

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
 Data File : BF135128.D
 Acq On : 07 Sep 2023 11:35
 Operator : CG\JU
 Sample : 04283-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-1

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135128.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.134	4	8	16	rVB	40899	57898	2.32%	0.300%
2	5.004	492	496	506	rVB	73977	90406	3.62%	0.469%
3	5.404	559	564	567	rBV	2386548	2154650	86.27%	11.171%
4	6.410	730	735	738	rBV	2042683	2140245	85.70%	11.096%
5	6.604	763	768	777	rBV	83283	85012	3.40%	0.441%
6	6.769	793	796	800	rBV	731808	573087	22.95%	2.971%
7	7.334	887	892	895	rBV	1510223	1390302	55.67%	7.208%
8	8.051	1009	1014	1017	rBV	893483	796336	31.89%	4.129%
9	9.128	1192	1197	1200	rBV	2613892	2497503	100.00%	12.948%
10	9.245	1214	1217	1221	rBV	38925	34669	1.39%	0.180%
11	9.810	1308	1313	1316	rBV	948563	885237	35.44%	4.590%
12	10.016	1344	1348	1351	rBV2	28064	33218	1.33%	0.172%
13	10.604	1443	1448	1451	rBV	1674263	1559060	62.42%	8.083%
14	10.751	1468	1473	1478	rVB3	63467	82439	3.30%	0.427%
15	10.892	1494	1497	1502	rBV	18783	26339	1.05%	0.137%
16	10.951	1502	1507	1515	rVB3	57502	107141	4.29%	0.555%
17	11.298	1562	1566	1569	rBV	958155	869015	34.80%	4.505%
18	11.704	1629	1635	1638	rVB	23772	26976	1.08%	0.140%
19	11.839	1652	1658	1666	rBV	719839	915765	36.67%	4.748%
20	11.986	1679	1683	1686	rBV	194816	237444	9.51%	1.231%
21	12.016	1686	1688	1700	rVB	122055	166113	6.65%	0.861%
22	12.522	1771	1774	1777	rBV2	30626	50410	2.02%	0.261%
23	12.628	1787	1792	1803	rBV	415604	469474	18.80%	2.434%
24	12.898	1834	1838	1842	rBV	2347051	2124823	85.08%	11.016%
25	12.951	1842	1847	1854	rVB2	40453	81713	3.27%	0.424%
26	13.133	1870	1878	1883	rVB4	29778	55438	2.22%	0.287%
27	13.745	1975	1982	1989	rBV2	50256	99934	4.00%	0.518%
28	13.810	1989	1993	2004	rBV	155486	195730	7.84%	1.015%
29	13.957	2013	2018	2022	rBV	605914	581408	23.28%	3.014%
30	14.133	2045	2048	2053	rBV	26997	32885	1.32%	0.170%
31	14.345	2080	2084	2091	rVB3	56369	90230	3.61%	0.468%
32	14.439	2097	2100	2108	rVB4	24465	33474	1.34%	0.174%
33	14.874	2168	2174	2181	rVB2	57661	106059	4.25%	0.550%
34	15.427	2263	2268	2272	rBV	434491	566530	22.68%	2.937%
35	16.121	2381	2386	2396	rVB2	29125	71168	2.85%	0.369%

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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-BF083023.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

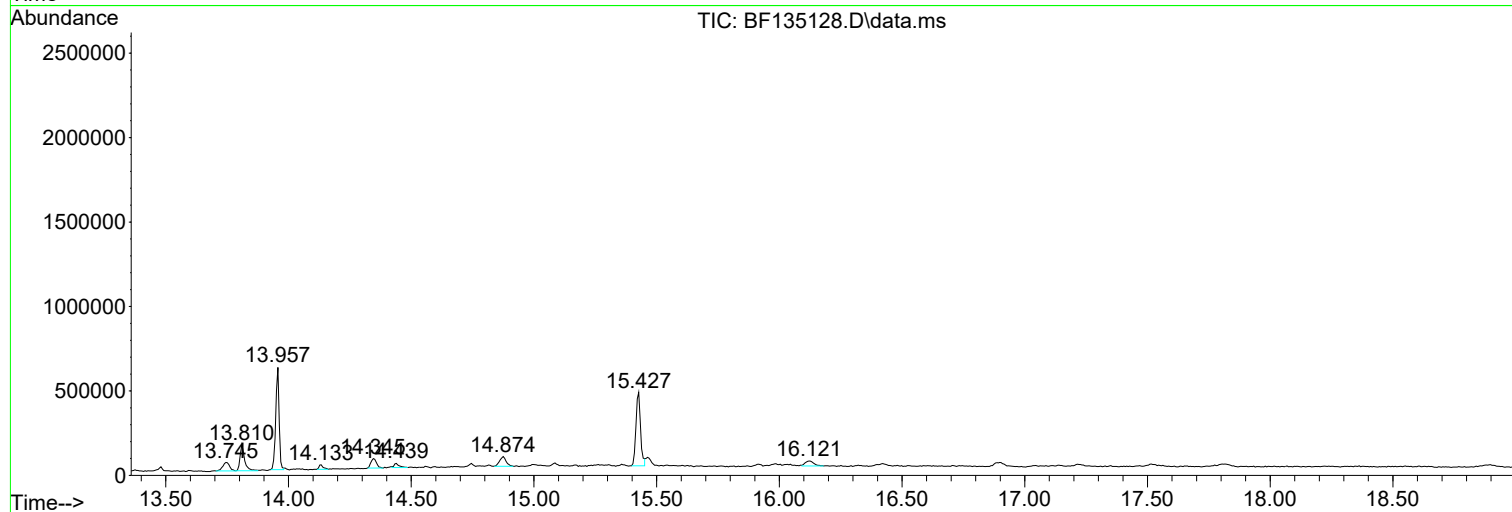
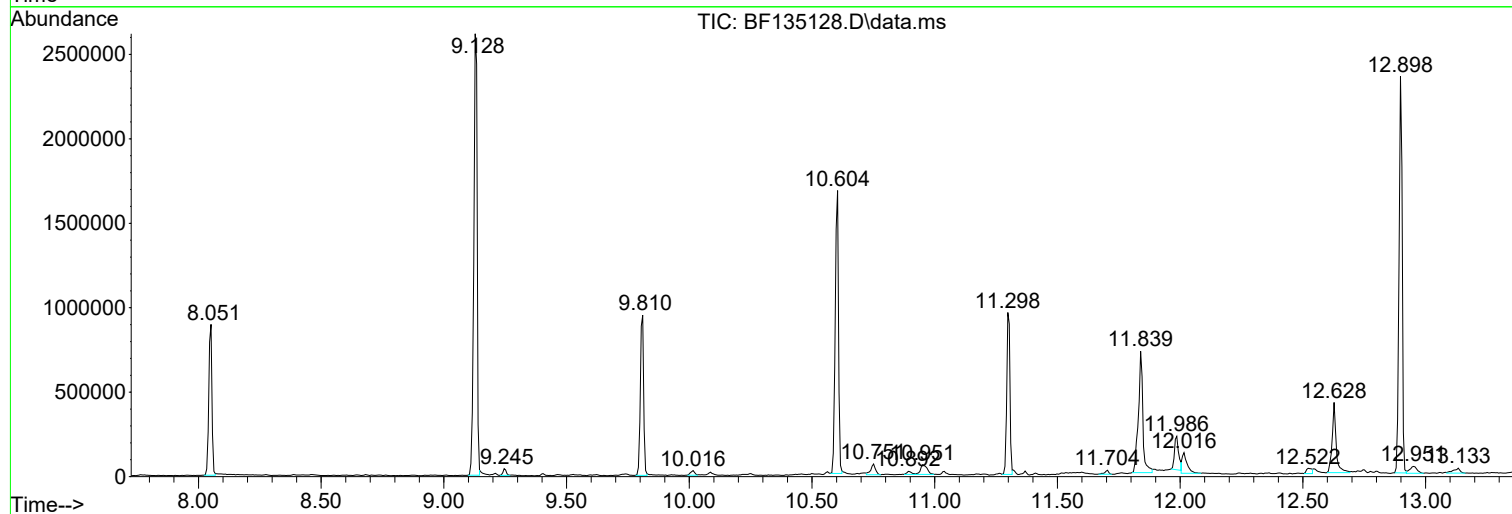
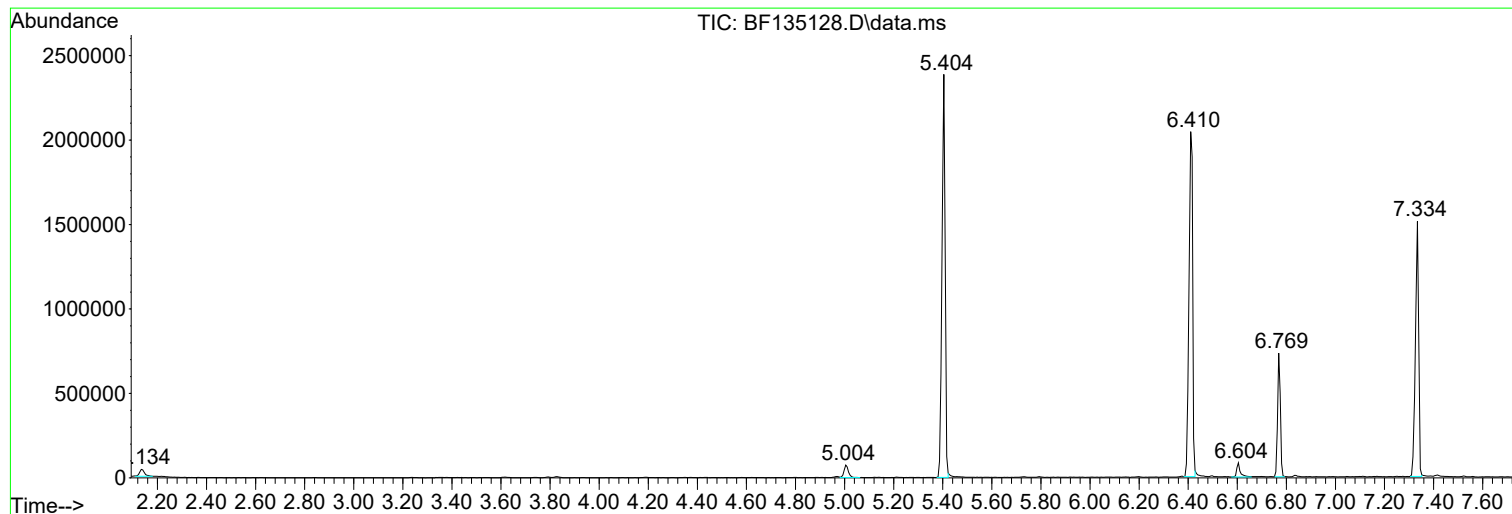
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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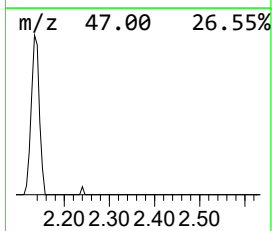
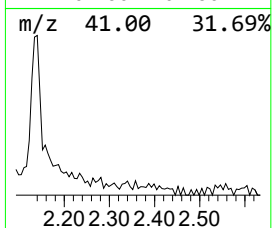
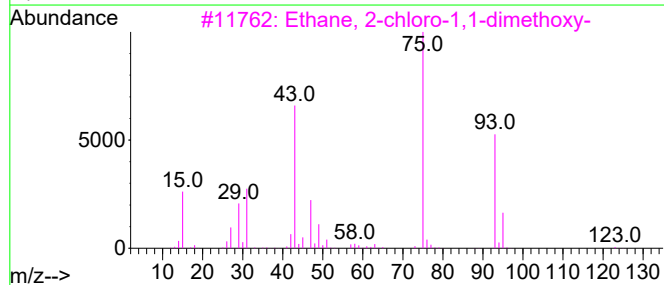
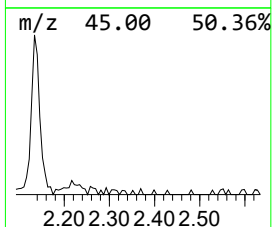
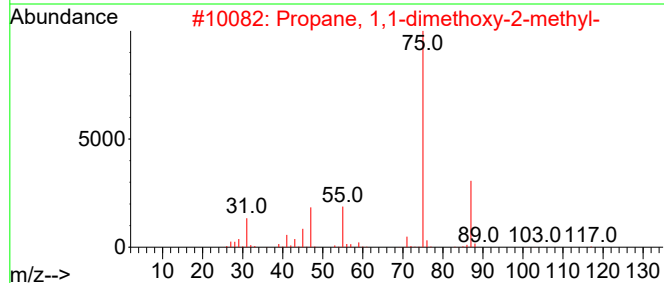
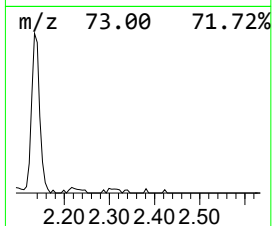
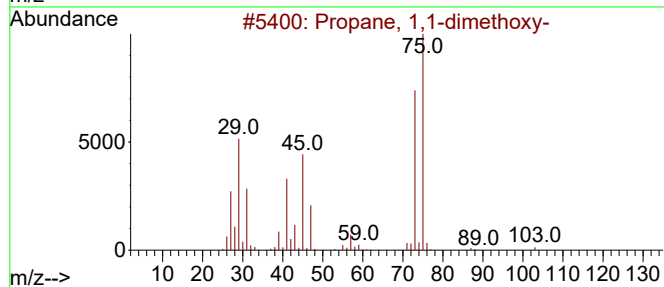
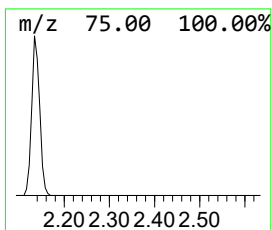
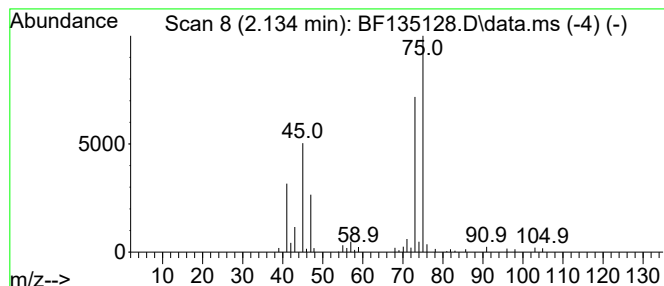
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.134	2.02 ng	57898	1,4-Dichlorobenzene-d4	6.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	74
2			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	40
3			Ethane, 2-chloro-1,1-dimethoxy-	124	C4H9ClO2	000097-97-2	33
4			2,5-Diethoxy-3-methyl-3-vinylhexane	214	C13H26O2	1000432-70-8	22
5			Dimethyl-cyano-phosphine	87	C3H6NP	031641-57-3	10



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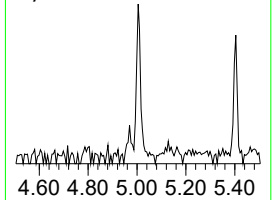
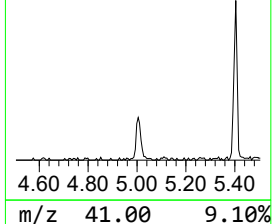
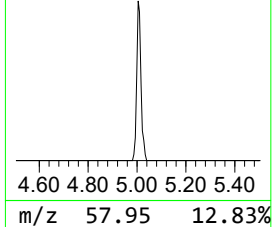
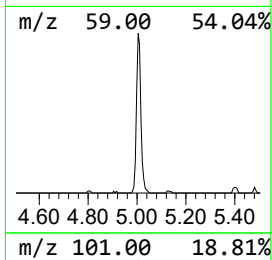
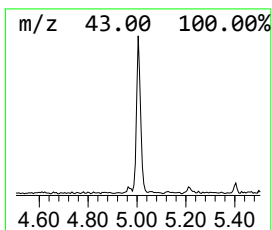
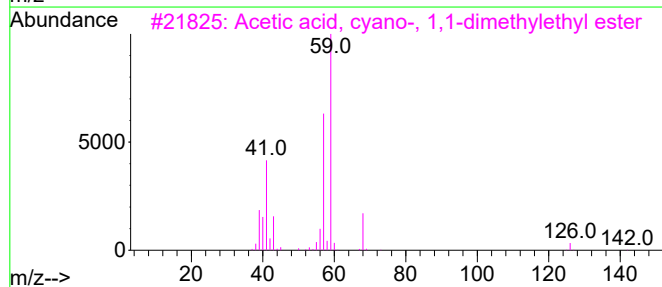
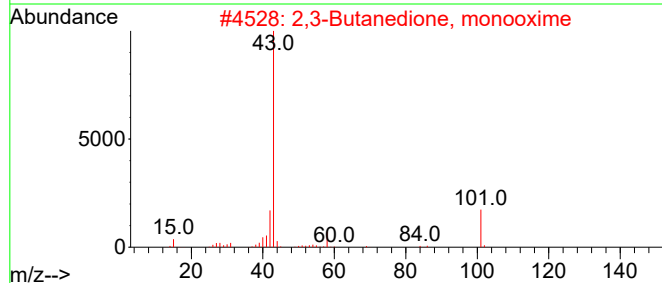
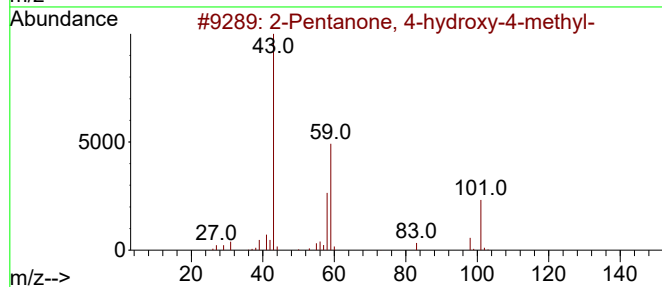
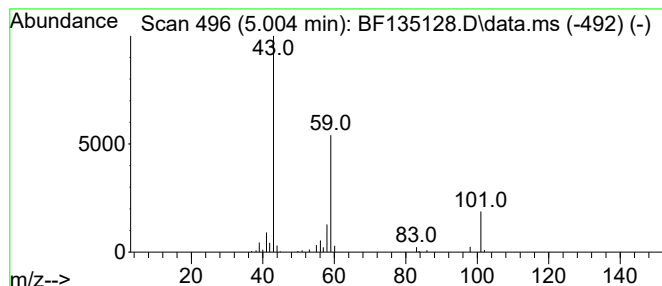
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.004	3.16 ng	90406	1,4-Dichlorobenzene-d4	6.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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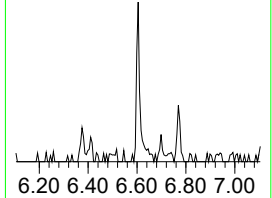
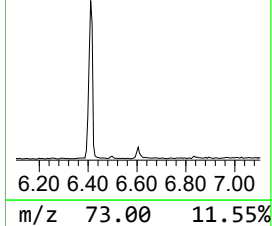
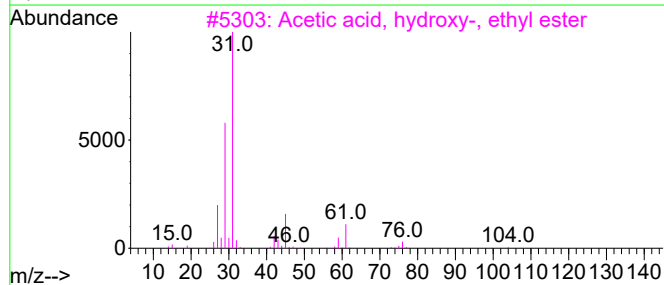
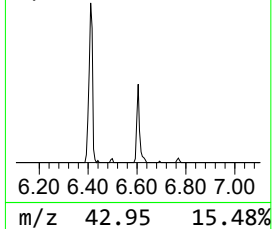
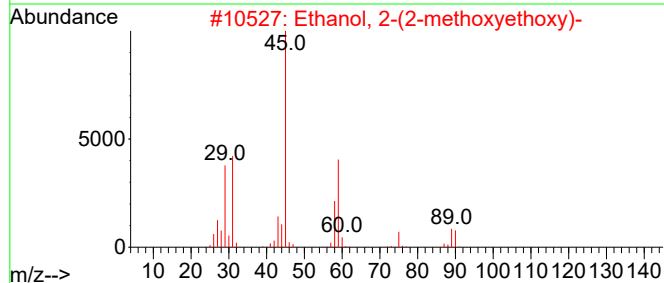
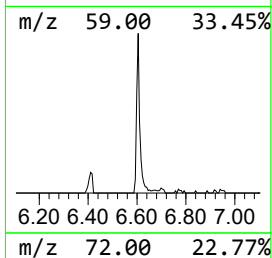
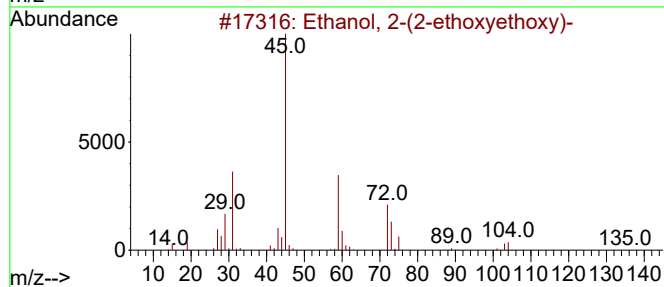
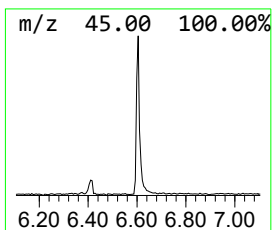
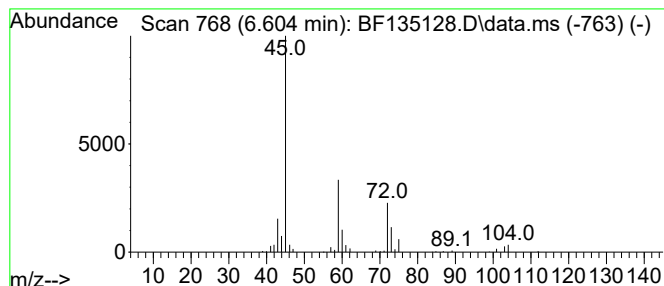
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.604	2.97 ng	85012	1,4-Dichlorobenzene-d4	6.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
2			Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	42
3			Acetic acid, hydroxy-, ethyl ester	104	C4H8O3	000623-50-7	40
4			Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	39
5			2-Pentanol	88	C5H12O	006032-29-7	38



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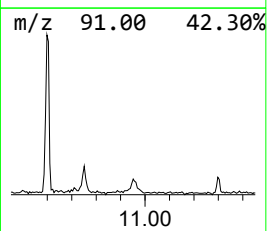
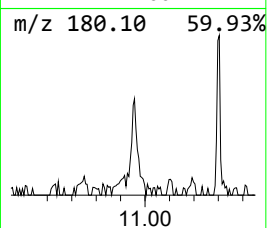
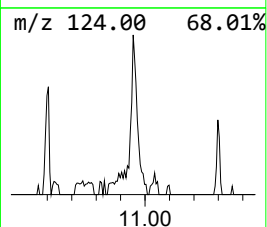
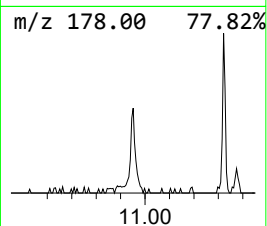
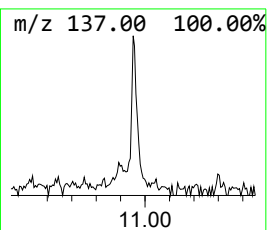
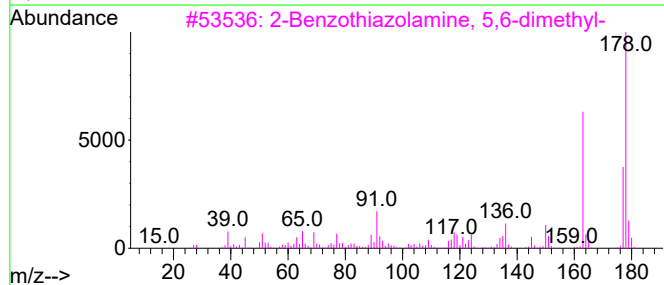
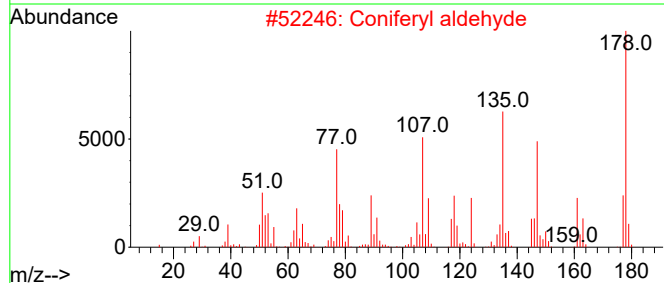
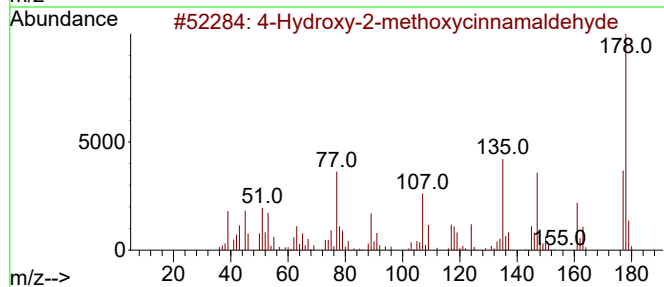
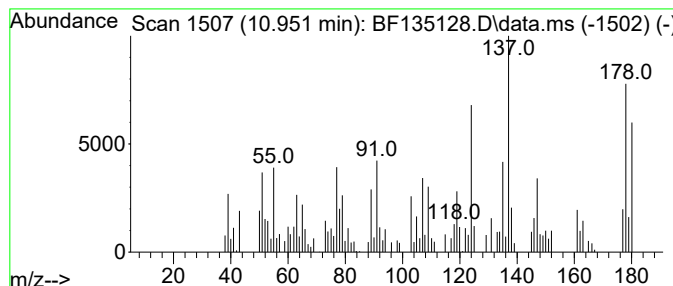
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 unknown10.951 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.951	2.47 ng	107141	Phenanthrene-d10	11.298

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Hydroxy-2-methoxycinnamaldehyde	178	C10H10O3	127321-19-1	46
2		Coniferyl aldehyde	178	C10H10O3	000458-36-6	42
3		2-Benzothiazolamine, 5,6-dimethyl-	178	C9H10N2S	029927-08-0	35
4		3,7-Benzofurandiol, 2,3-dihydro-...	180	C10H12O3	017781-15-6	30
5		2-Oxo-2,3-dihydro-1H-benzimidazo...	178	C8H6N2O3	023814-14-4	25



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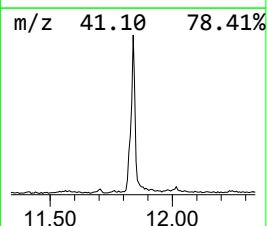
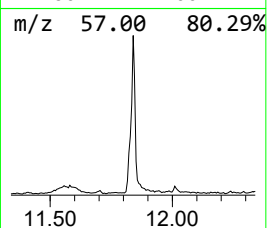
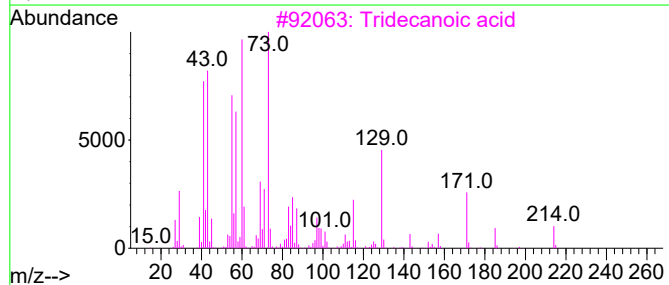
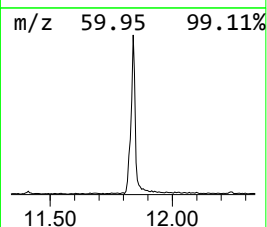
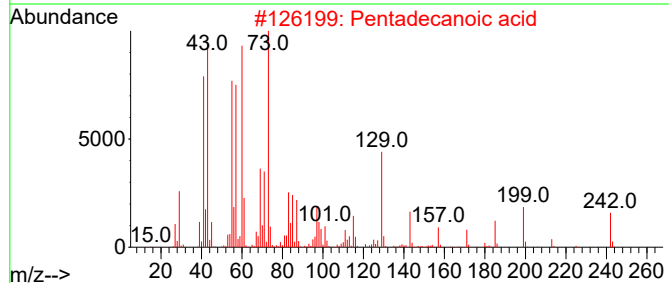
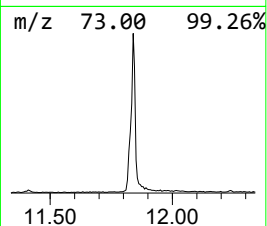
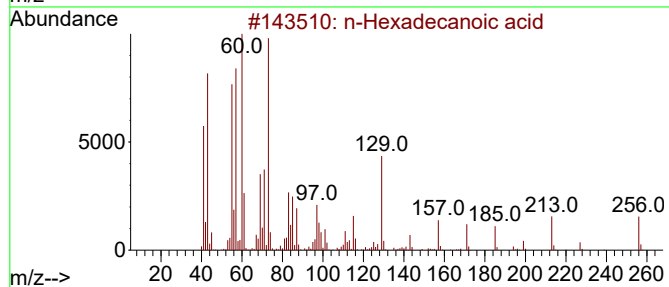
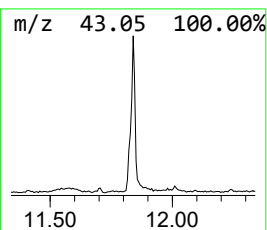
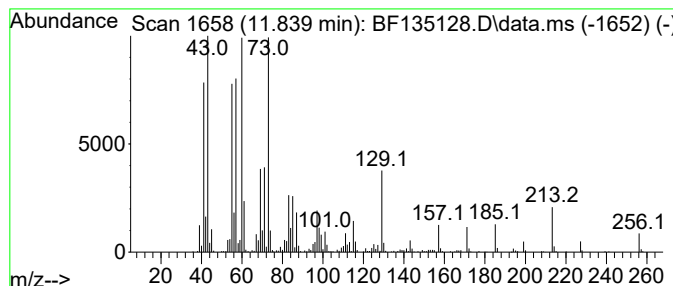
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.839	21.08 ng	915765	Phenanthrene-d10	11.298

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	91
3			Tridecanoic acid	214	C13H26O2	000638-53-9	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
Data File : BF135128.D
Acq On : 07 Sep 2023 11:35
Operator : CG\JU
Sample : 04283-01
Misc :
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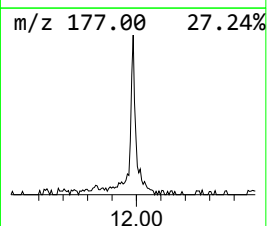
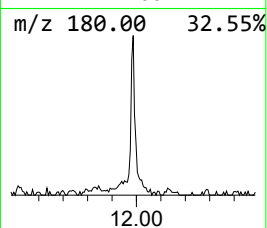
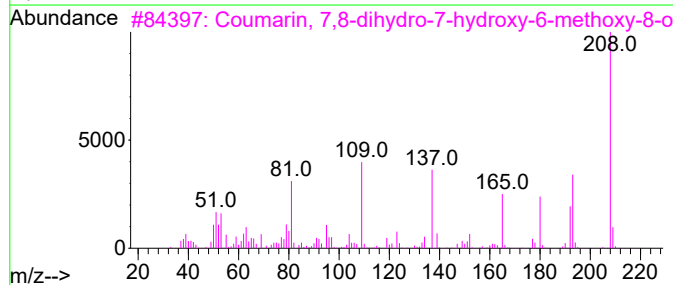
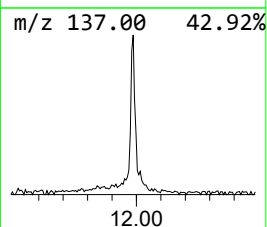
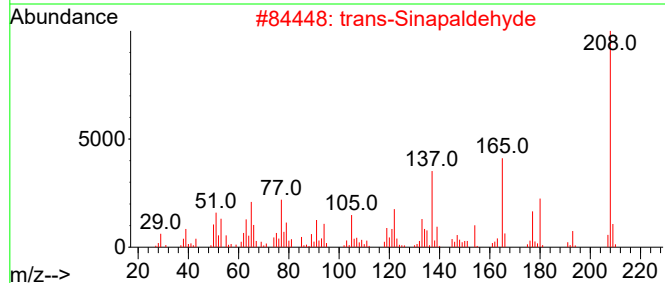
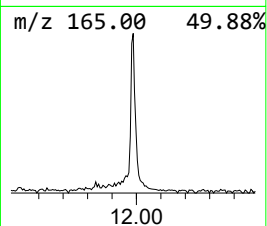
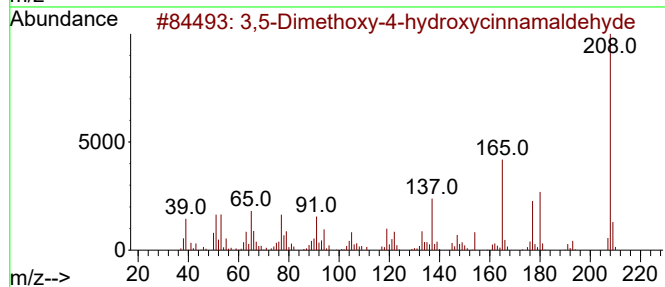
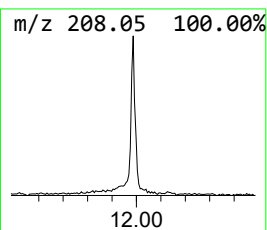
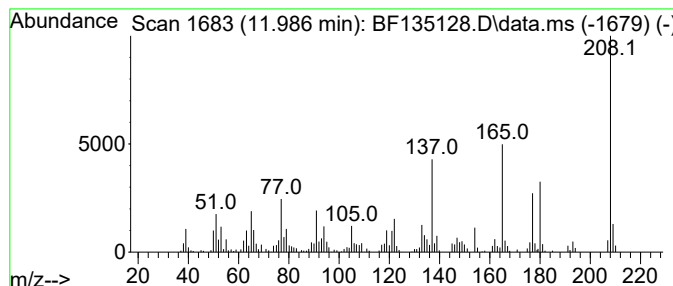
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.986	5.46 ng	237444	Phenanthrene-d10	11.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Dimethoxy-4-hydroxycinnamal... 208	C11H12O4	087345-53-7	98	
2			trans-Sinapaldehyde 208	C11H12O4	004206-58-0	93	
3			Coumarin, 7,8-dihydro-7-hydroxy-... 208	C10H8O5	1000263-13-6	58	
4			2H-1-Benzopyran-2-one, 7,8-dihyd... 208	C10H8O5	000574-84-5	52	
5			2,5-Dimethoxycinnamic acid 208	C11H12O4	010538-51-9	46	



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
Data File : BF135128.D
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Sample : 04283-01
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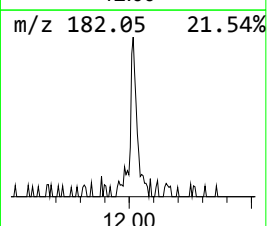
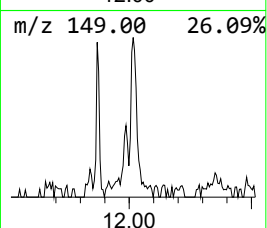
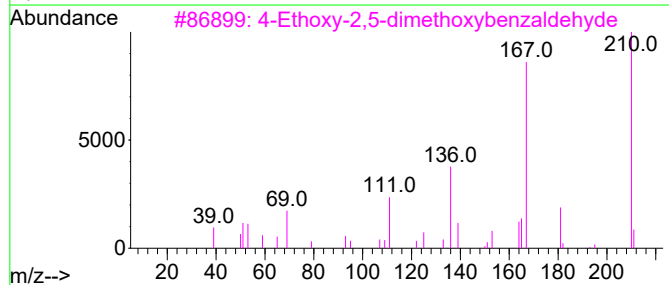
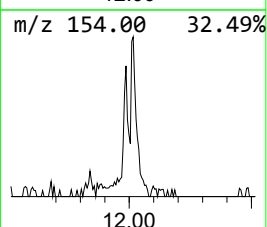
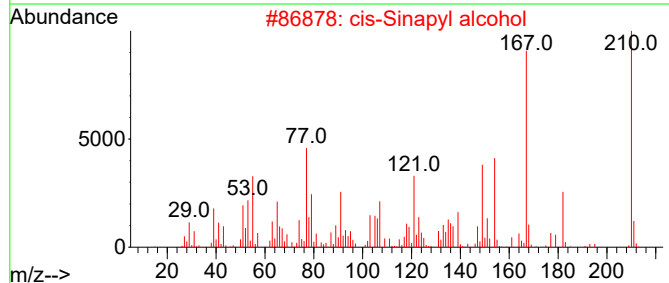
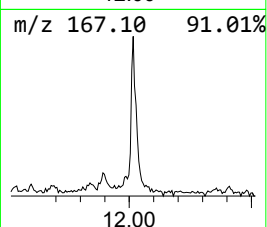
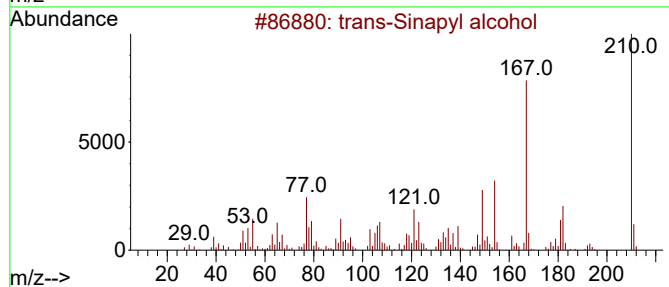
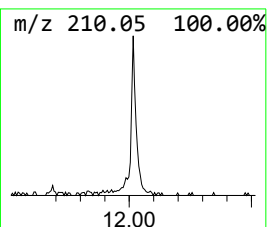
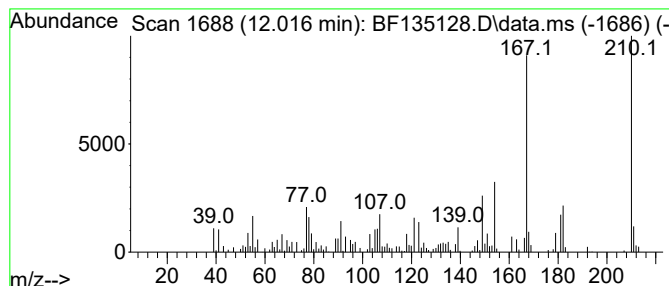
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 trans-Sinapyl alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.016	3.82 ng	166113	Phenanthrene-d10	11.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Sinapyl alcohol	210	C11H14O4	020675-96-1	95
2			cis-Sinapyl alcohol	210	C11H14O4	104330-63-4	91
3			4-Ethoxy-2,5-dimethoxybenzaldehyde	210	C11H14O4	1000121-30-2	53
4			2-Butenal, 3-methyl-, dibutylhyd...	210	C13H26N2	075268-13-2	50
5			4'-(.beta.-Chloroethyl)acetophenone	182	C10H11ClO	069614-95-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
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Sample : 04283-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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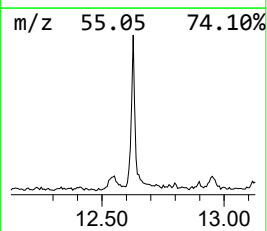
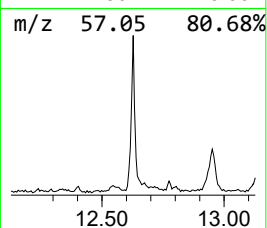
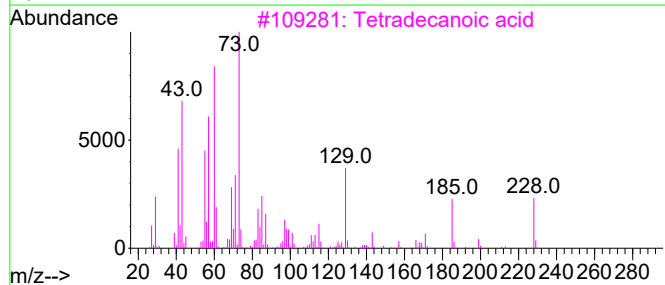
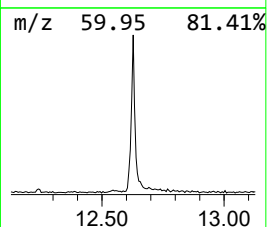
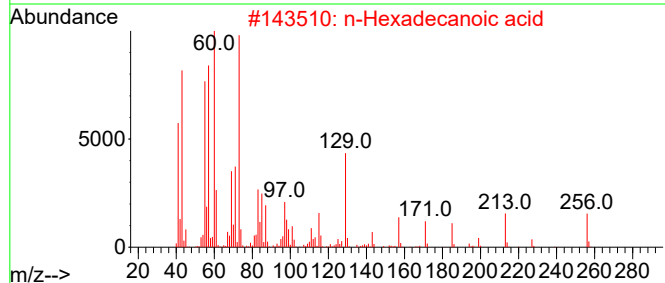
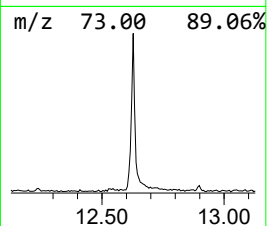
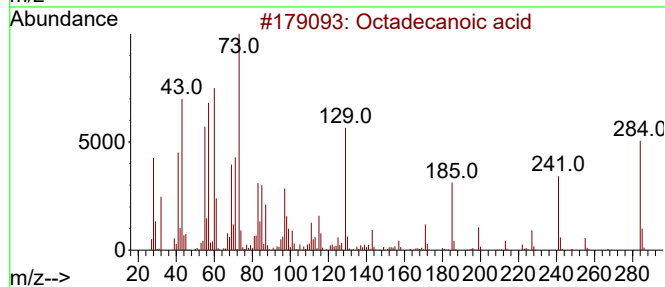
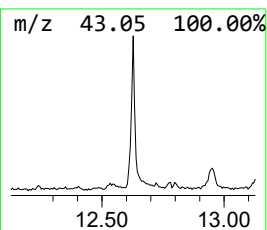
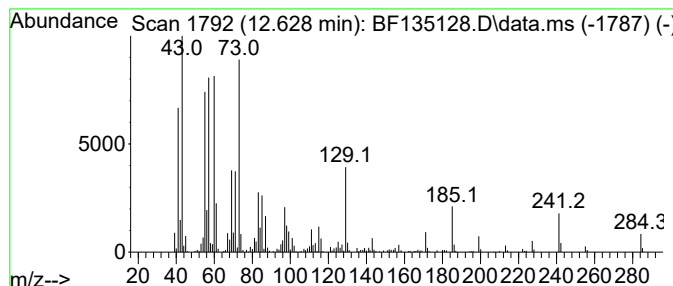
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.628	10.80 ng	469474	Phenanthrene-d10	11.298

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	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72



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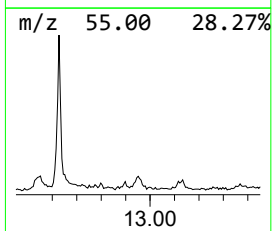
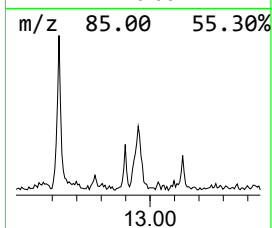
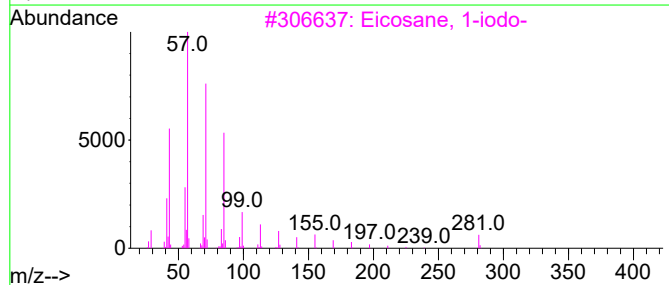
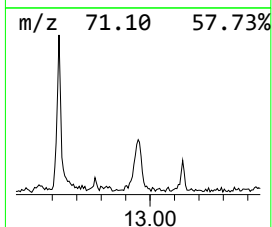
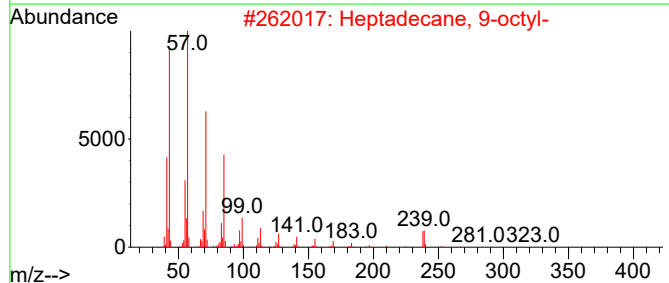
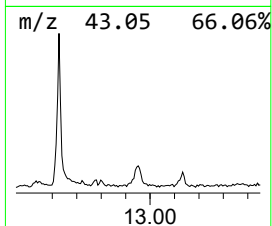
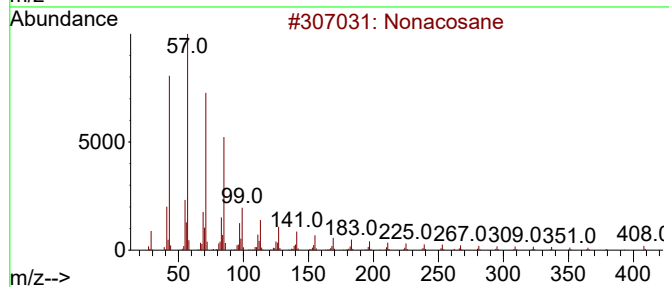
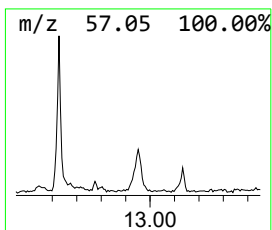
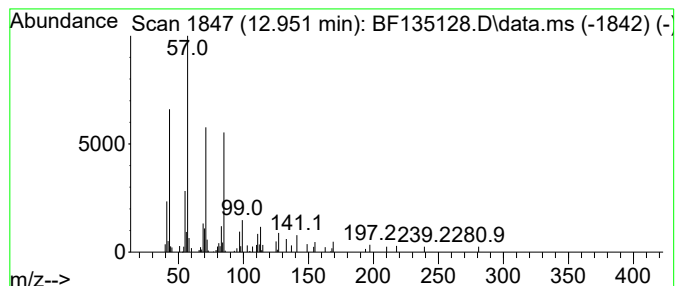
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Nonacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.951	2.81 ng	81713	Chrysene-d12		13.957	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonacosane		408	C29H60	000630-03-5	90
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	86
3	Eicosane, 1-iodo-		408	C20H41I	1000406-31-8	86
4	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	80
5	Octacosane, 2-methyl-		408	C29H60	001560-98-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
Data File : BF135128.D
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Sample : 04283-01
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Instrument :
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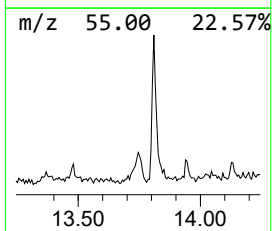
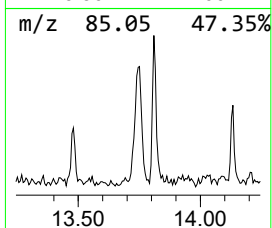
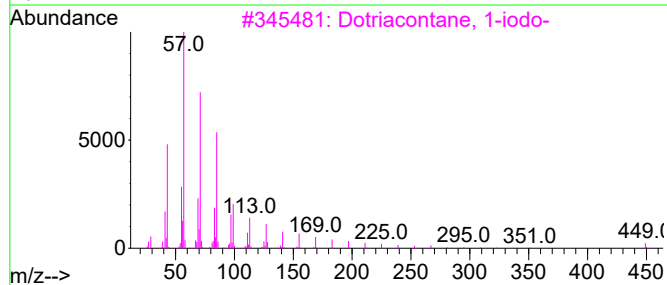
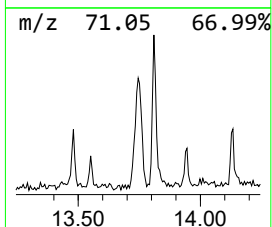
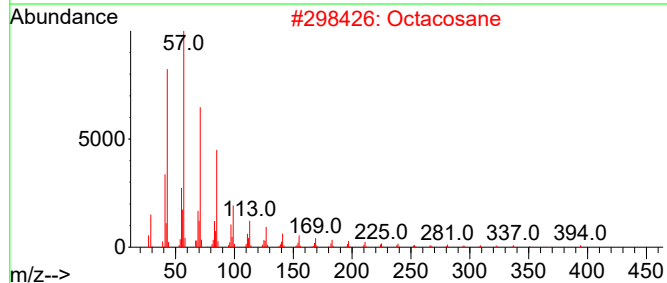
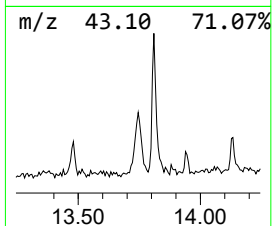
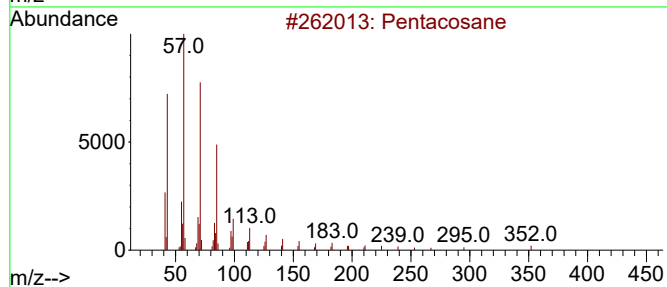
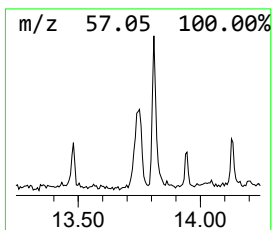
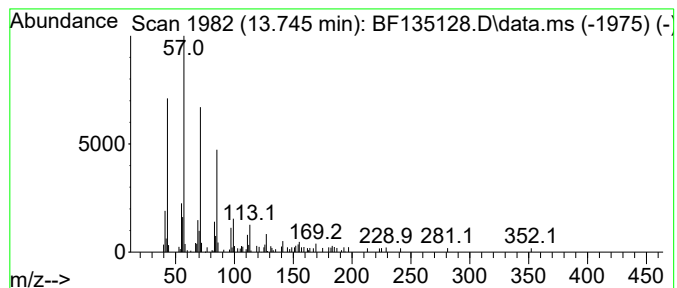
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Pentacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.745	3.44 ng	99934	Chrysene-d12	13.957

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentacosane	352	C25H52	000629-99-2	95
2		Octacosane	394	C28H58	000630-02-4	91
3		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Hentriacontane	437	C31H64	000630-04-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
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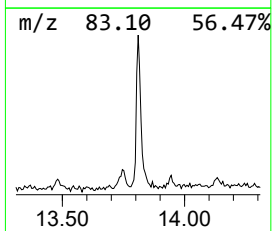
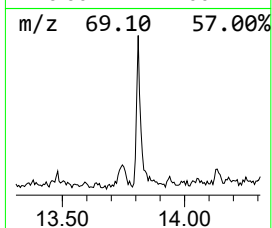
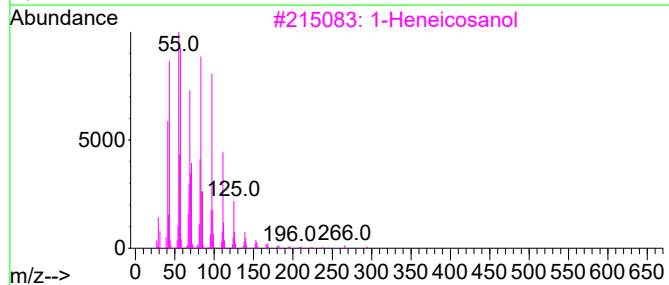
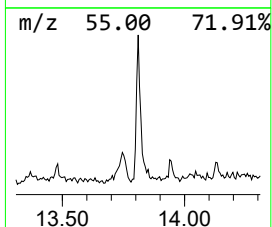
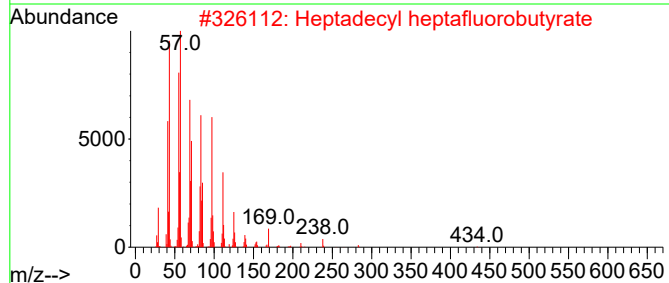
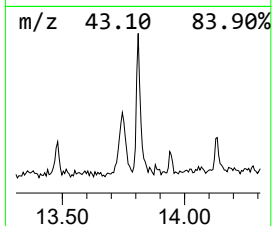
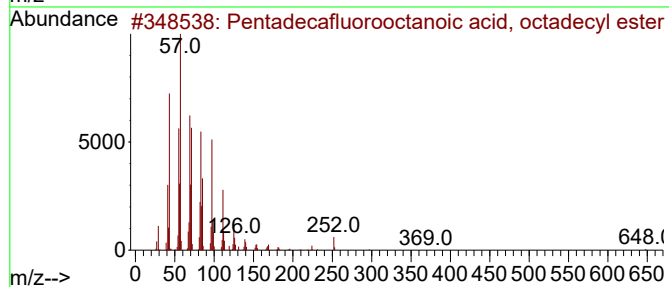
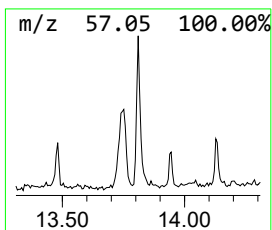
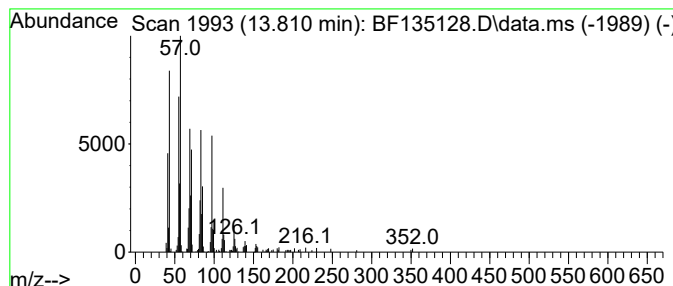
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentadecafluorooctanoic aci... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	6.73 ng	195730	Chrysene-d12	13.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			1-Heneicosanol	312	C21H44O	015594-90-8	93
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
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ALS Vial : 4 Sample Multiplier: 1

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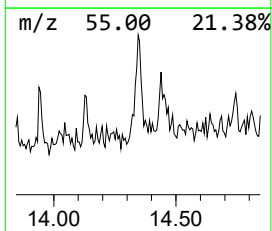
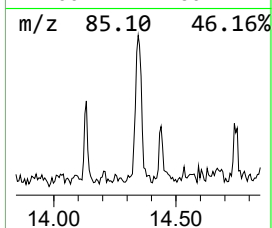
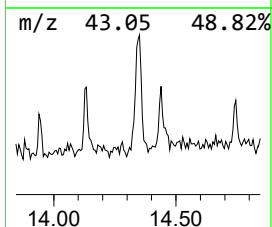
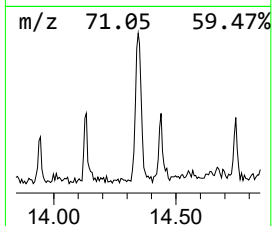
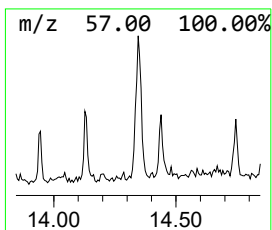
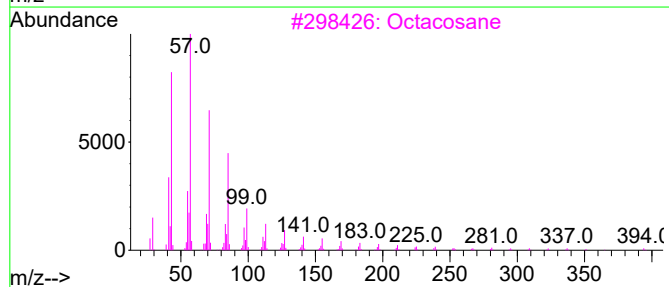
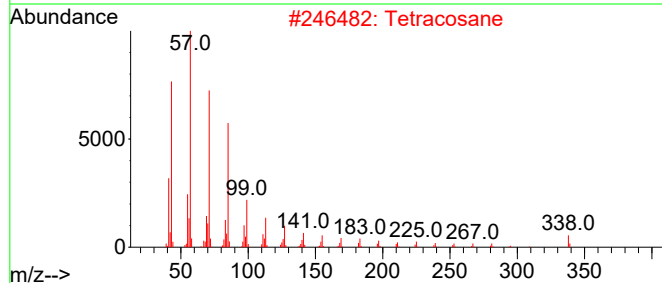
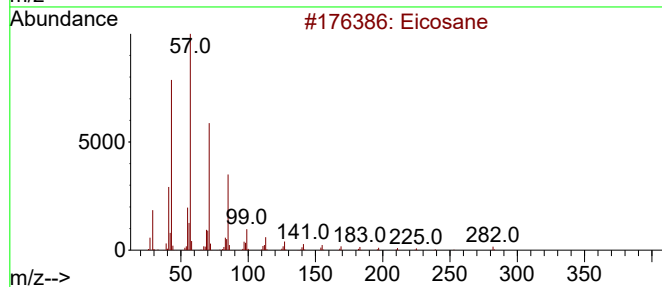
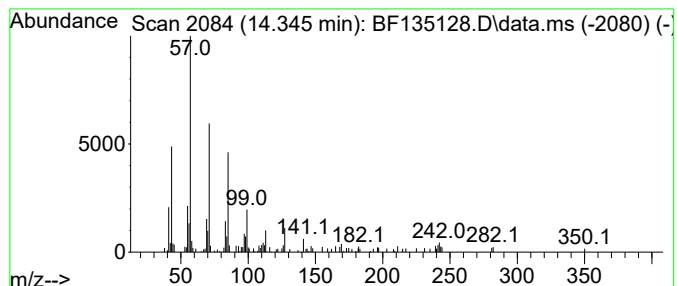
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.345	3.10 ng	90230	Chrysene-d12	13.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	91
2	Tetracosane			338	C24H50	000646-31-1	91
3	Octacosane			394	C28H58	000630-02-4	90
4	Docosane, 1-iodo-			436	C22H45I	1000406-31-9	87
5	Tetracosane, 1-iodo-			464	C24H49I	1000406-32-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
Data File : BF135128.D
Acq On : 07 Sep 2023 11:35
Operator : CG\JU
Sample : 04283-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-1

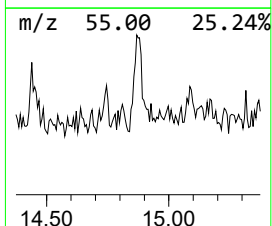
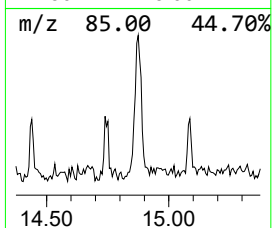
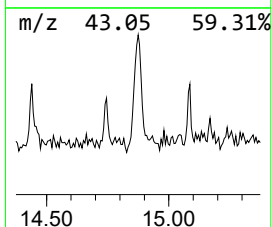
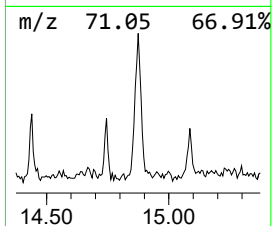
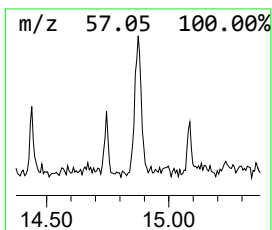
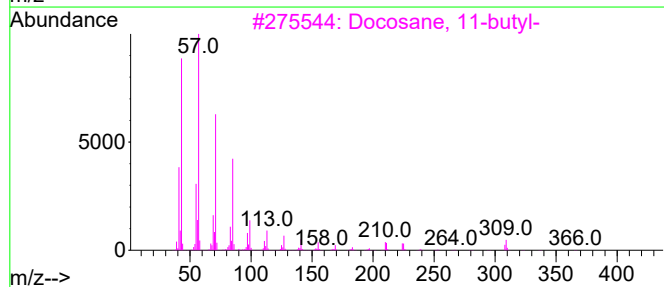
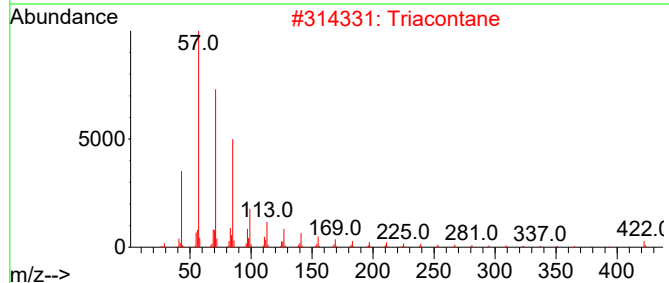
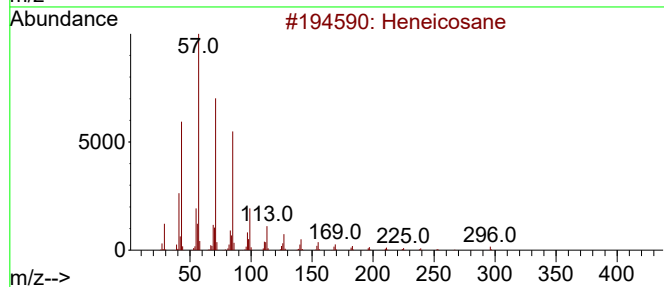
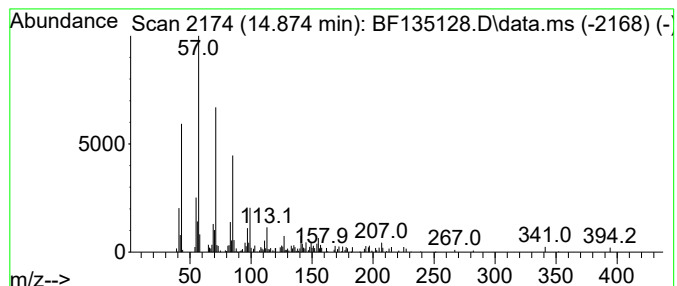
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.874	3.74 ng	106059	Perylene-d12	15.427

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			Docosane, 11-butyl-	366	C26H54	013475-76-8	87
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
5			Nonadecane, 2-methyl-	282	C20H42	001560-86-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
Data File : BF135128.D
Acq On : 07 Sep 2023 11:35
Operator : CG\JU
Sample : 04283-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-1

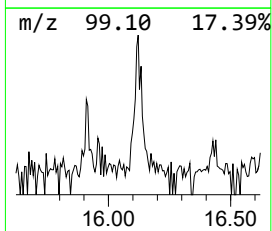
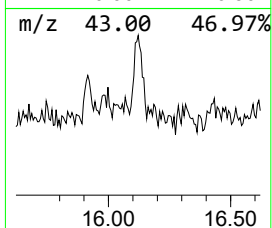
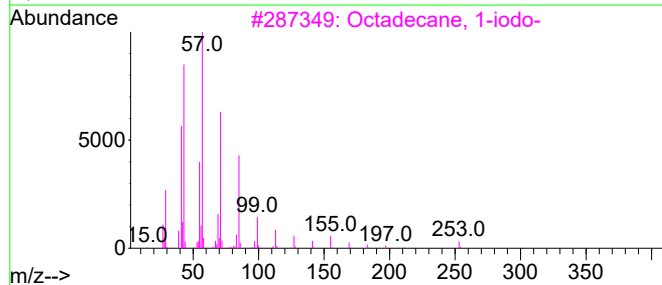
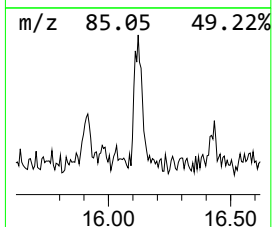
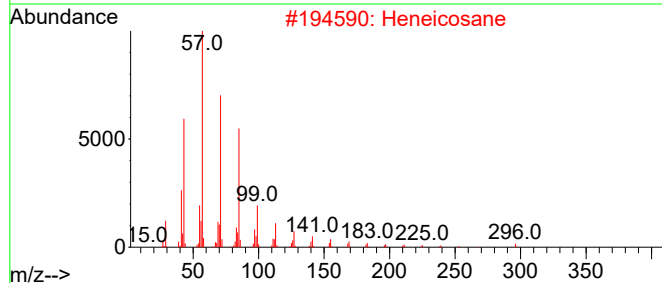
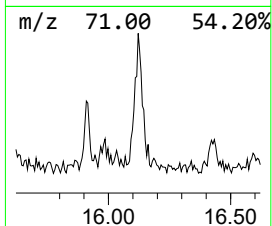
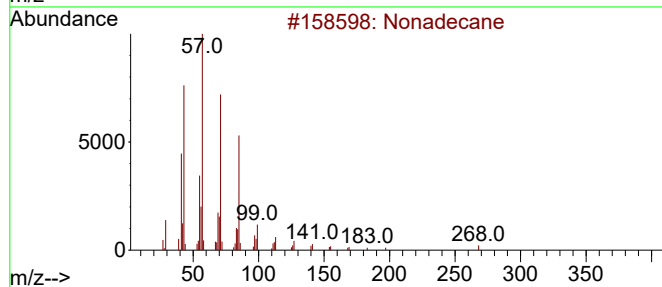
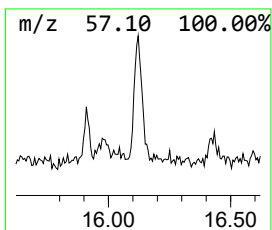
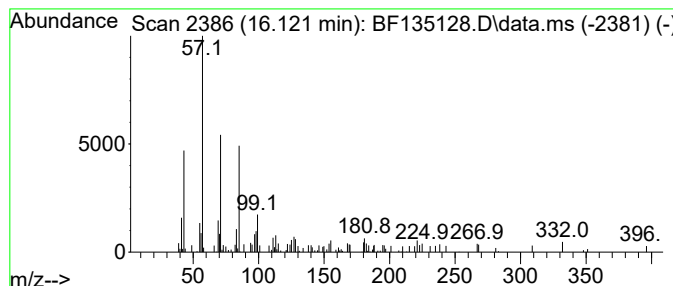
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.121	2.51 ng	71168	Perylene-d12	15.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	80
2			Heneicosane	296	C21H44	000629-94-7	64
3			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	64
4			Pentacosane	352	C25H52	000629-99-2	64
5			4-Methylheneicosane	310	C22H46	025117-29-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090723\
Data File : BF135128.D
Acq On : 07 Sep 2023 11:35
Operator : CG\JU
Sample : 04283-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.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
Propane, 1,1-di...	2.134	2.0	ng	57898	1	6.769	573087	20.0
2-Pentanone, 4-...	5.004	3.2	ng	90406	1	6.769	573087	20.0
Ethanol, 2-(2-e...	6.604	3.0	ng	85012	1	6.769	573087	20.0
unknown10.951	10.951	2.5	ng	107141	4	11.298	869015	20.0
n-Hexadecanoic ...	11.839	21.1	ng	915765	4	11.298	869015	20.0
3,5-Dimethoxy-4...	11.986	5.5	ng	237444	4	11.298	869015	20.0
trans-Sinapyl a...	12.016	3.8	ng	166113	4	11.298	869015	20.0
Octadecanoic acid	12.628	10.8	ng	469474	4	11.298	869015	20.0
Nonacosane	12.951	2.8	ng	81713	5	13.957	581408	20.0
Pentacosane	13.745	3.4	ng	99934	5	13.957	581408	20.0
Pentadecafluoro...	13.810	6.7	ng	195730	5	13.957	581408	20.0
Eicosane	14.345	3.1	ng	90230	5	13.957	581408	20.0
Heneicosane	14.874	3.7	ng	106059	6	15.427	566530	20.0
Nonadecane	16.121	2.5	ng	71168	6	15.427	566530	20.0