

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072125\
 Data File : BF143174.D
 Acq On : 21 Jul 2025 16:28
 Operator : RC/JU
 Sample : Q2651-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH 8-7

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-BF071725.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143174.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.310	12	20	27	rBV	676568	1059032	38.51%	5.319%
2	5.193	504	510	519	rBV	183250	264866	9.63%	1.330%
3	5.587	571	577	583	rBV	1734909	2300539	83.66%	11.554%
4	6.581	739	746	751	rBV	1772277	2393874	87.06%	12.023%
5	6.963	805	811	816	rVB	714409	870319	31.65%	4.371%
6	7.516	899	905	910	rBV	1190573	1510708	54.94%	7.587%
7	8.239	1023	1028	1033	rBV	855787	1096426	39.87%	5.506%
8	9.316	1205	1211	1216	rBV	2219842	2749730	100.00%	13.810%
9	9.998	1321	1327	1332	rBV	865546	1056588	38.43%	5.306%
10	10.704	1442	1447	1452	rBV	188320	226147	8.22%	1.136%
11	10.780	1455	1460	1470	rBV	1004163	1260943	45.86%	6.333%
12	11.016	1496	1500	1504	rVB	24871	30465	1.11%	0.153%
13	11.480	1574	1579	1585	rBV	651303	813800	29.60%	4.087%
14	11.704	1612	1617	1625	rBV2	16859	29064	1.06%	0.146%
15	12.004	1662	1668	1686	rBV	350911	530640	19.30%	2.665%
16	12.786	1797	1801	1809	rBV	62750	95324	3.47%	0.479%
17	13.063	1843	1848	1854	rBV	1489264	1943462	70.68%	9.761%
18	13.957	1996	2000	2011	rBV	160642	298153	10.84%	1.497%
19	14.121	2023	2028	2035	rBV	576376	745973	27.13%	3.746%
20	14.210	2039	2043	2052	rVV	45449	74333	2.70%	0.373%
21	14.986	2172	2175	2180	rVB	50228	64710	2.35%	0.325%
22	15.621	2278	2283	2289	rBV	308793	496393	18.05%	2.493%

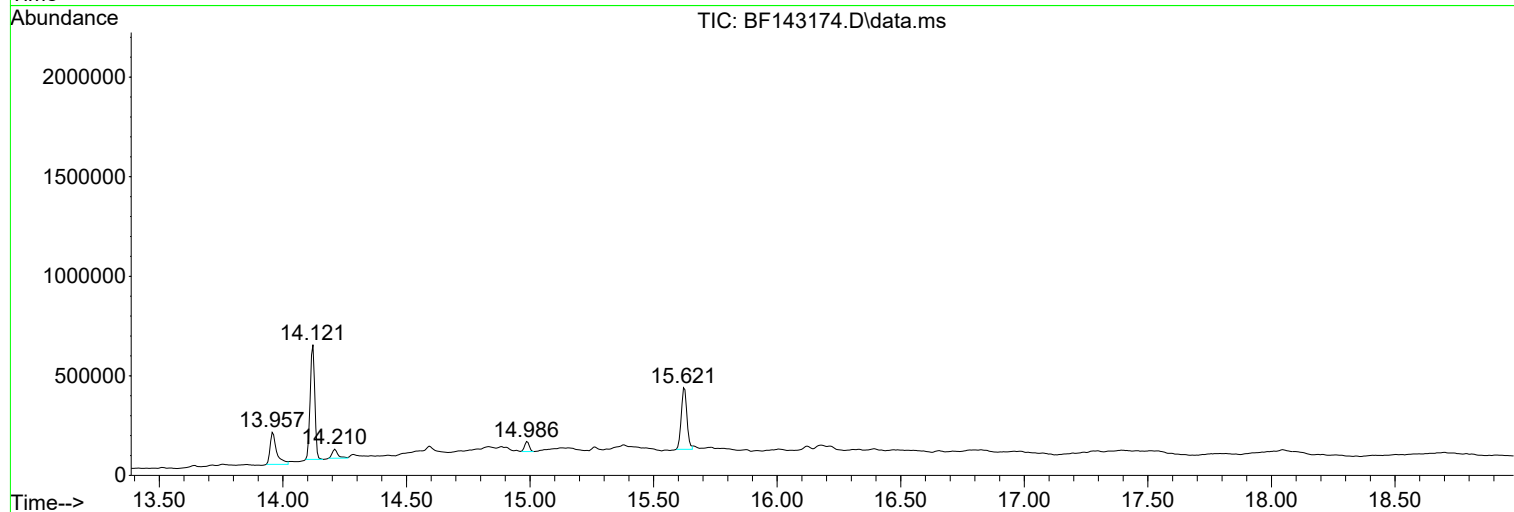
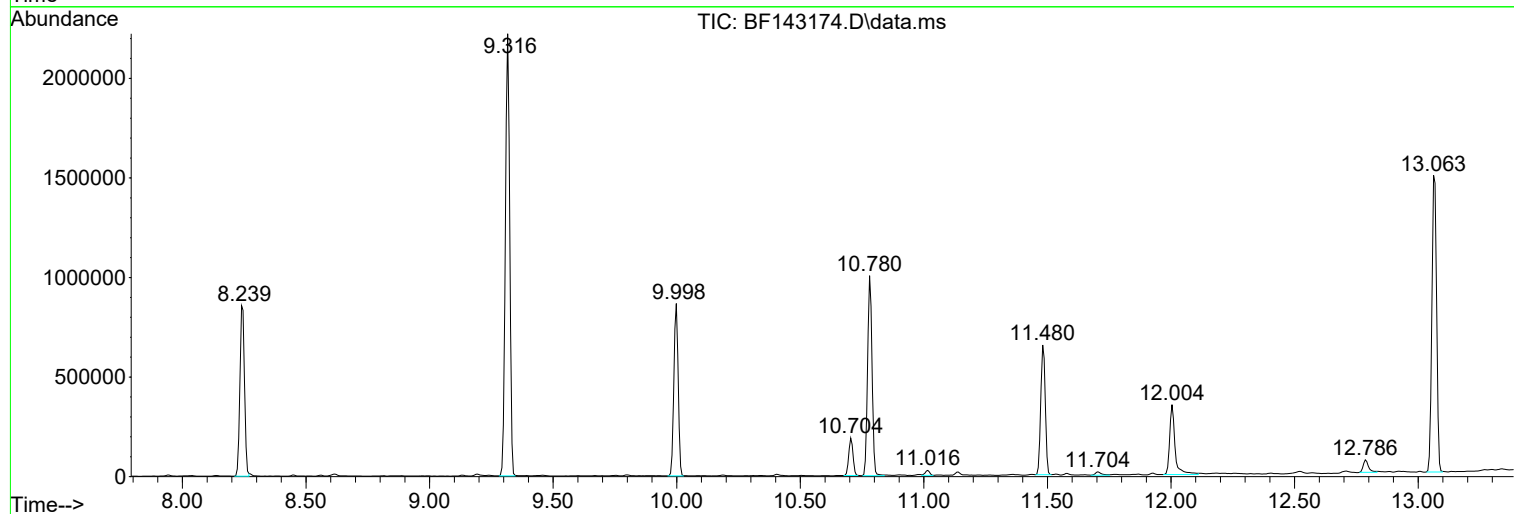
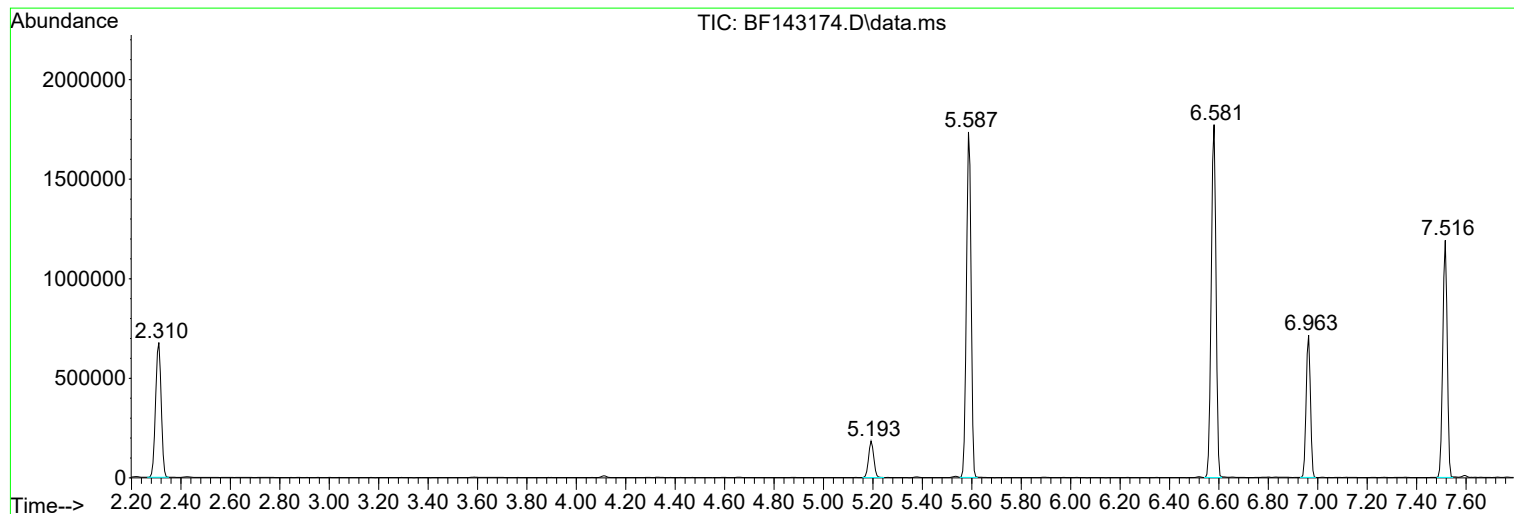
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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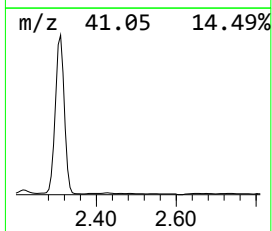
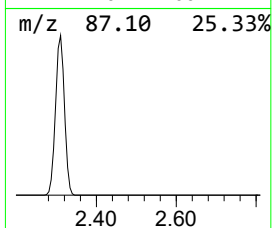
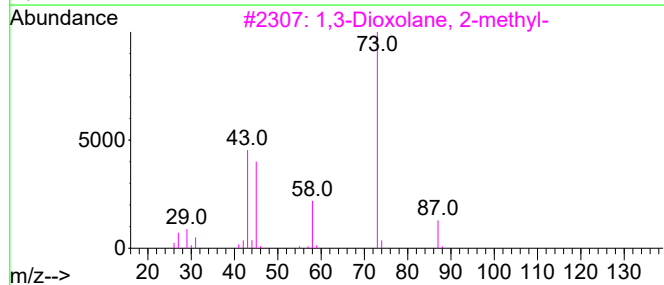
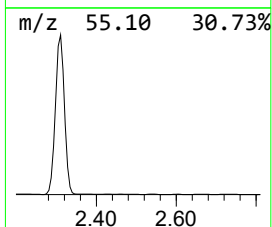
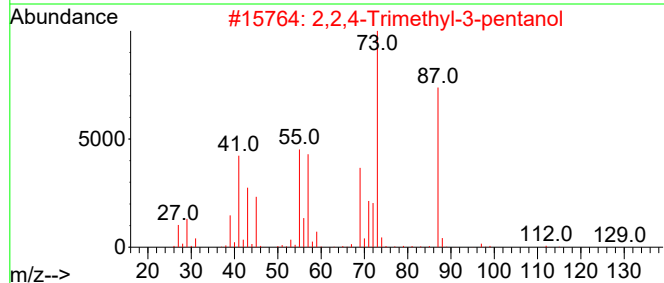
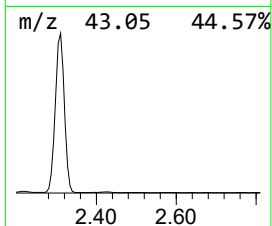
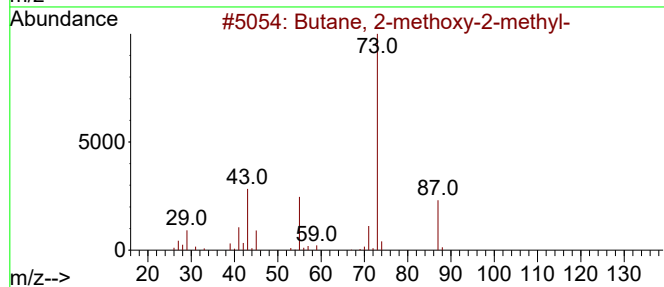
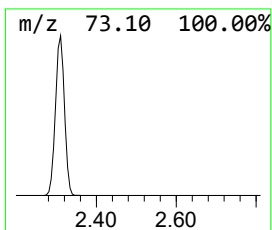
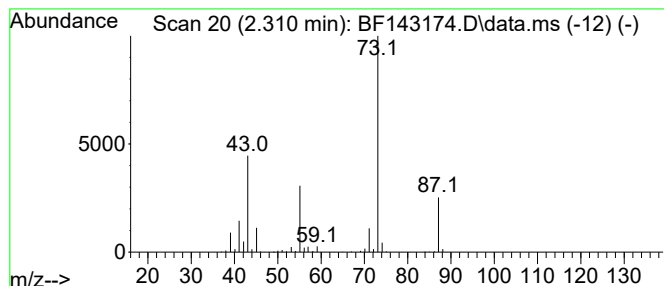
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.310	24.34 ng	1059030	1,4-Dichlorobenzene-d4	6.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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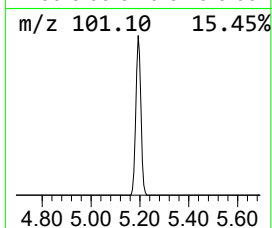
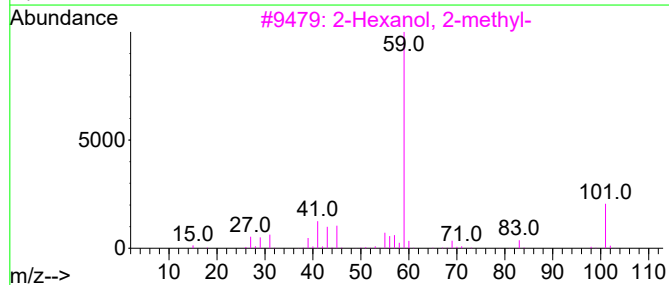
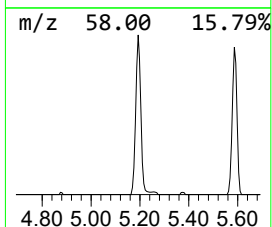
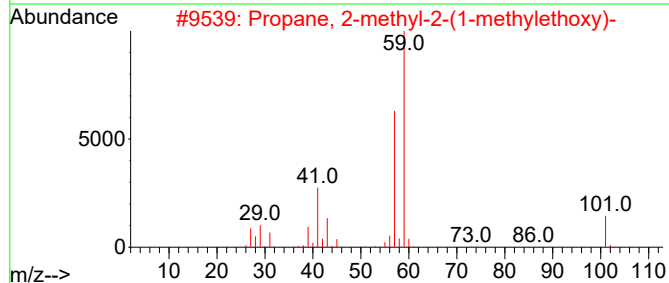
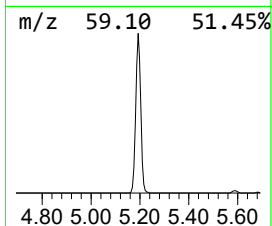
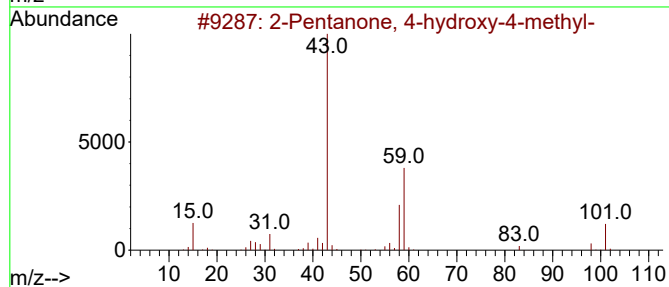
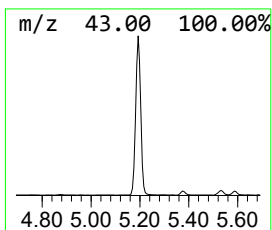
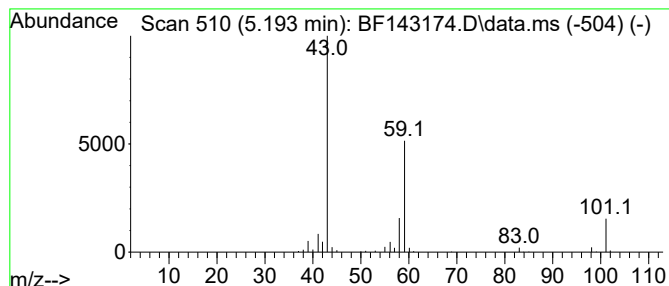
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.193	6.09 ng	264866	1,4-Dichlorobenzene-d4			6.963
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	59
2	Propane, 2-methyl-2-(1-methyleth...		116	C7H16O	017348-59-3	36
3	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	33
4	3-Hexanol, 4-methyl-		116	C7H16O	000615-29-2	33
5	Acetic acid, 1,1-dimethylethyl e...		116	C6H12O2	000540-88-5	28



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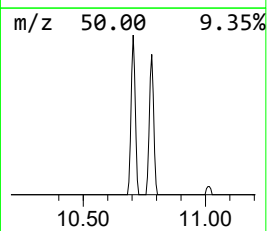
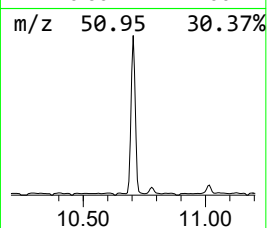
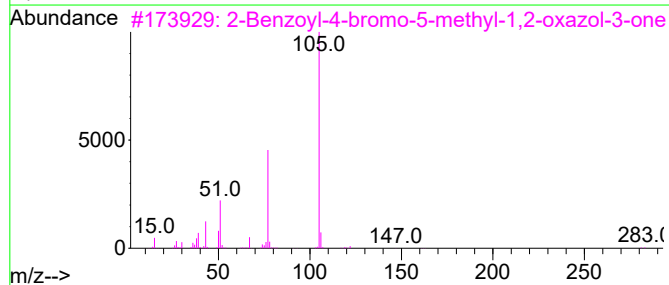
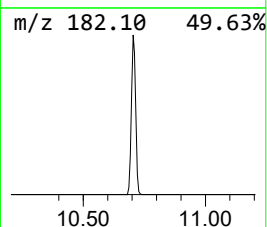
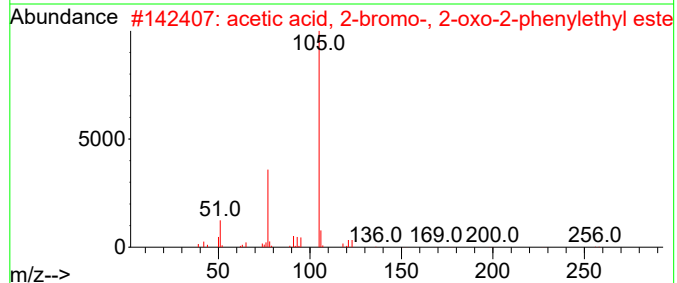
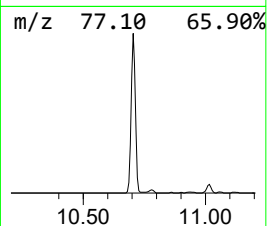
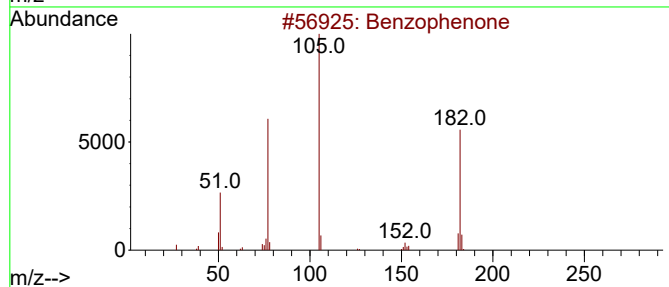
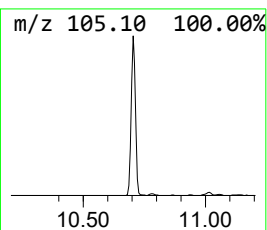
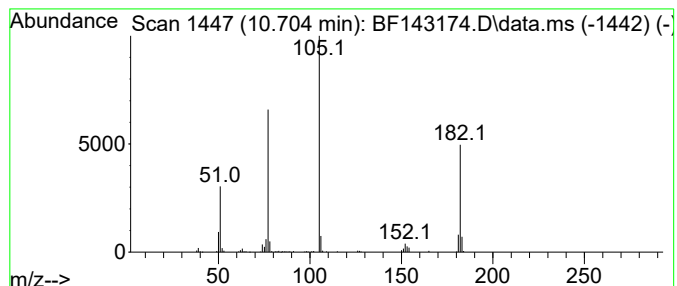
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.704	4.28 ng	226147	Acenaphthene-d10	9.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	50
3			2-Benzoyl-4-bromo-5-methyl-1,2-o...	281	C11H8BrNO3	1000444-92-3	50
4			1-Butanone, 1-phenyl-	148	C10H12O	000495-40-9	50
5			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	50



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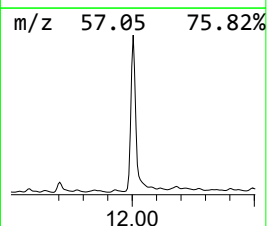
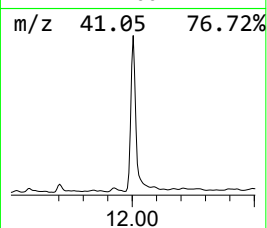
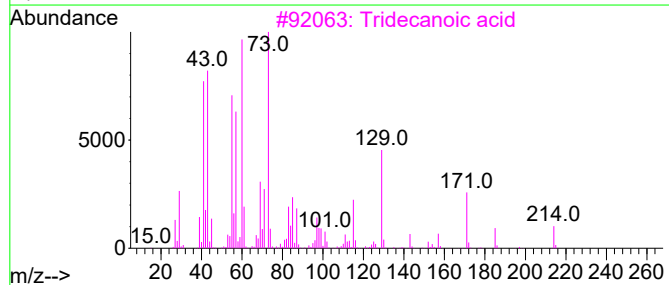
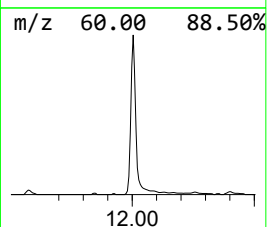
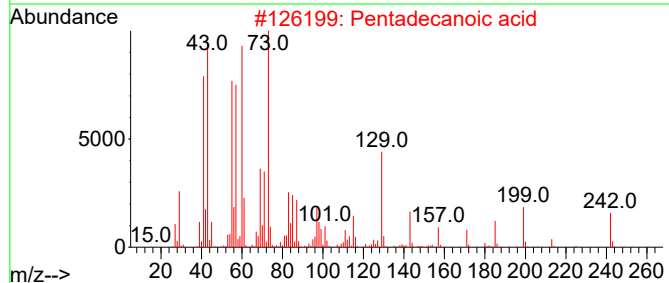
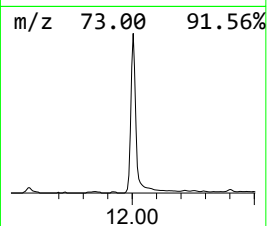
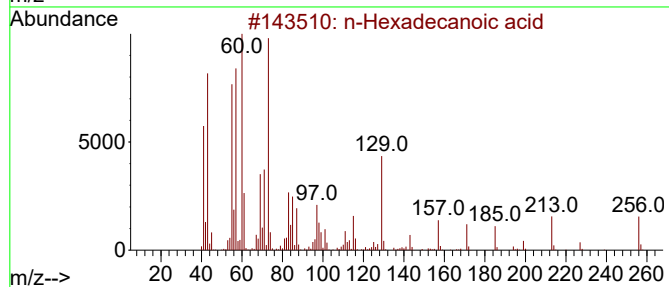
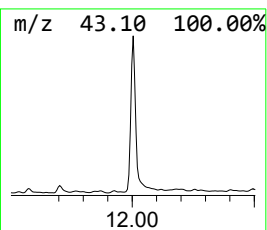
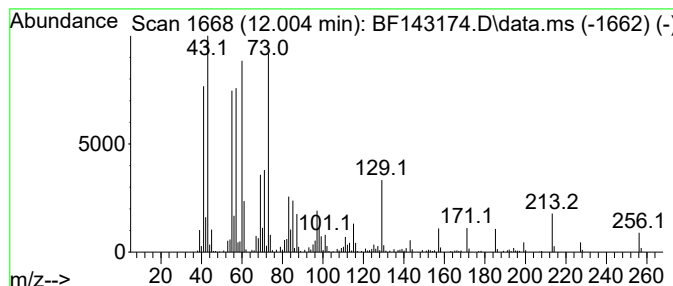
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	13.04 ng	530640	Phenanthrene-d10	11.480

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	90
3			Tridecanoic acid	214	C13H26O2	000638-53-9	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	76
5			Undecanoic acid	186	C11H22O2	000112-37-8	58



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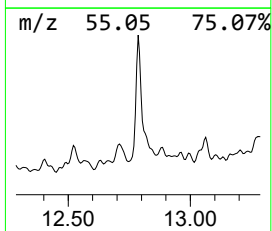
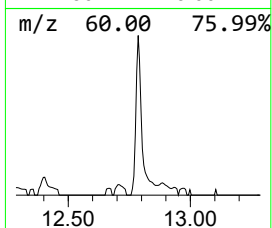
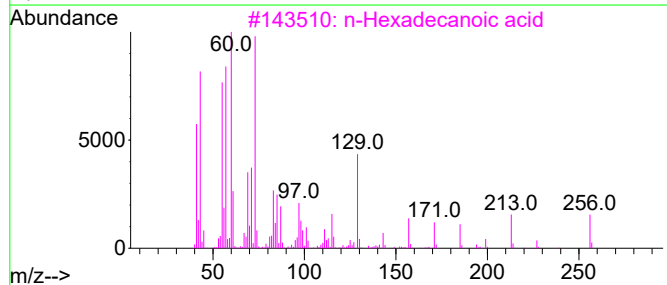
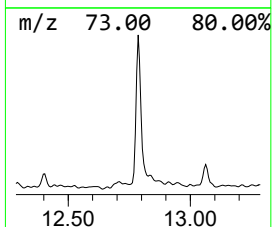
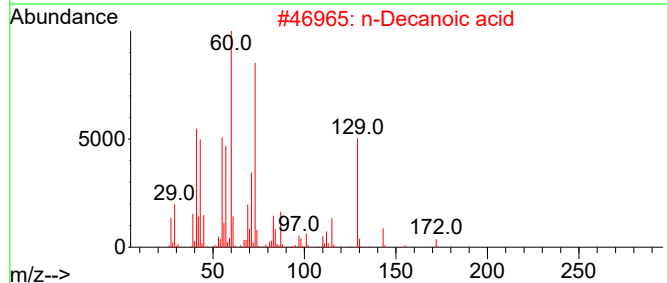
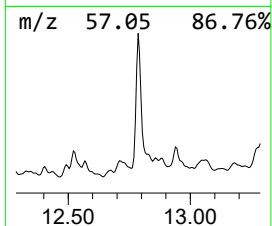
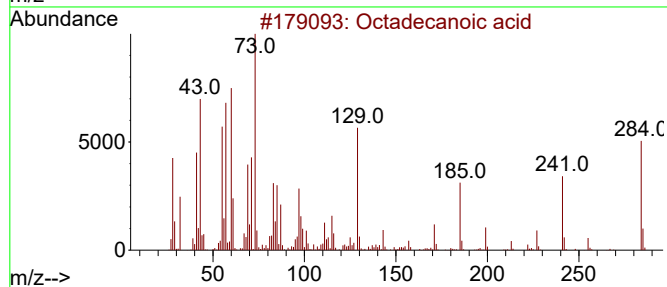
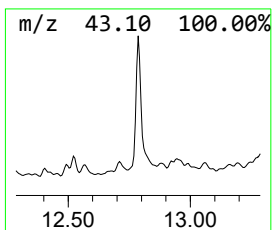
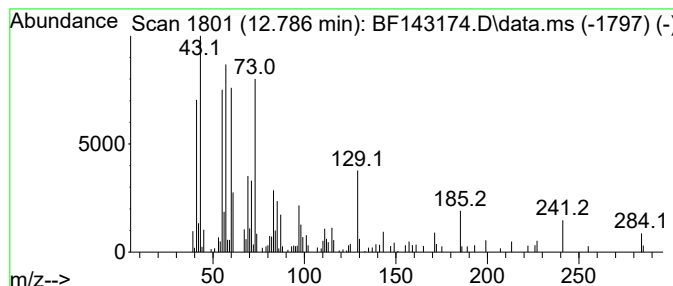
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.786	2.34 ng	95324	Phenanthrene-d10	11.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Decanoic acid	172	C10H20O2	000334-48-5	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	74



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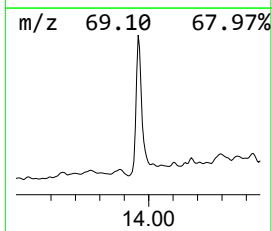
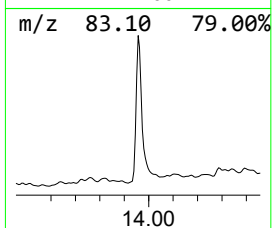
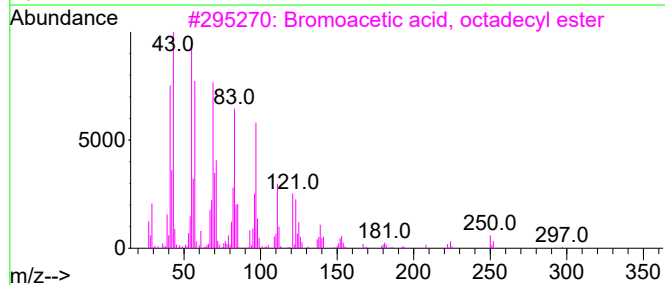
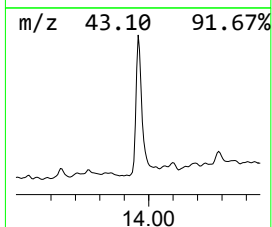
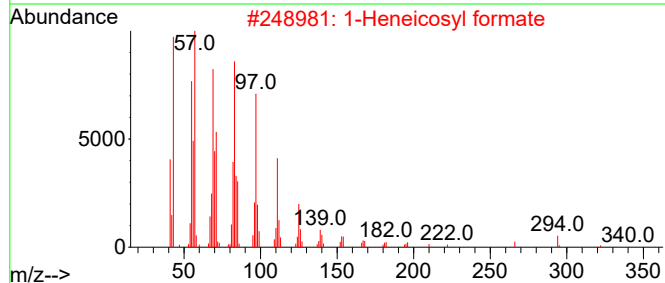
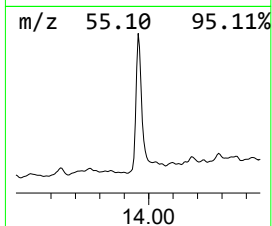
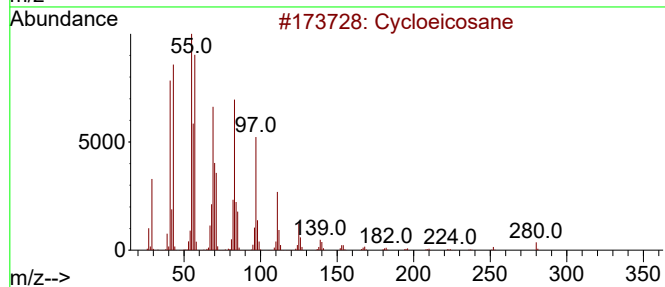
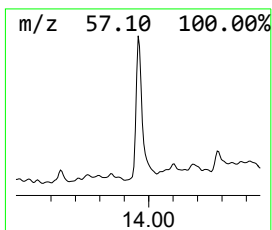
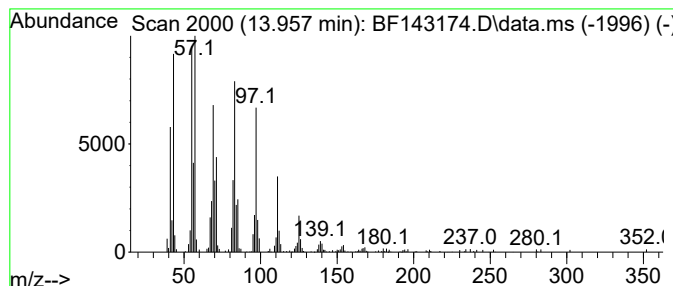
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cycloeicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.957	7.99 ng	298153	Chrysene-d12	14.121

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloeicosane	280	C20H40	000296-56-0	97
2			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	94
4			Octacosanol	410	C28H58O	000557-61-9	94
5			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072125\
Data File : BF143174.D
Acq On : 21 Jul 2025 16:28
Operator : RC/JU
Sample : Q2651-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH 8-7

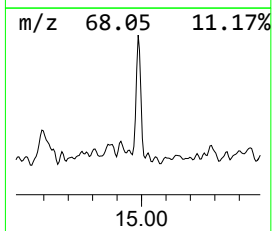
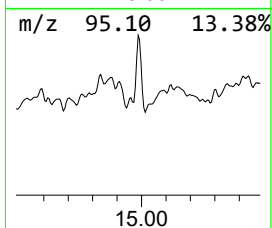
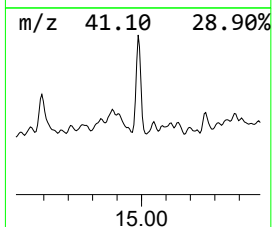
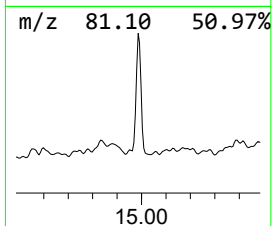
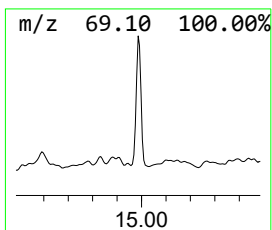
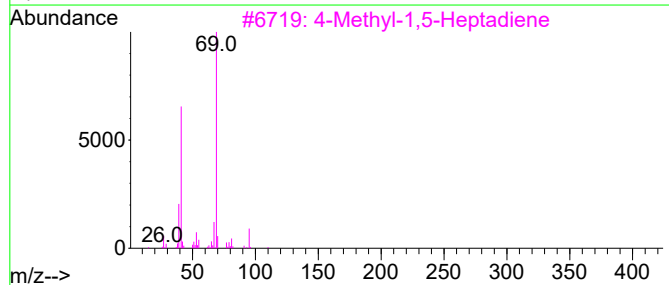
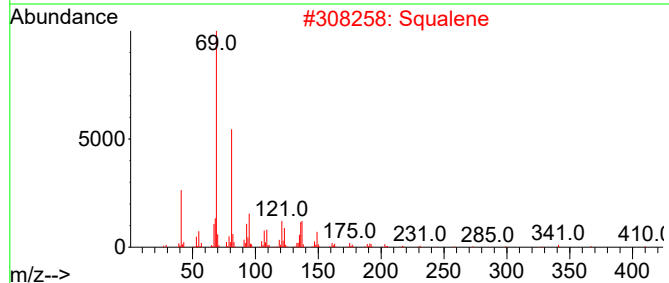
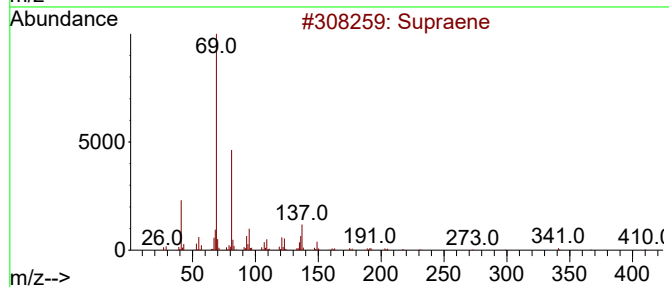
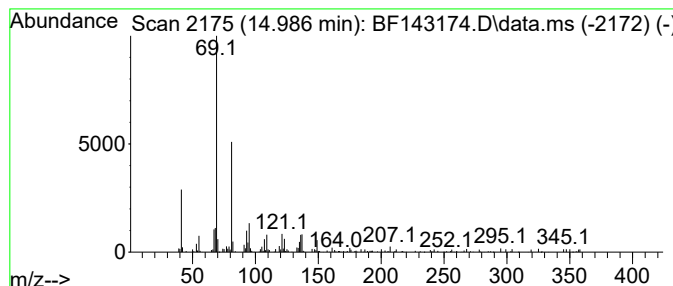
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.986	2.61 ng	64710	Perylene-d12	15.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	74
2			Squalene	410	C30H50	000111-02-4	72
3			4-Methyl-1,5-Heptadiene	110	C8H14	000998-94-7	64
4			5-Bromo-1-pentanol	166	C5H11BrO	034626-51-2	59
5			trans-Farnesol	222	C15H26O	000106-28-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072125\
Data File : BF143174.D
Acq On : 21 Jul 2025 16:28
Operator : RC/JU
Sample : Q2651-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH 8-7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.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
Butane, 2-metho...	2.310	24.3	ng	1059030	1	6.963	870319	20.0
2-Pentanone, 4-...	5.193	6.1	ng	264866	1	6.963	870319	20.0
Benzophenone	10.704	4.3	ng	226147	3	9.998	1056590	20.0
n-Hexadecanoic ...	12.004	13.0	ng	530640	4	11.480	813800	20.0
Octadecanoic acid	12.786	2.3	ng	95324	4	11.480	813800	20.0
Cycloeicosane	13.957	8.0	ng	298153	5	14.121	745973	20.0
Supraene	14.986	2.6	ng	64710	6	15.621	496393	20.0