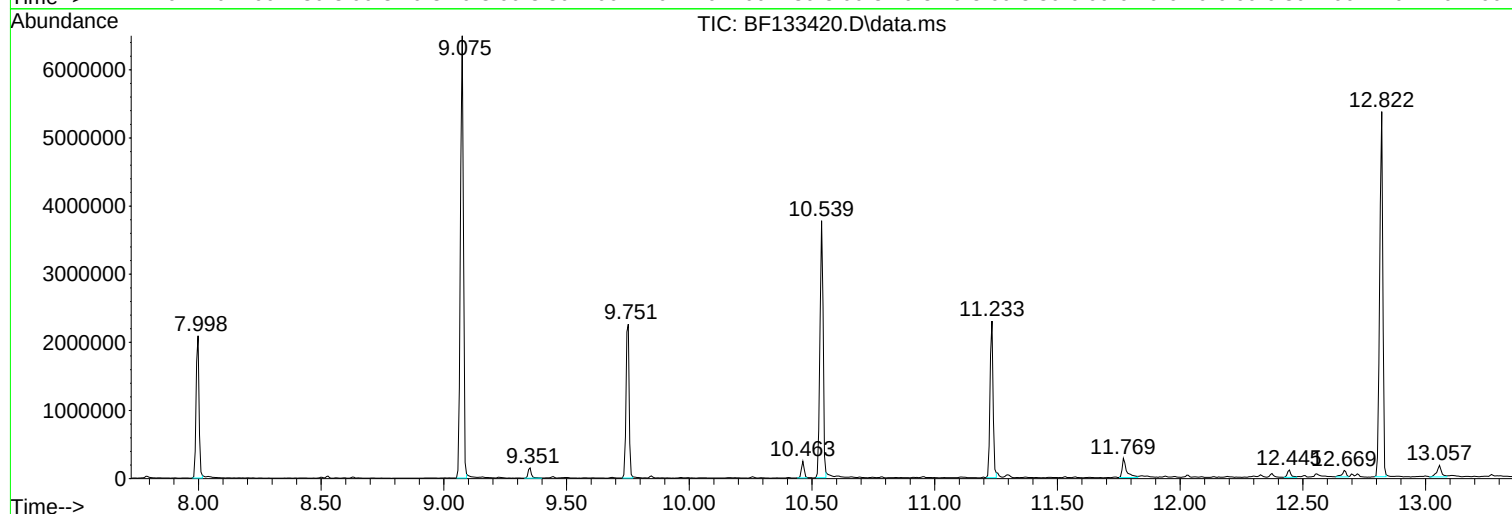
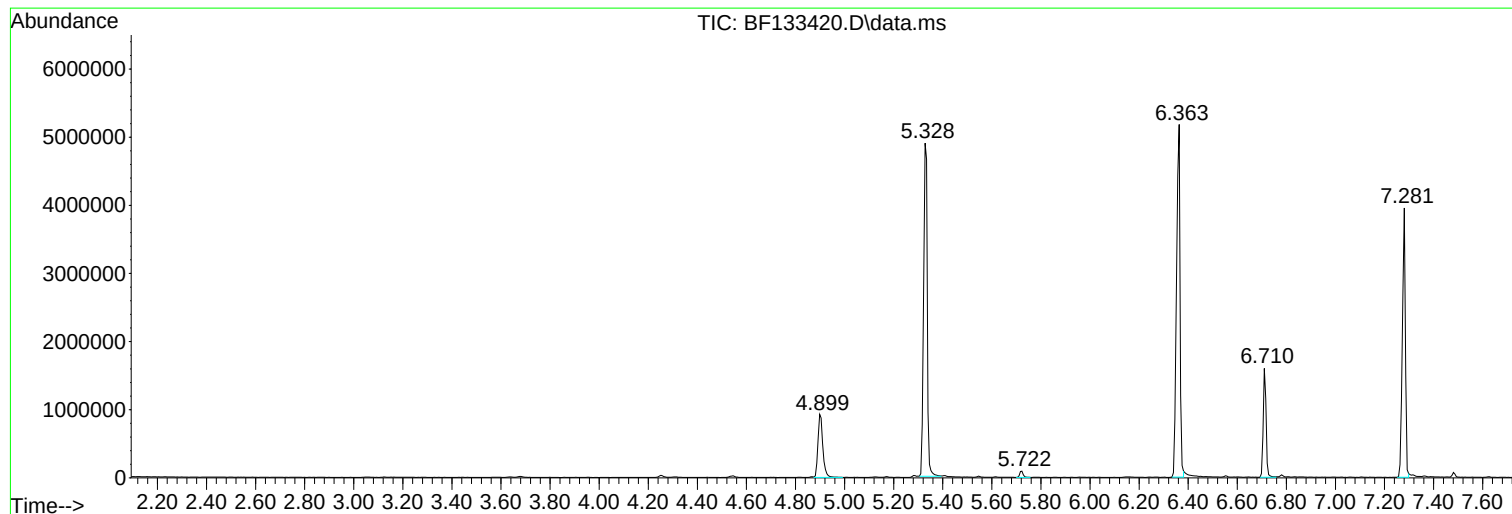
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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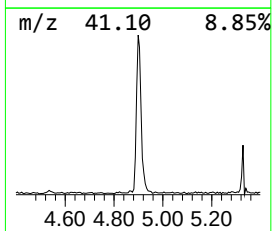
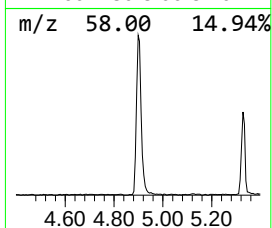
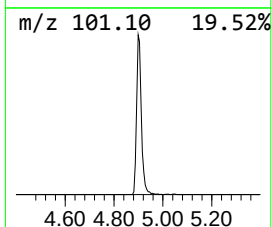
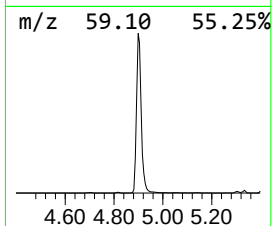
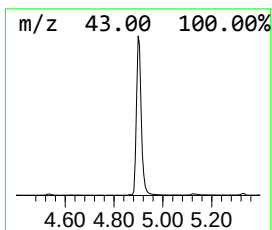
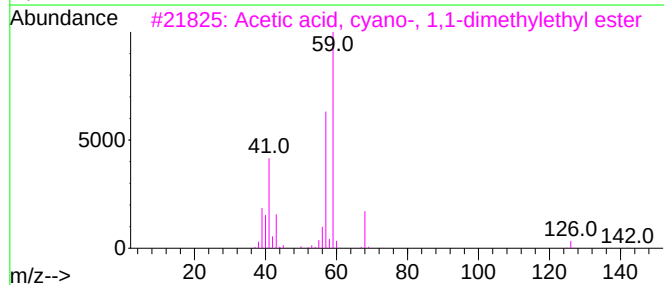
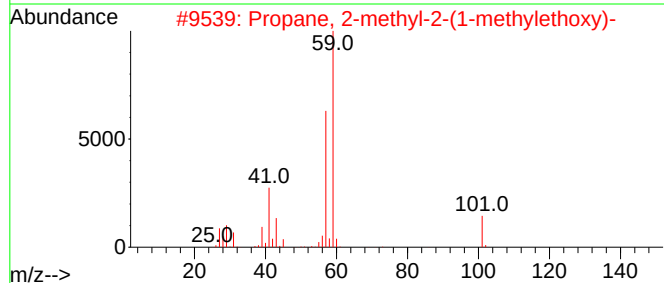
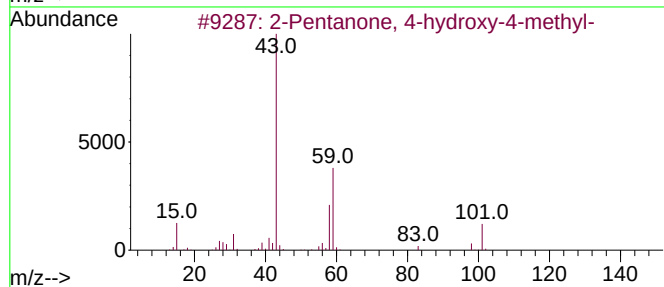
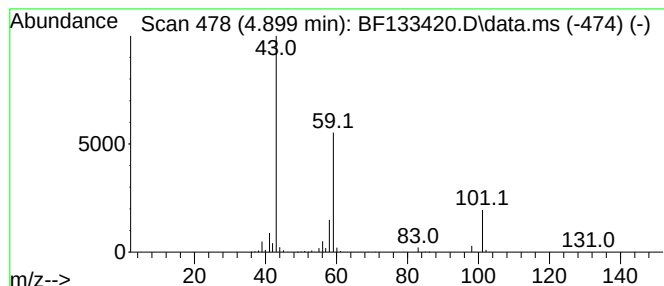
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.899	19.47 ng	1268190	1,4-Dichlorobenzene-d4	6.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		



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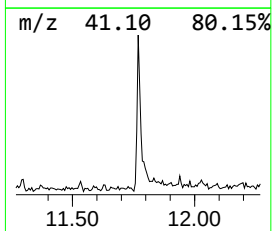
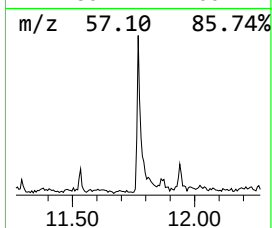
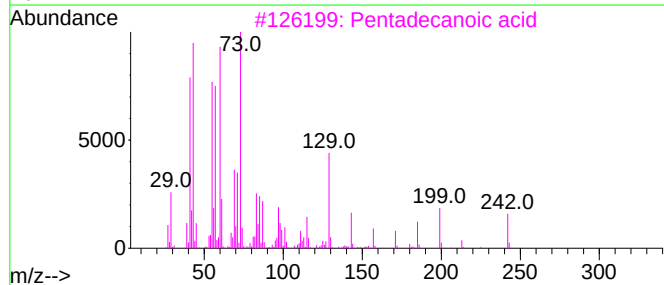
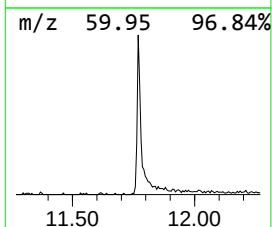
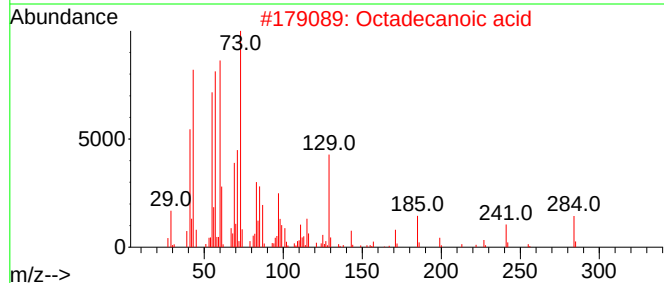
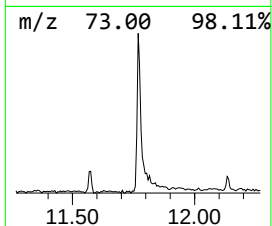
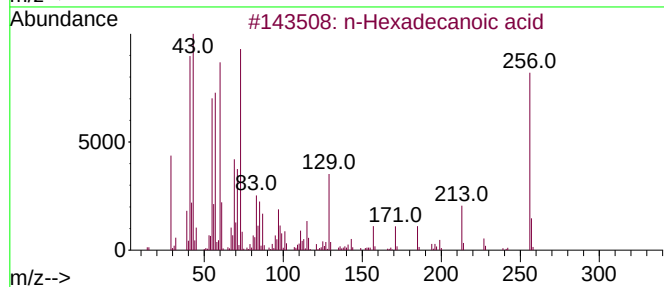
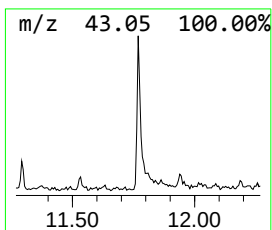
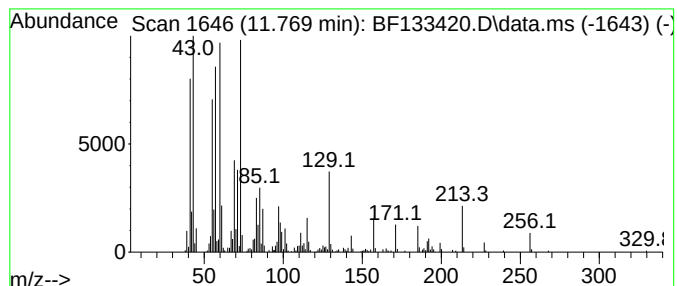
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.769	3.56 ng	355323	Phenanthrene-d10	11.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	93
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	90
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	83



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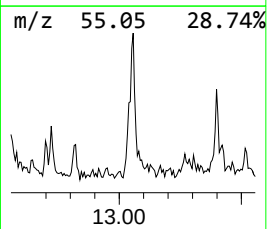
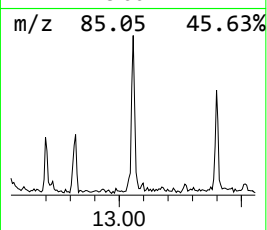
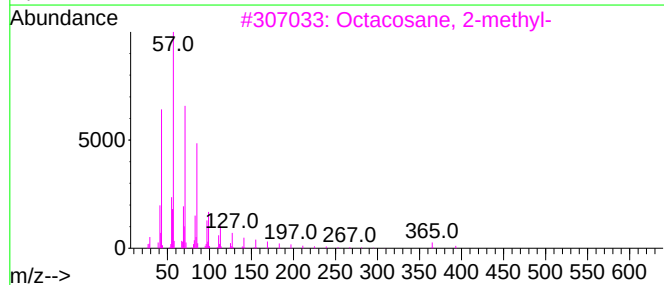
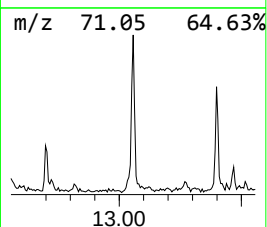
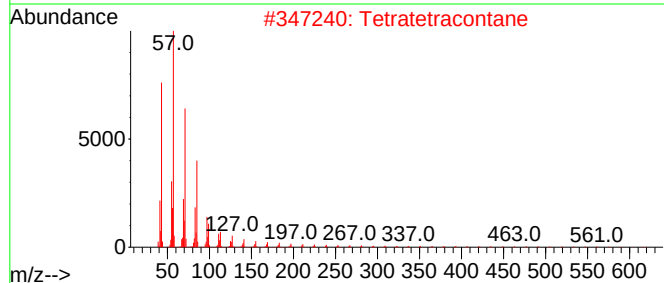
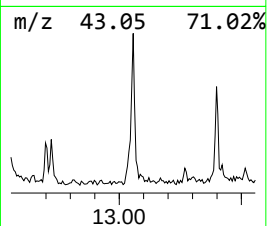
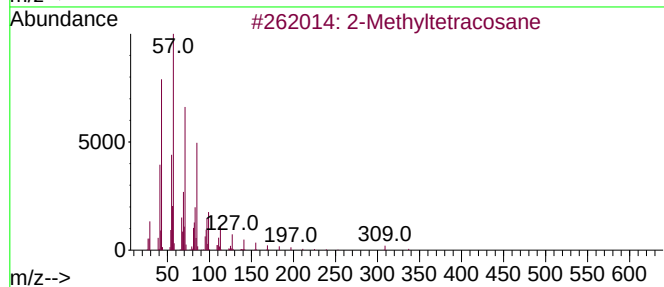
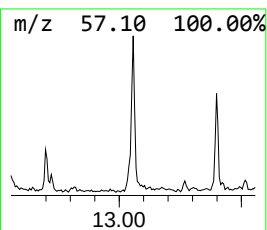
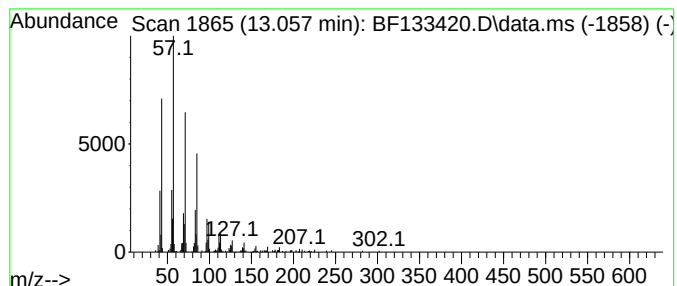
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Methyltetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.057	3.67 ng	218997	Chrysene-d12	13.868

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methyltetracosane	352	C25H52	001560-78-7	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
4		Hexacosane	366	C26H54	000630-01-3	87
5		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	87



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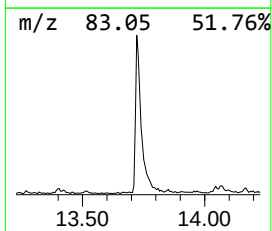
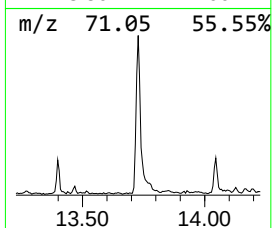
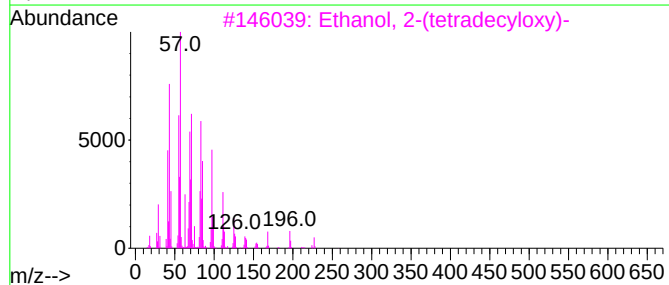
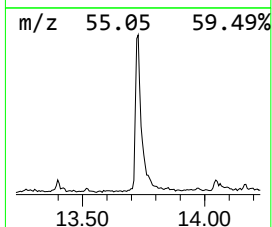
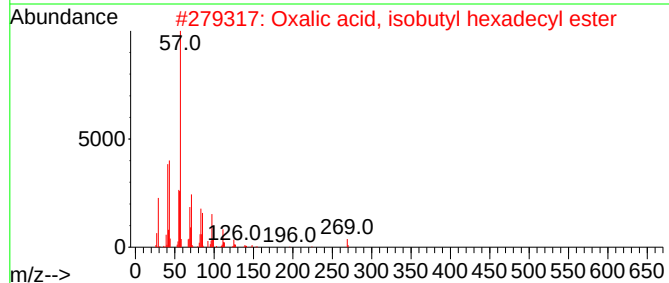
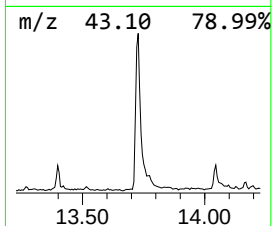
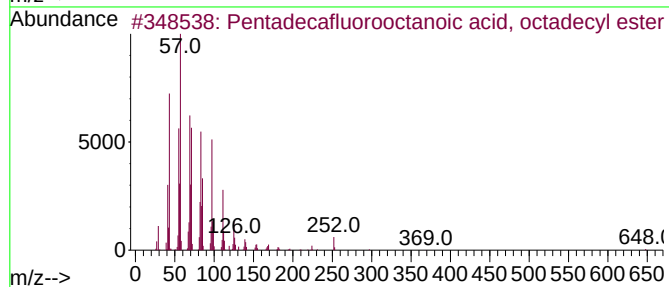
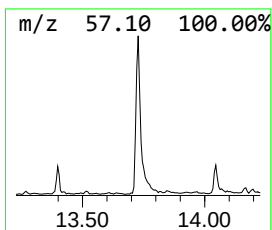
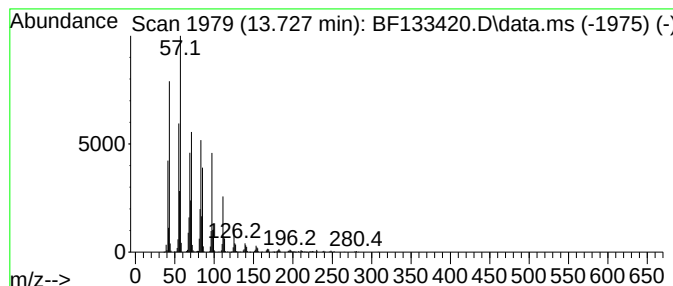
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Peak Number 4 Pentadecafluorooctanoic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.727	23.22 ng	1386890	Chrysene-d12	13.868

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
4			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	76
5			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	76



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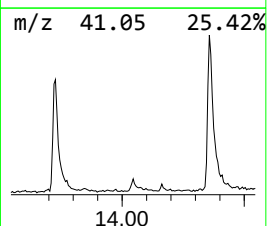
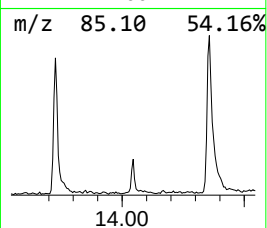
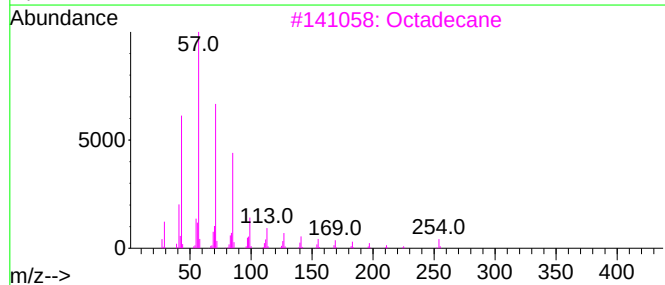
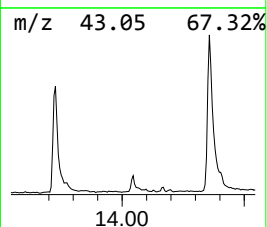
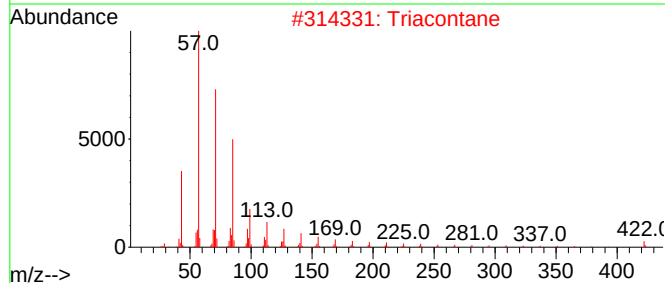
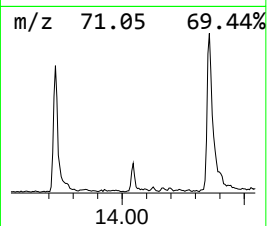
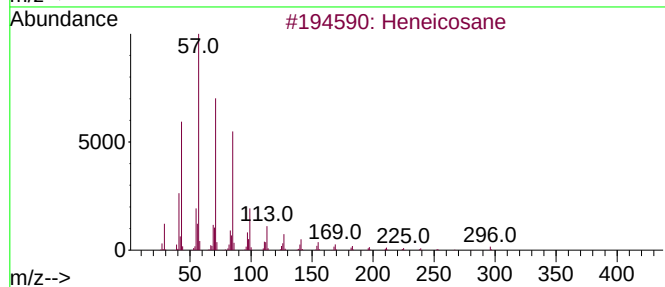
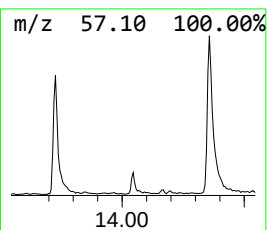
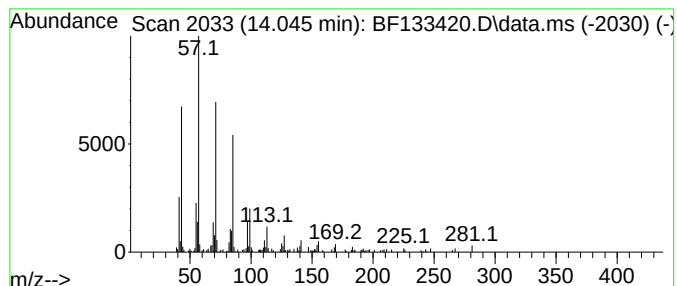
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.045	2.37 ng	141450	Chrysene-d12	13.868

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	90
2			Triacontane	422	C30H62	000638-68-6	90
3			Octadecane	254	C18H38	000593-45-3	86
4			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
Acq On : 17 May 2023 18:49
Operator : CG\JU
Sample : 02812-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WG-0516-B5-0-10

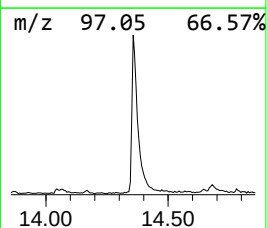
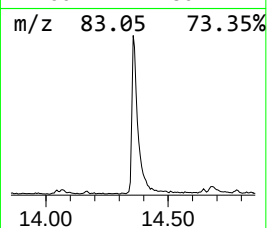
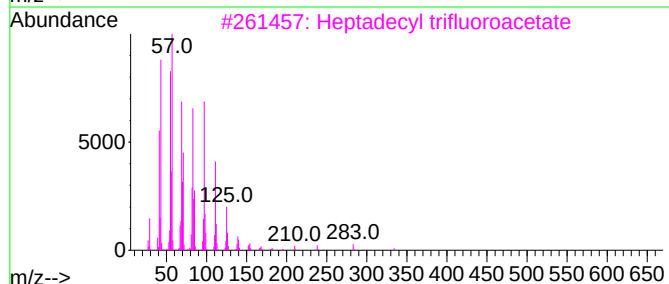
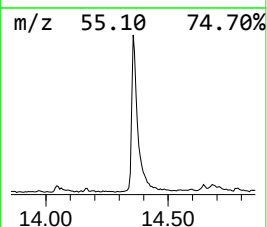
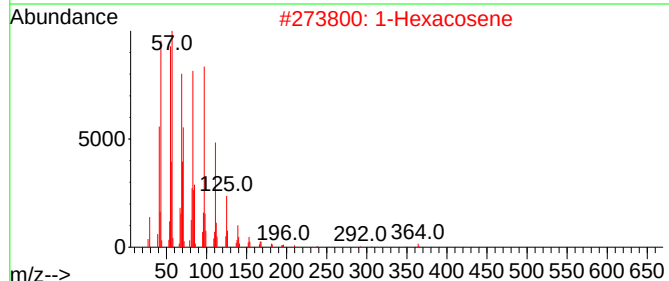
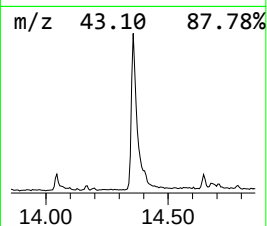
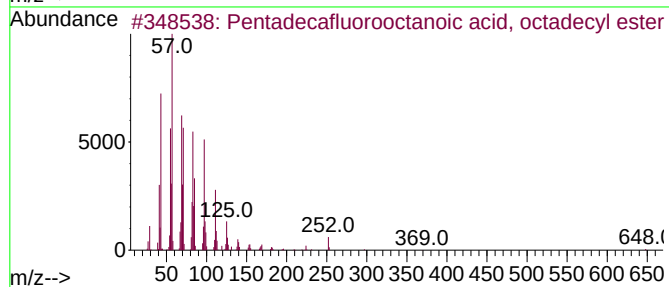
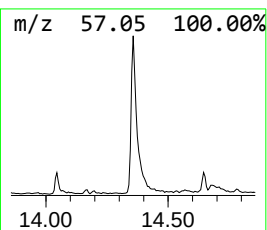
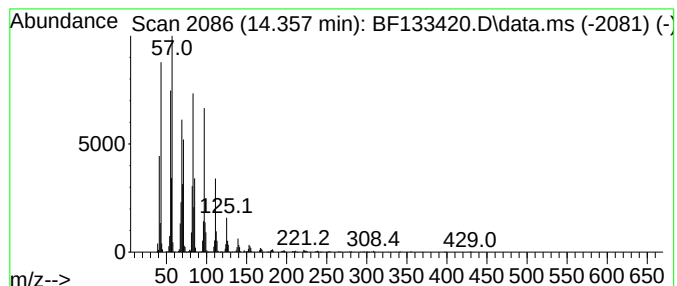
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1-Hexacosene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.357	41.69 ng	2489430	Chrysene-d12	13.868

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			1-Hexacosene	364	C26H52	018835-33-1	93
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
4			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5			1-Hexacosanol	382	C26H54O	000506-52-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
Acq On : 17 May 2023 18:49
Operator : CG\JU
Sample : 02812-02
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ClientSampleId :
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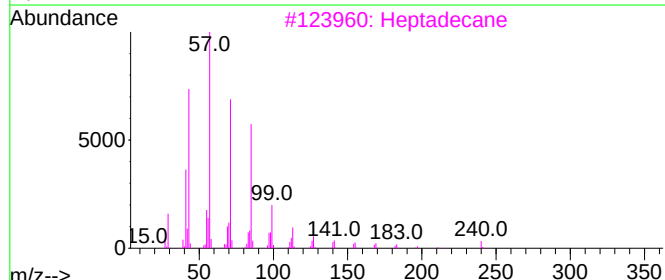
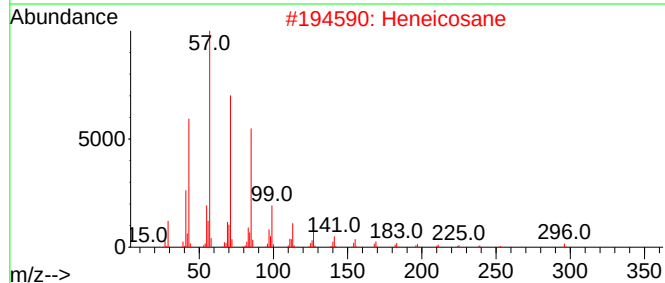
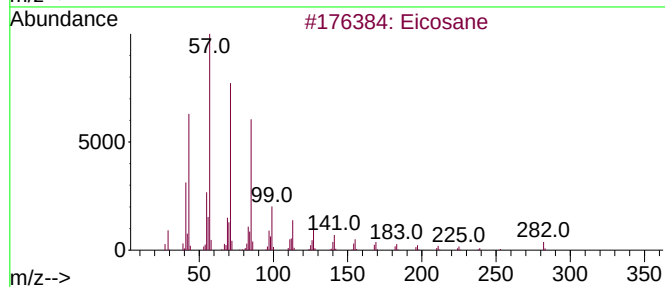
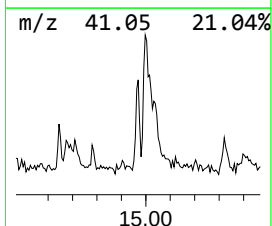
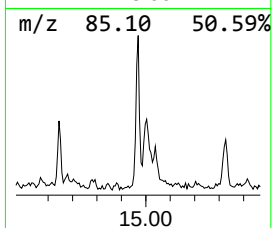
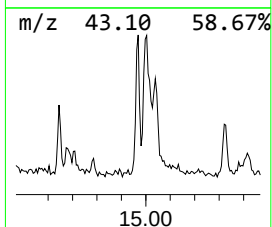
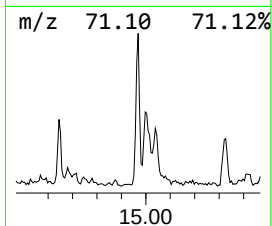
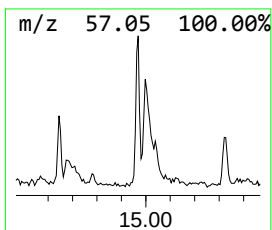
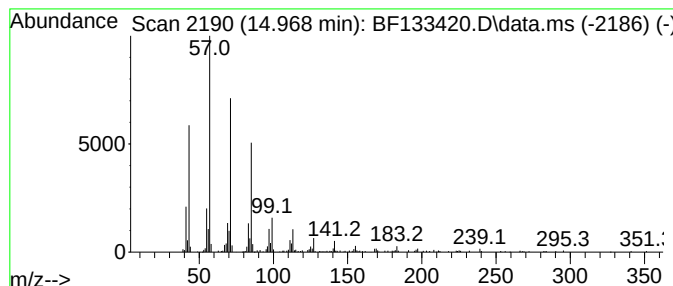
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.969	2.88 ng	205965	Perylene-d12	15.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Heneicosane	296	C21H44	000629-94-7	91
3			Heptadecane	240	C17H36	000629-78-7	91
4			Hentriacontane	437	C31H64	000630-04-6	91
5			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
Acq On : 17 May 2023 18:49
Operator : CG\JU
Sample : 02812-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

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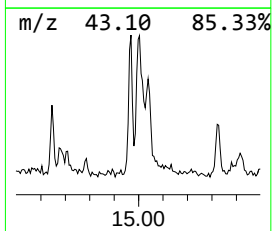
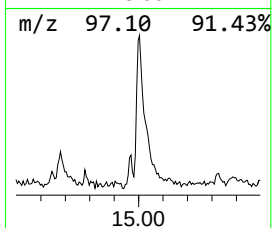
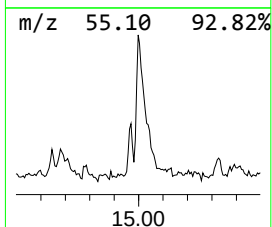
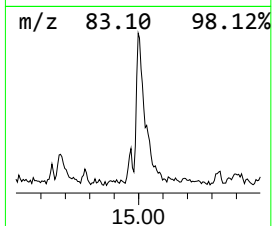
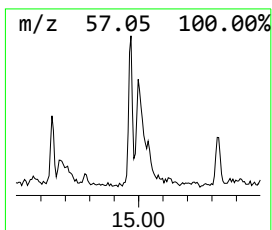
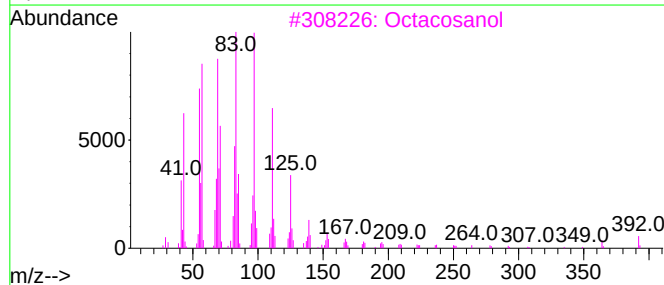
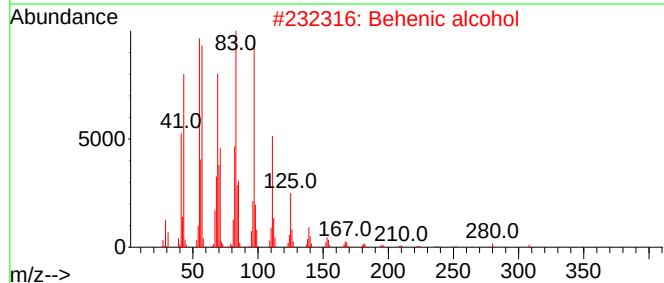
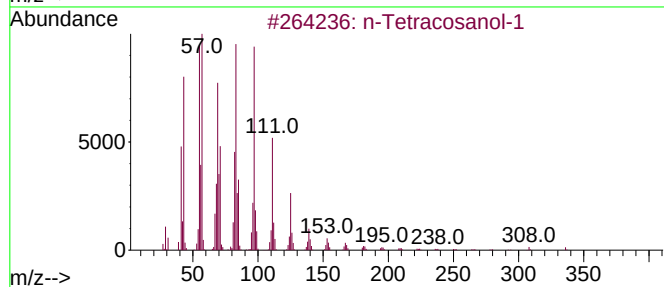
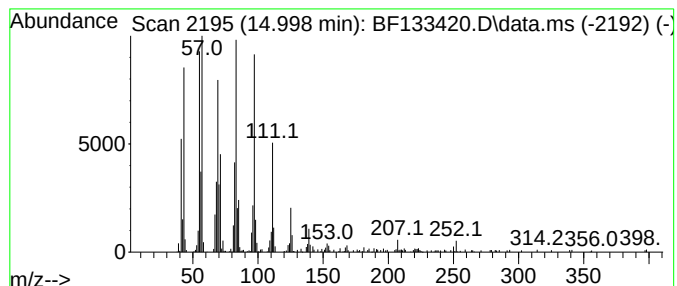
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Tetracosanol-1 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.998	5.75 ng	410780	Perylene-d12	15.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
2		Behenic alcohol	326	C22H46O	000661-19-8	95
3		Octacosanol	410	C28H58O	000557-61-9	94
4		1-Hexacosene	364	C26H52	018835-33-1	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
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Sample : 02812-02
Misc :
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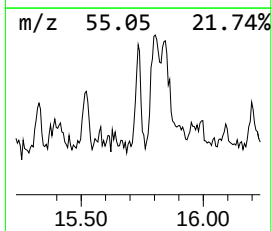
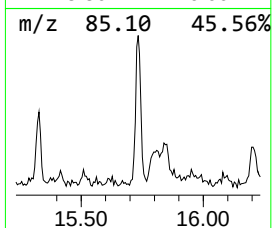
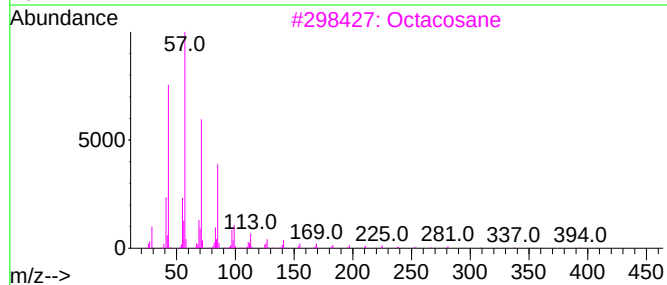
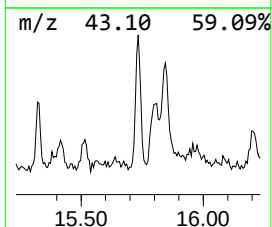
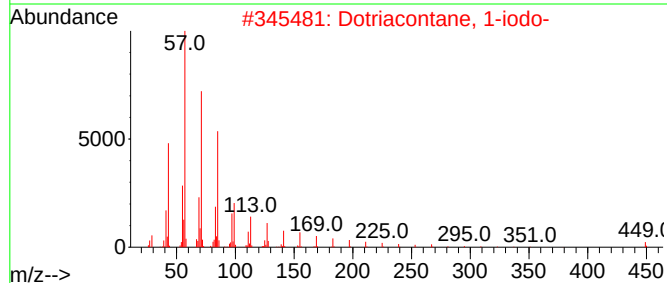
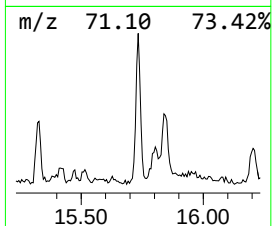
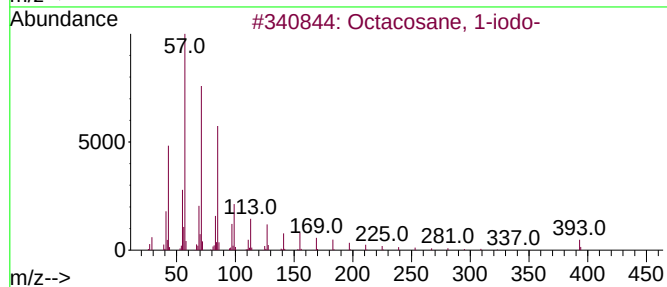
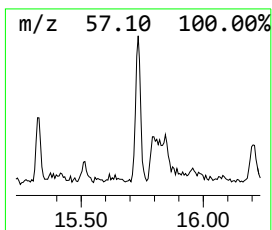
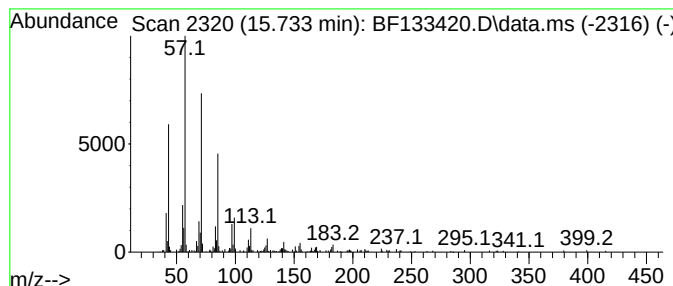
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octacosane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.733	2.65 ng	189310	Perylene-d12	15.292	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane, 1-iodo-	520	C28H57I	1000406-32-2	91
2	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3	Octacosane	394	C28H58	000630-02-4	91
4	Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	90
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
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Instrument :
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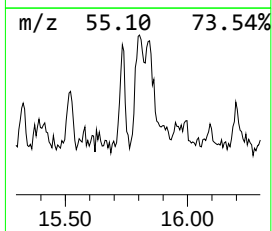
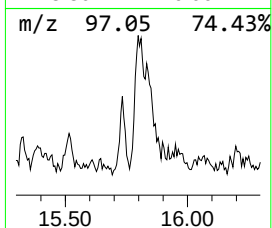
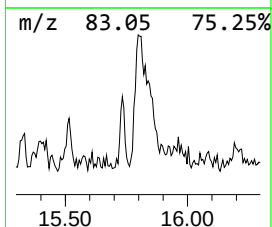
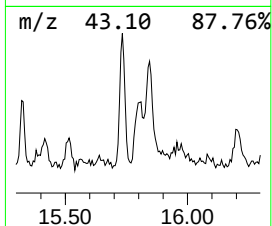
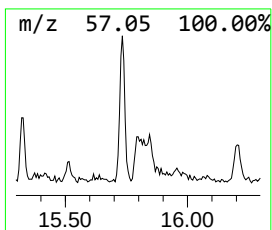
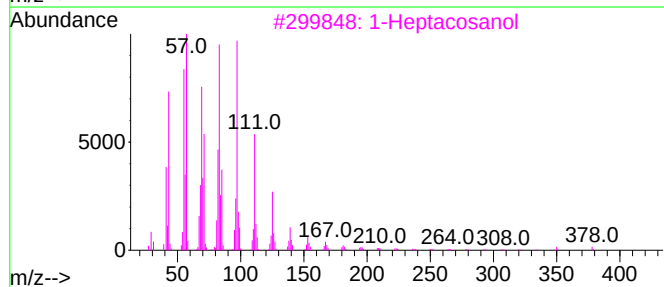
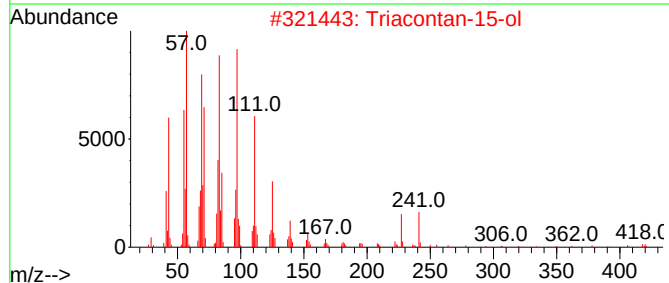
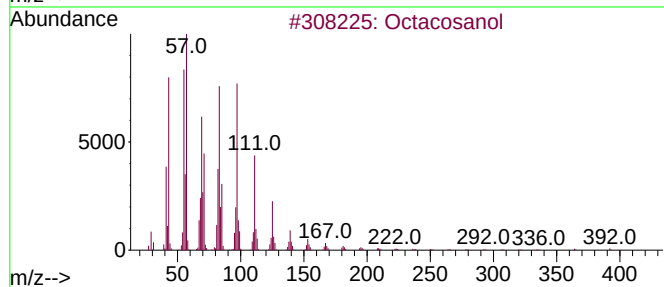
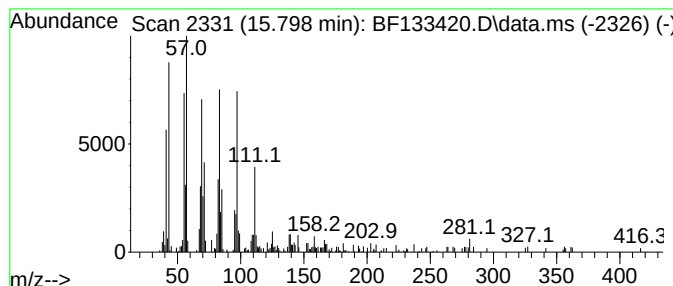
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octacosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.798	2.90 ng	207472	Perylene-d12	15.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosanol	410	C28H58O	000557-61-9	91
2			Triacontan-15-ol	438	C30H62O	031849-26-0	90
3			1-Heptacosanol	396	C27H56O	002004-39-9	90
4			1-Hexacosanol	382	C26H54O	000506-52-5	87
5			Heptacosyl acetate	438	C29H58O2	1000351-78-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
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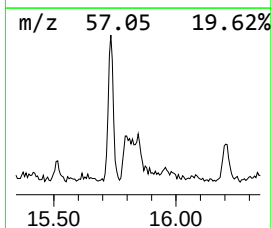
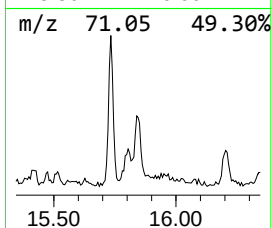
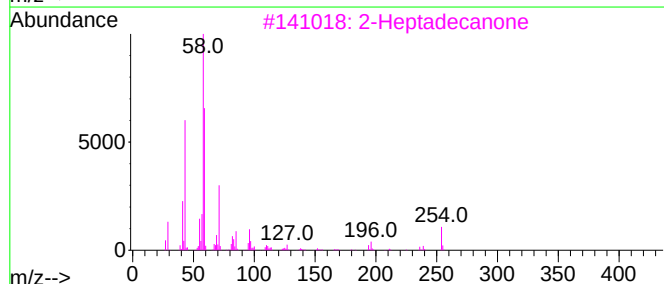
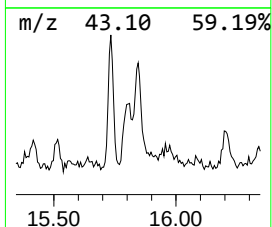
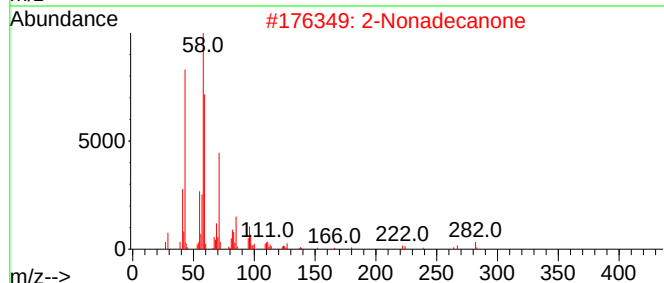
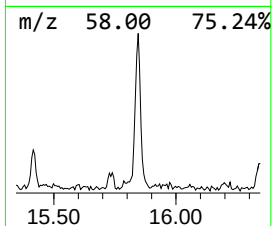
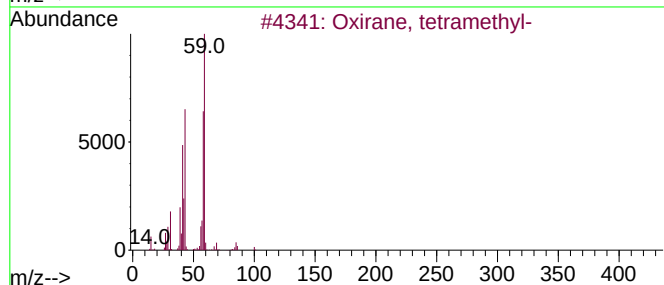
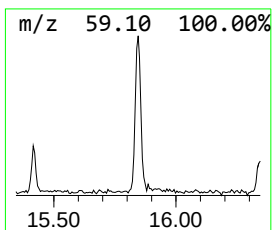
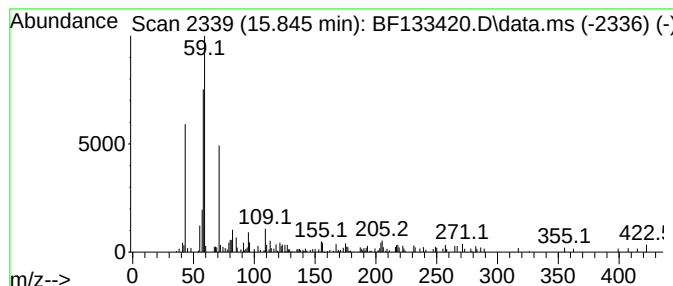
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Oxirane, tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.845	3.15 ng	225234	Perylene-d12	15.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, tetramethyl-	100	C6H12O	005076-20-0	50
2			2-Nonadecanone	282	C19H38O	000629-66-3	42
3			2-Heptadecanone	254	C17H34O	002922-51-2	38
4			2-Tetradecanone	212	C14H28O	002345-27-9	38
5			Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
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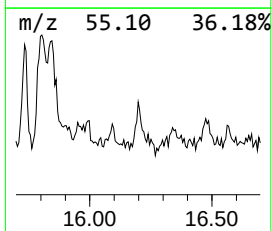
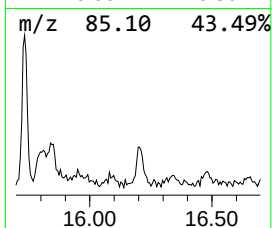
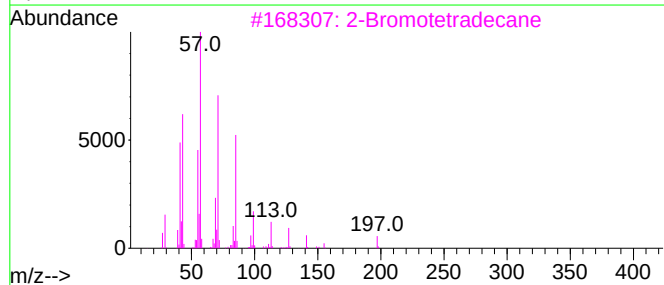
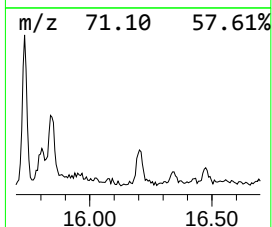
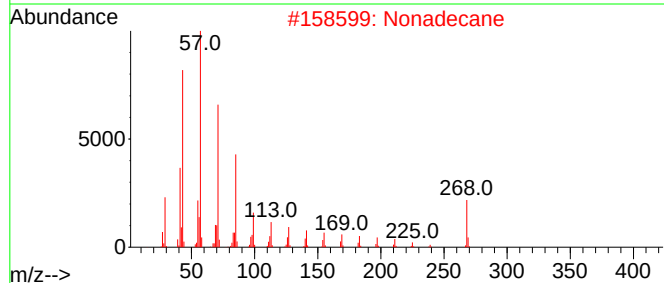
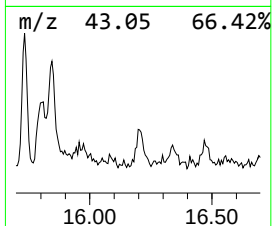
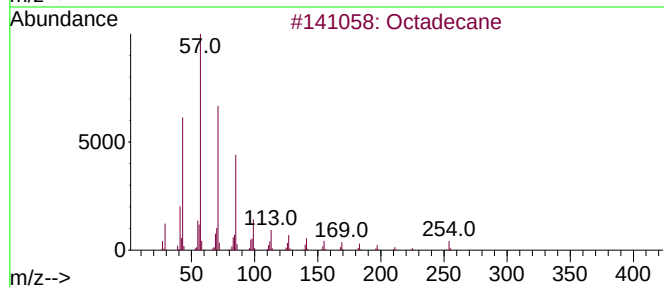
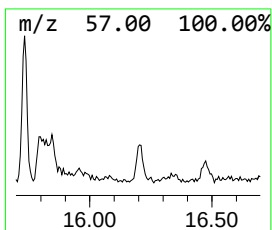
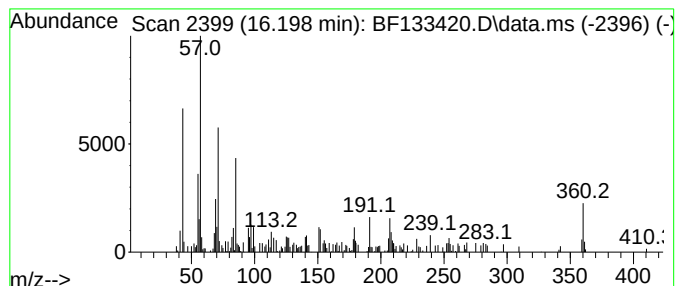
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.198	2.17 ng	155106	Perylene-d12	15.292

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	78
2		Nonadecane	268	C19H40	000629-92-5	60
3		2-Bromotetradecane	276	C14H29Br	074036-95-6	53
4		10-Methylnonadecane	282	C20H42	056862-62-5	53
5		Hentriacontane	437	C31H64	000630-04-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
Acq On : 17 May 2023 18:49
Operator : CG\JU
Sample : 02812-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WG-0516-B5-0-10

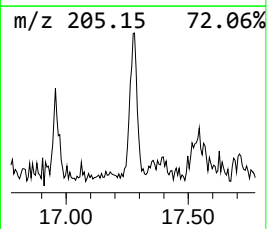
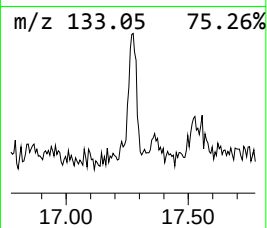
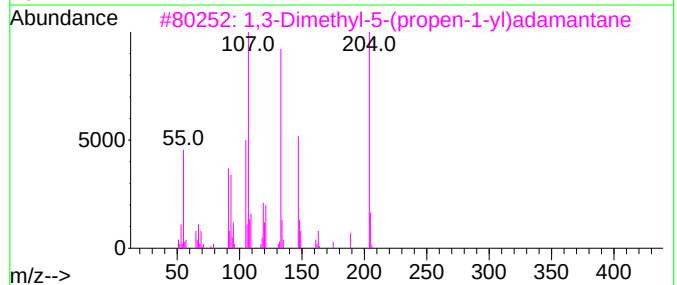
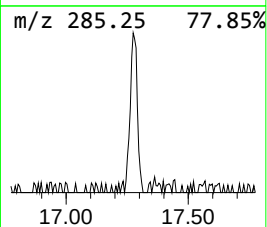
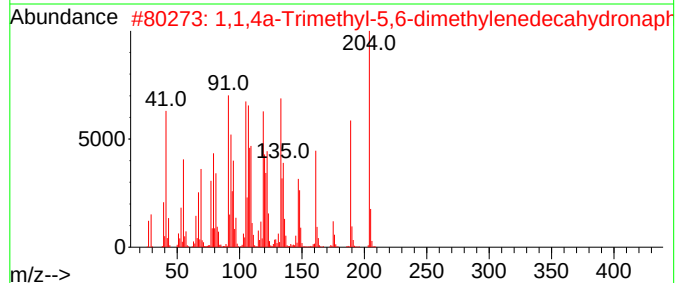
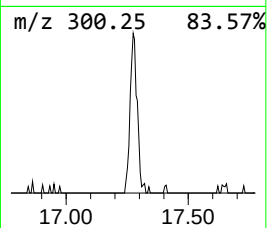
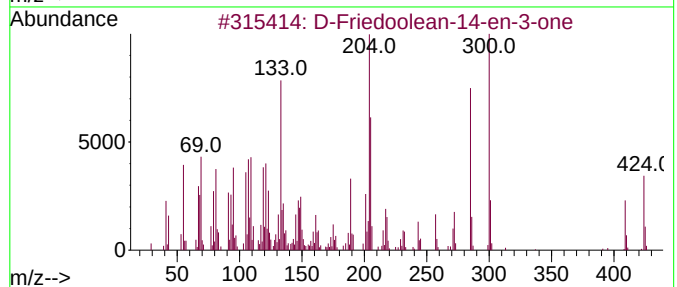
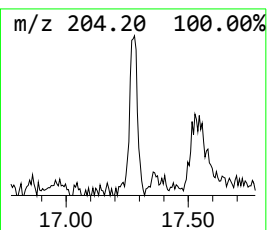
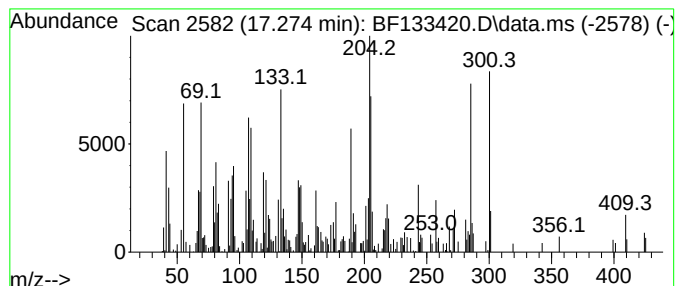
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 D-Friedoolean-14-en-3-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.274	3.54 ng	253151	Perylene-d12	15.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D-Friedoolean-14-en-3-one	424	C30H48O	000514-07-8	95
2			1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	49
3			1,3-Dimethyl-5-(propen-1-yl)adam...	204	C15H24	057040-48-9	46
4			2(1H)-Phenanthrene, 3,4,4a,9,1...	300	C20H28O2	000472-37-7	41
5			Phenanthrene, 1,2,3,4,4a,9,10,10...	300	C21H32O	010064-26-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
Acq On : 17 May 2023 18:49
Operator : CG\JU
Sample : 02812-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

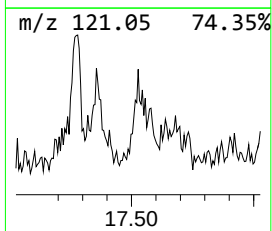
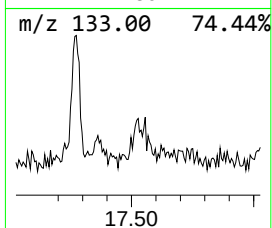
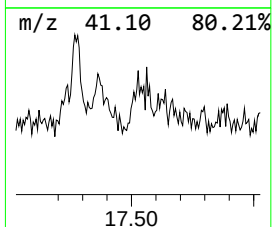
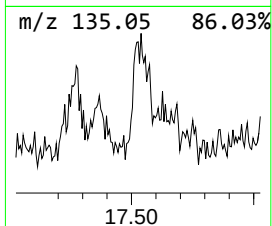
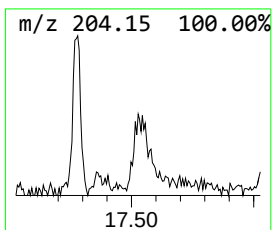
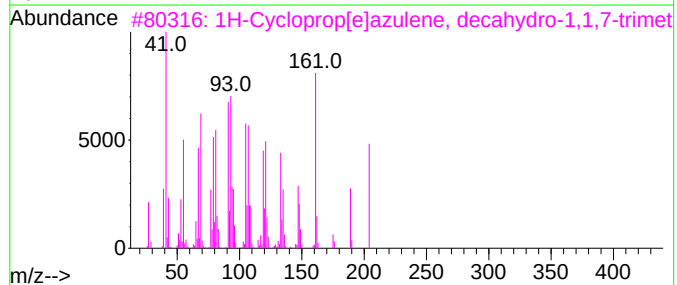
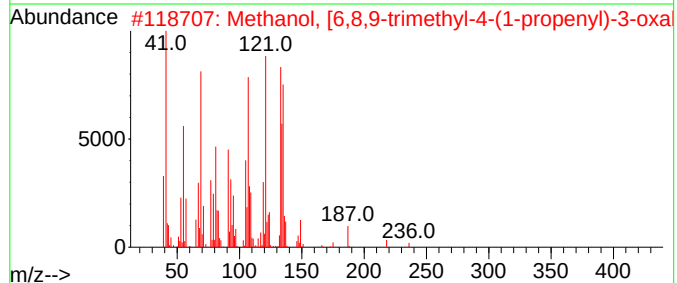
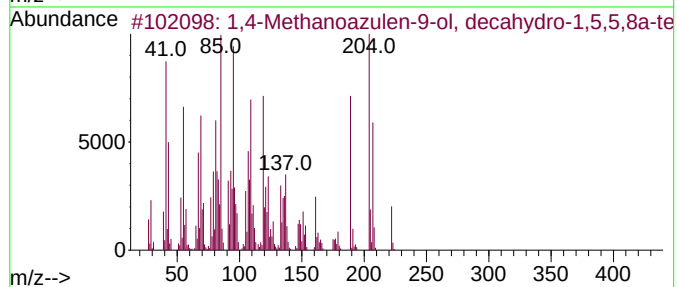
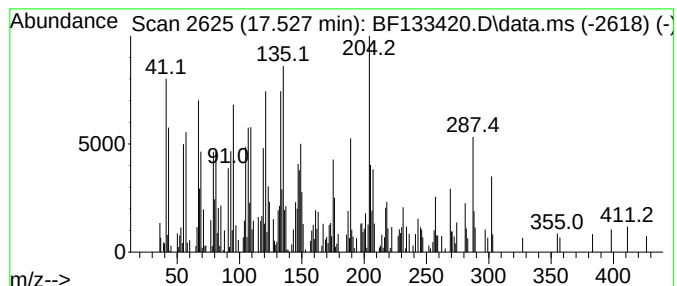
Instrument :
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ClientSampleId :
WG-0516-B5-0-10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1,4-Methanoazulen-9-ol, dec... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.527	2.56 ng	182801	Perylene-d12	15.292	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Methanoazulen-9-ol, decahydr...	222	C15H26O	000465-24-7	56
2	Methanol, [6,8,9-trimethyl-4-(1-...	236	C15H24O2	1000277-60-9	50
3	1H-Cycloprop[e]azulene, decahydr...	204	C15H24	072747-25-2	47
4	2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	45
5	Dihydrotachysterol	398	C28H46O	000067-96-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051723\
Data File : BF133420.D
Acq On : 17 May 2023 18:49
Operator : CG\JU
Sample : 02812-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WG-0516-B5-0-10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.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
2-Pentanone, 4-...	4.899	19.5	ng	1268190	1	6.710	1302870	20.0
n-Hexadecanoic ...	11.769	3.6	ng	355323	4	11.233	1993760	20.0
2-Methyltetraaco...	13.057	3.7	ng	218997	5	13.868	1194310	20.0
Pentadecafluoro...	13.727	23.2	ng	1386890	5	13.868	1194310	20.0
Heneicosane	14.045	2.4	ng	141450	5	13.868	1194310	20.0
1-Hexacosene	14.357	41.7	ng	2489430	5	13.868	1194310	20.0
Eicosane	14.969	2.9	ng	205965	6	15.292	1428750	20.0
n-Tetracosanol-1	14.998	5.8	ng	410780	6	15.292	1428750	20.0
Octacosane, 1-i...	15.733	2.6	ng	189310	6	15.292	1428750	20.0
Octacosanol	15.798	2.9	ng	207472	6	15.292	1428750	20.0
Oxirane, tetram...	15.845	3.1	ng	225234	6	15.292	1428750	20.0
Octadecane	16.198	2.2	ng	155106	6	15.292	1428750	20.0
D-Friedoolean-1...	17.274	3.5	ng	253151	6	15.292	1428750	20.0
1,4-Methanoazul...	17.527	2.6	ng	182801	6	15.292	1428750	20.0