

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
 Data File : BF131095.D
 Acq On : 09 Nov 2022 19:32
 Operator : CG\JU
 Sample : N5437-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 596C-3

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-BF110922.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131095.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.181	10	16	21	rVB	1915623	2381068	75.50%	11.965%
2	4.463	398	404	406	rBV	50725	54909	1.74%	0.276%
3	4.493	406	409	416	rVB	87779	93606	2.97%	0.470%
4	4.904	475	479	485	rBV	42143	42588	1.35%	0.214%
5	5.110	509	514	521	rVB	167835	188895	5.99%	0.949%
6	5.504	577	581	592	rBV	1404577	1262780	40.04%	6.345%
7	6.510	748	752	763	rVB	855555	780816	24.76%	3.924%
8	6.893	813	817	820	rBV	644621	586034	18.58%	2.945%
9	7.457	907	913	916	rBV	1894446	1889341	59.91%	9.494%
10	8.175	1030	1035	1038	rBV	916723	754945	23.94%	3.794%
11	9.251	1212	1218	1221	rBV	3420346	3153655	100.00%	15.847%
12	9.934	1328	1334	1337	rBV	1079577	867051	27.49%	4.357%
13	10.128	1361	1367	1372	rBV5	22315	36616	1.16%	0.184%
14	10.728	1463	1469	1472	rBV	2214797	2076555	65.85%	10.435%
15	11.428	1584	1588	1591	rBV	920980	856692	27.17%	4.305%
16	11.945	1670	1676	1681	rBV	164957	183303	5.81%	0.921%
17	12.345	1736	1744	1746	rBV3	44346	93687	2.97%	0.471%
18	12.733	1806	1810	1813	rBV3	26480	35909	1.14%	0.180%
19	12.839	1821	1828	1836	rBV5	42652	109206	3.46%	0.549%
20	13.016	1853	1858	1861	rBV	3210954	2869428	90.99%	14.419%
21	13.633	1956	1963	1970	rBV	58486	120207	3.81%	0.604%
22	13.986	2019	2023	2029	rBV4	32320	64594	2.05%	0.325%
23	14.069	2034	2037	2041	rVV	644088	524308	16.63%	2.635%
24	14.163	2049	2053	2059	rVB	64243	67663	2.15%	0.340%
25	14.239	2060	2066	2073	rVB2	63289	113757	3.61%	0.572%
26	14.757	2150	2154	2160	rBV	59127	93802	2.97%	0.471%
27	15.321	2245	2250	2257	rVB	45366	79499	2.52%	0.399%
28	15.563	2286	2291	2297	rVB	383107	473140	15.00%	2.378%
29	15.968	2356	2360	2367	rBV3	25636	46468	1.47%	0.234%

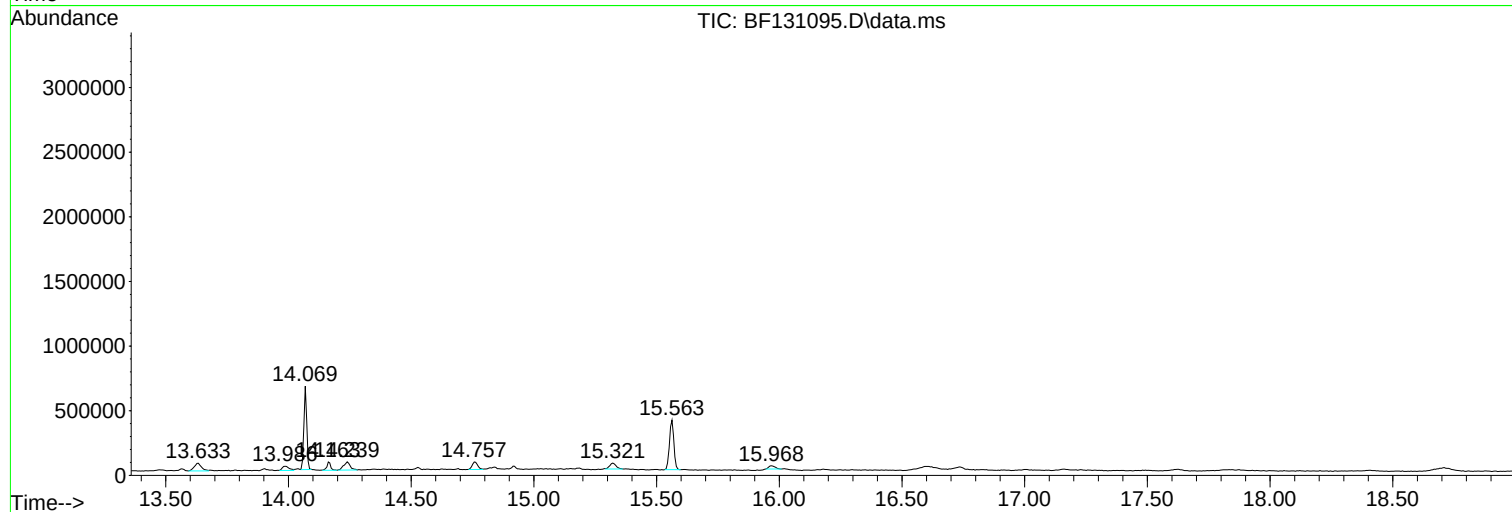
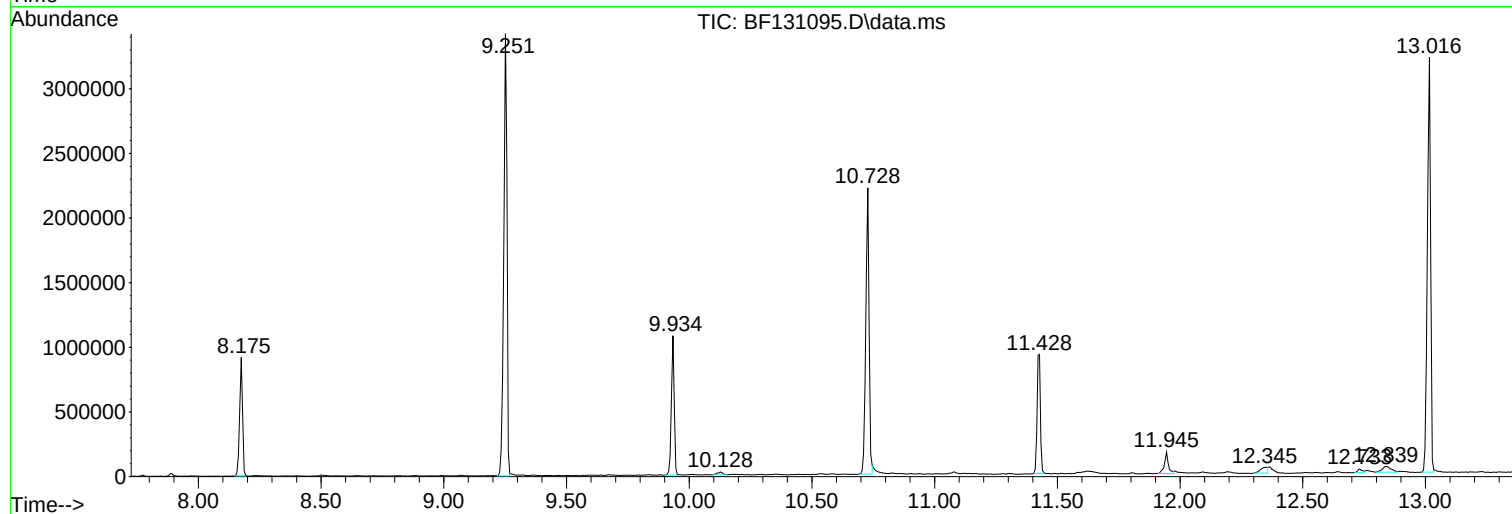
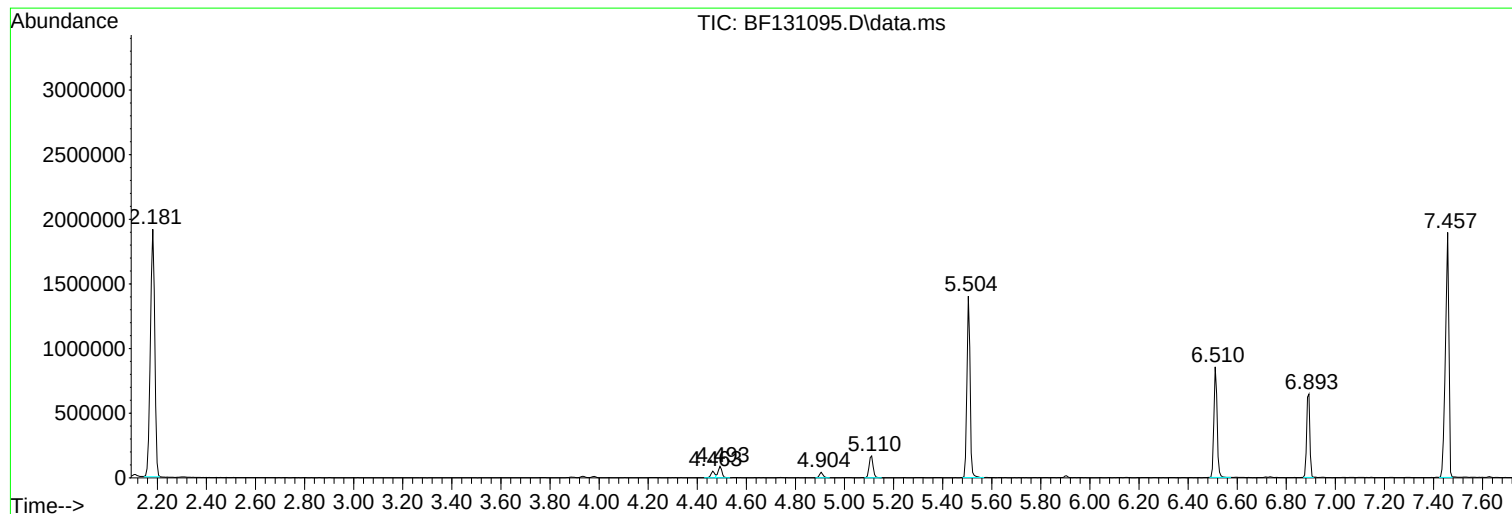
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Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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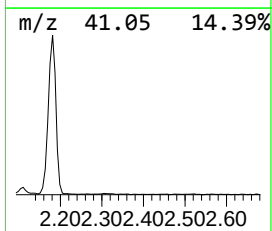
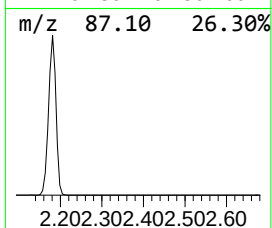
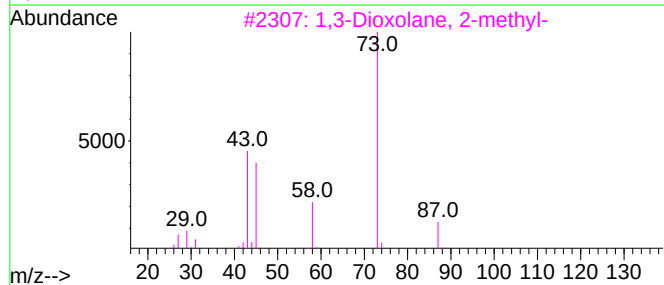
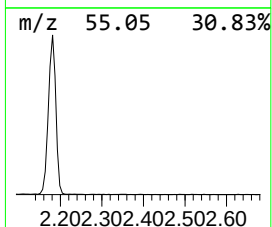
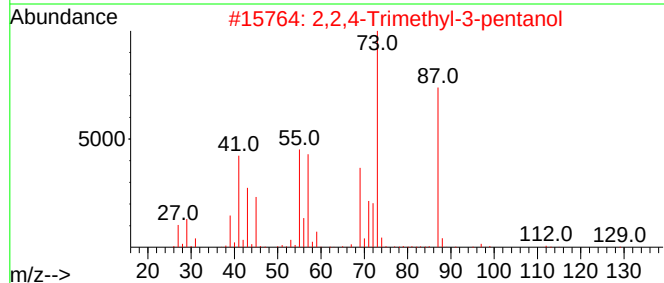
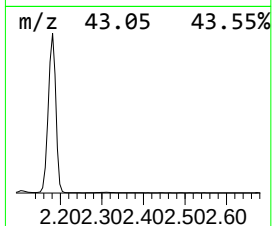
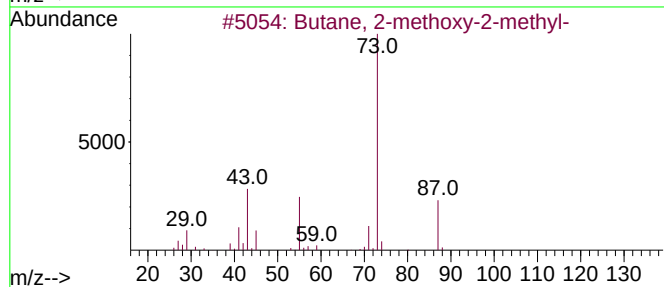
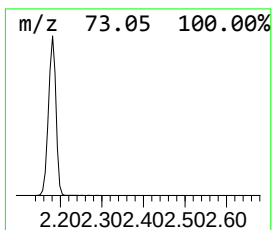
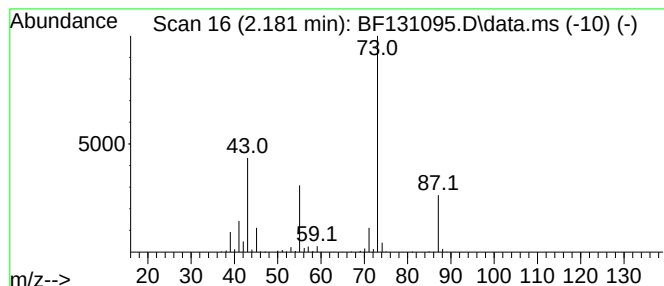
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.181	81.26 ng	2381070	1,4-Dichlorobenzene-d4	6.893

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	39
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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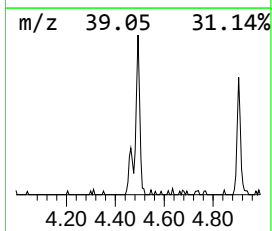
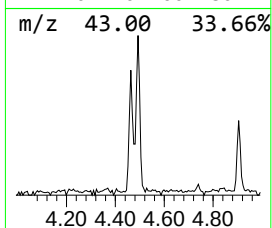
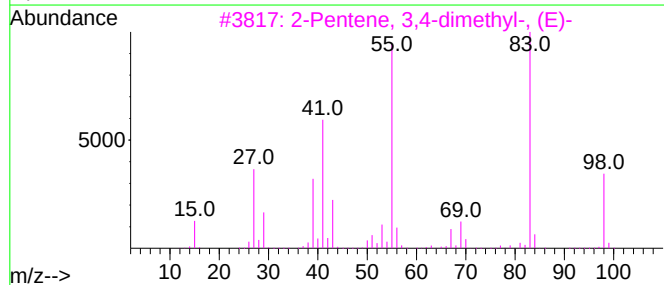
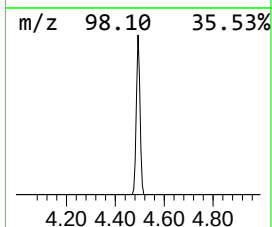
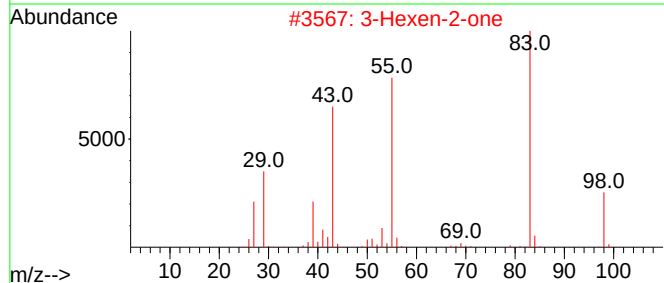
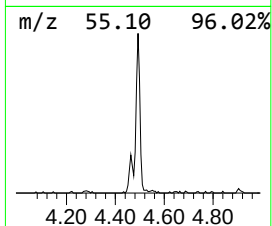
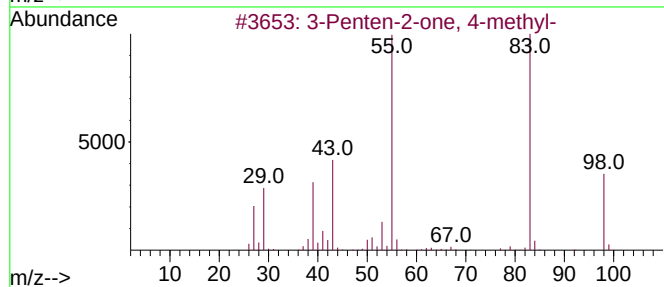
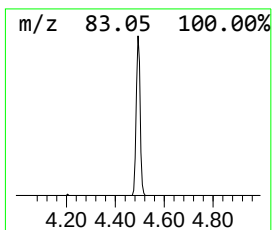
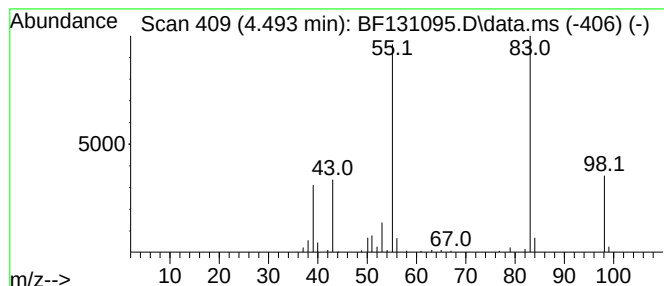
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.493	3.19 ng	93606	1,4-Dichlorobenzene-d4	6.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2			3-Hexen-2-one	98	C6H10O	000763-93-9	91
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	90
5			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86



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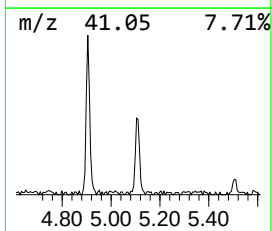
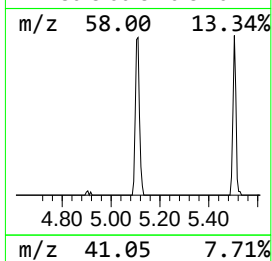
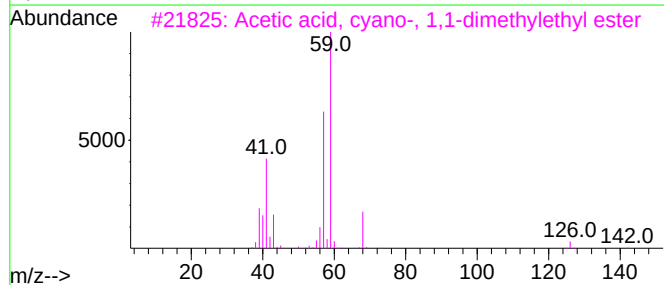
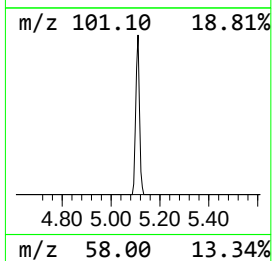
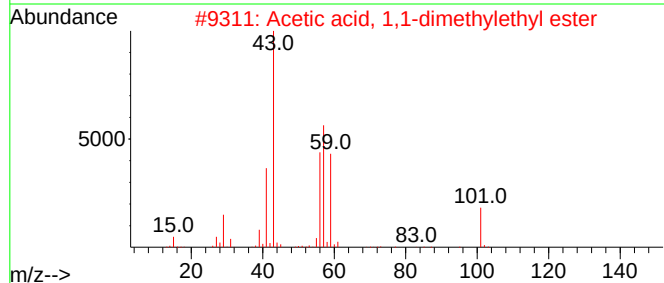
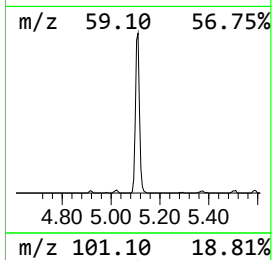
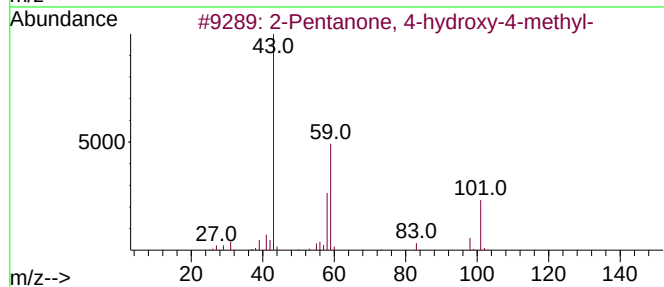
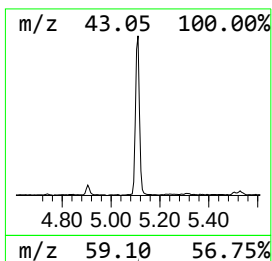
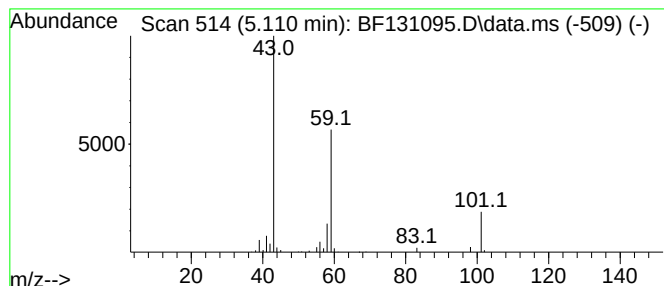
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.110	6.45 ng	188895	1,4-Dichlorobenzene-d4	6.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	37		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		



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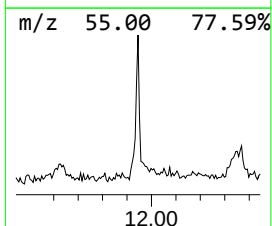
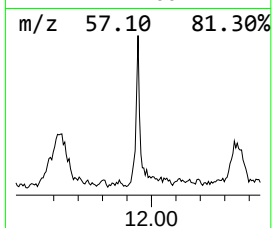
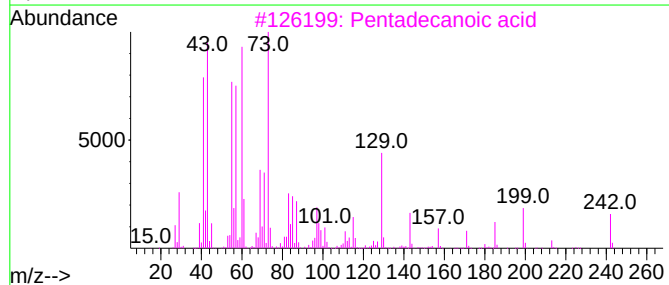
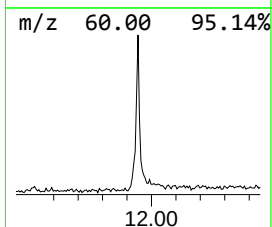
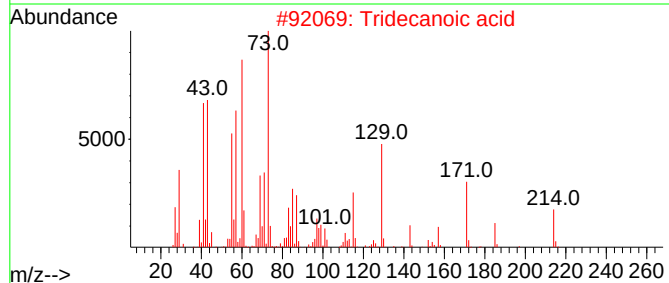
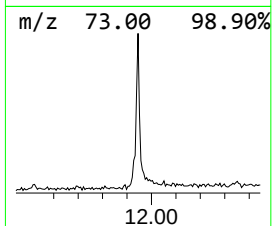
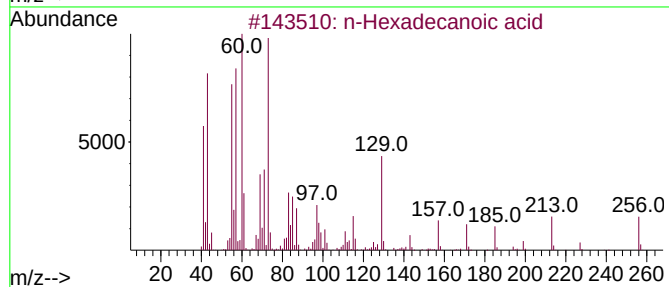
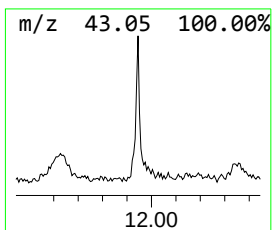
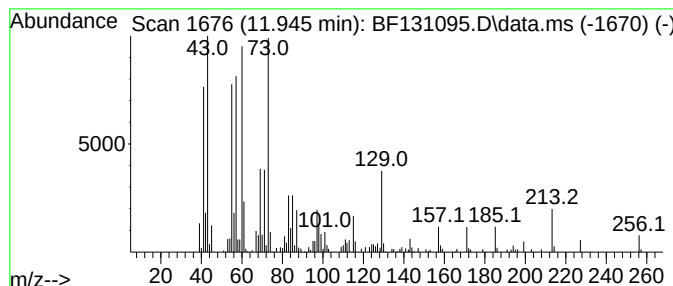
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.945	4.28 ng	183303	Phenanthrene-d10	11.428	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	96
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5	Dodecanoic acid	200	C12H24O2	000143-07-7	72



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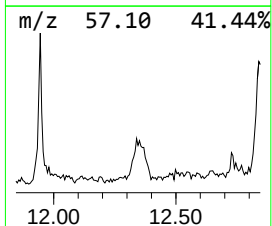
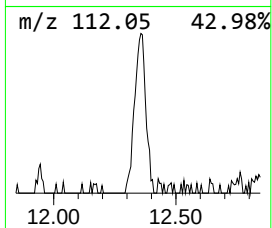
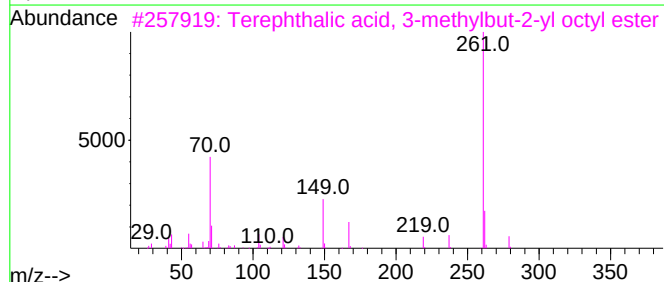
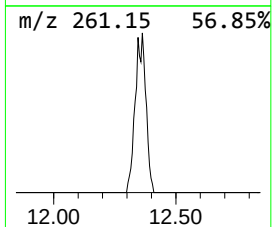
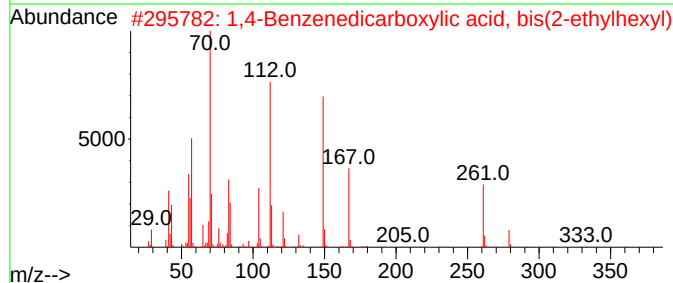
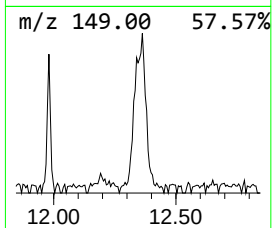
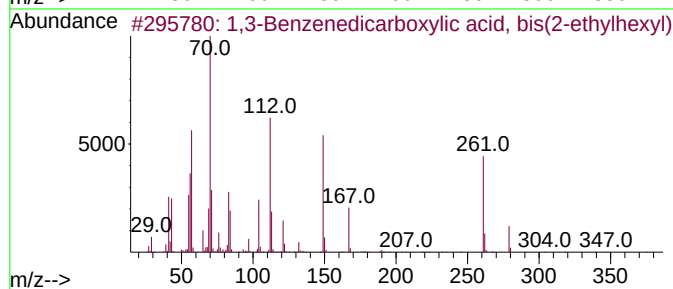
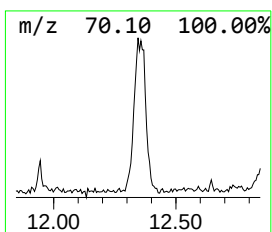
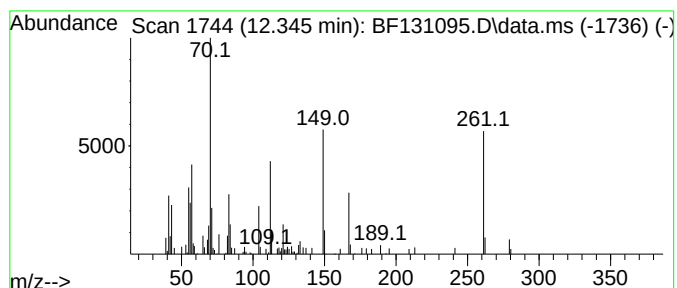
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Peak Number 5 1,3-Benzenedicarboxylic aci... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.345	2.19 ng	93687	Phenanthrene-d10	11.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	53
2			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	49
3			Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	47
4			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	22
5			Terephthalic acid, di(4-octyl) e...	390	C24H38O4	1000323-74-2	22



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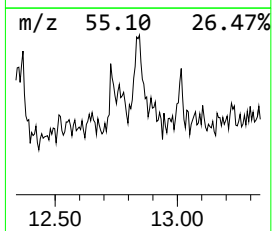
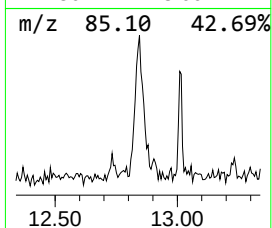
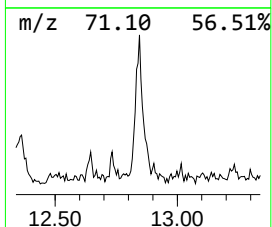
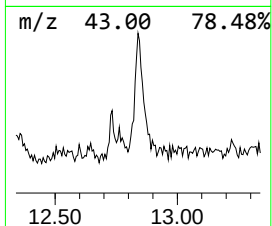
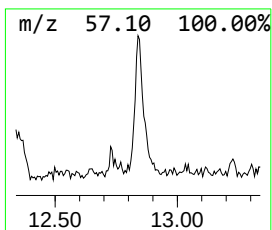
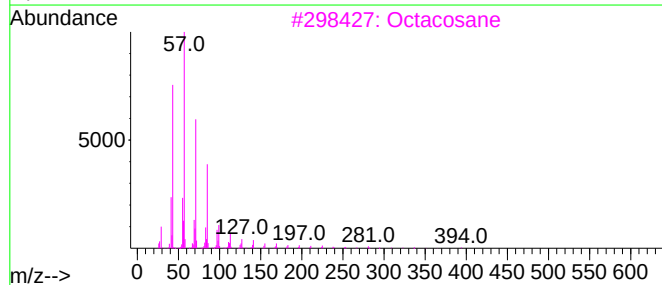
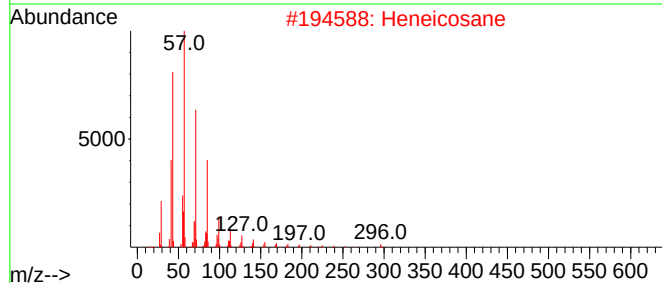
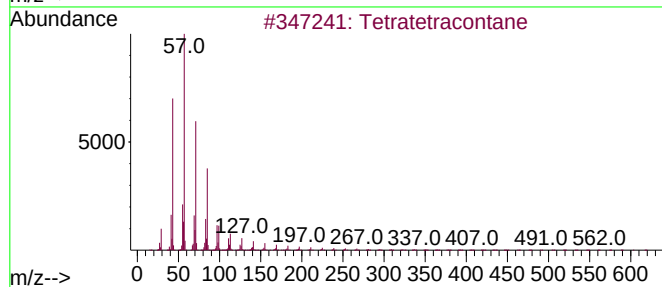
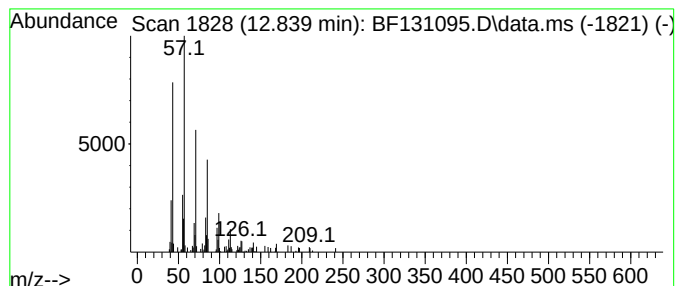
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetratetracontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.839	4.17 ng	109206	Chrysene-d12	14.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	91
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
Data File : BF131095.D
Acq On : 09 Nov 2022 19:32
Operator : CG\JU
Sample : N5437-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
596C-3

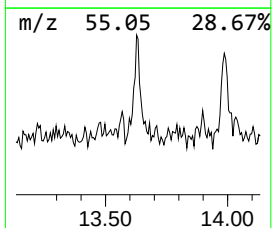
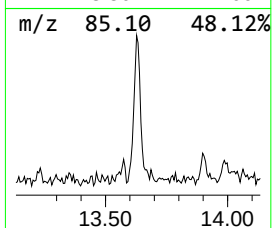
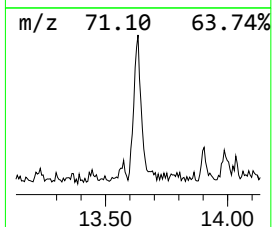
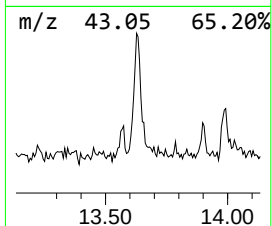
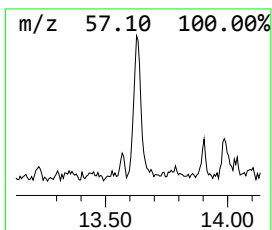
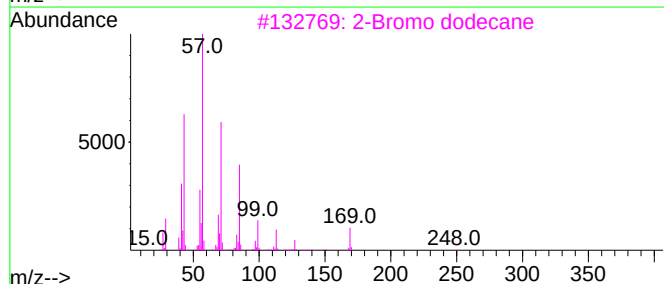
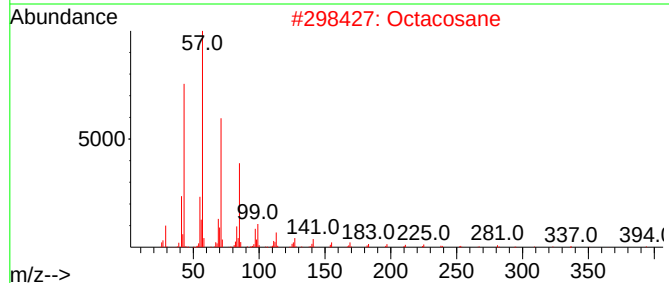
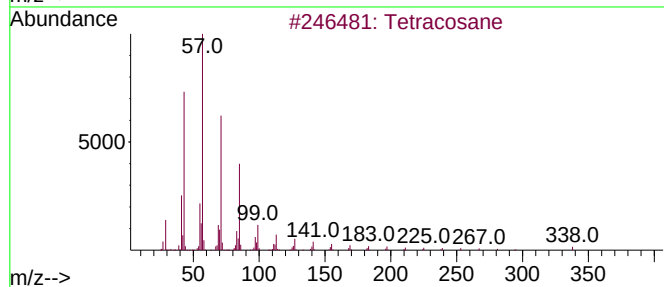
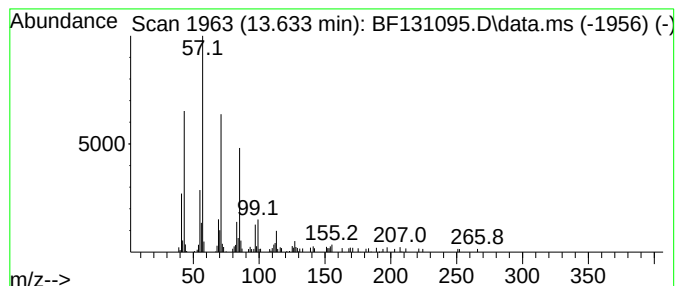
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.633	4.59 ng	120207	Chrysene-d12	14.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	87
2			Octacosane	394	C28H58	000630-02-4	87
3			2-Bromo dodecane	248	C12H25Br	013187-99-0	87
4			Pentacosane	352	C25H52	000629-99-2	86
5			Heptadecane	240	C17H36	000629-78-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
Data File : BF131095.D
Acq On : 09 Nov 2022 19:32
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Sample : N5437-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
596C-3

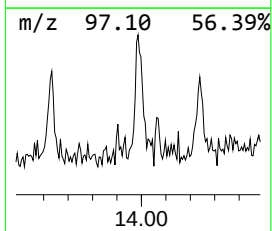
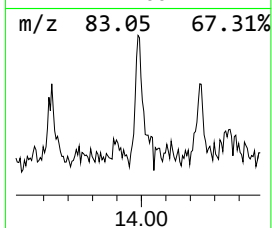
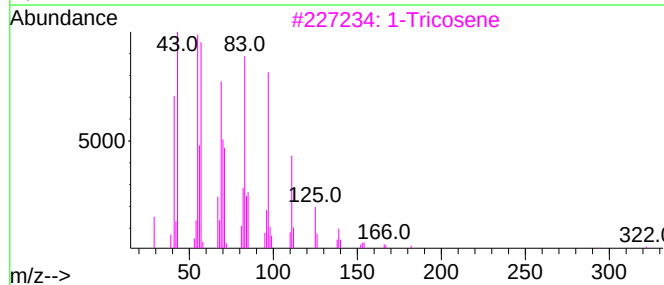
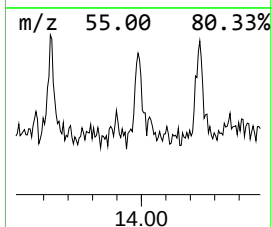
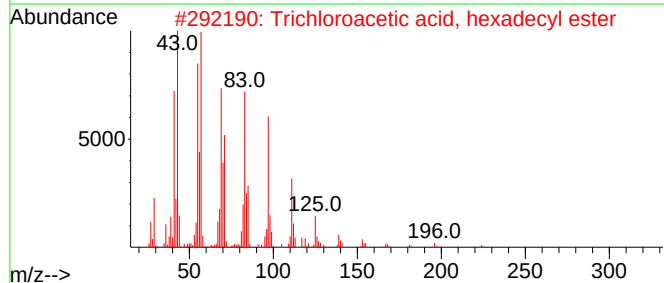
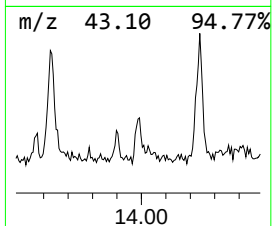
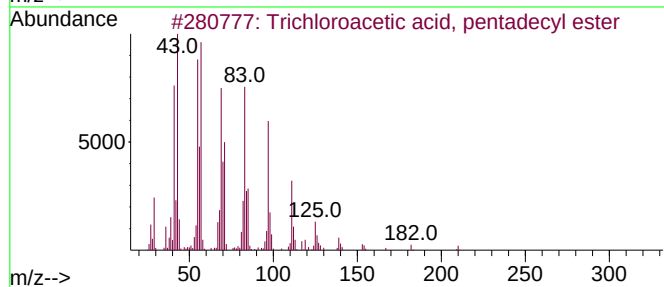
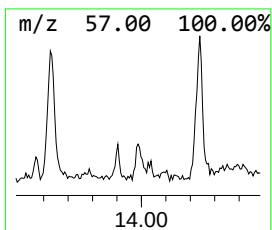
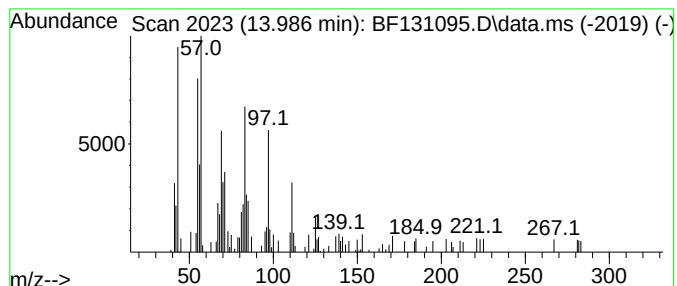
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Trichloroacetic acid, penta... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.986	2.46 ng	64594	Chrysene-d12	14.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	93
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
3			1-Tricosene	322	C23H46	018835-32-0	87
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	87
5			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
Data File : BF131095.D
Acq On : 09 Nov 2022 19:32
Operator : CG\JU
Sample : N5437-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
596C-3

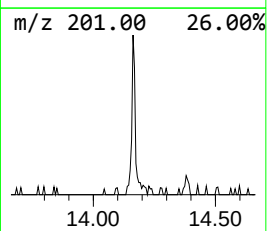
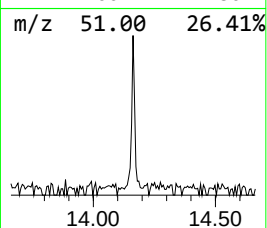
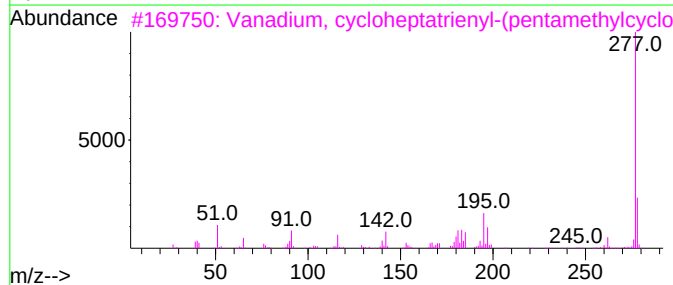
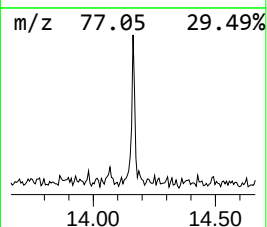
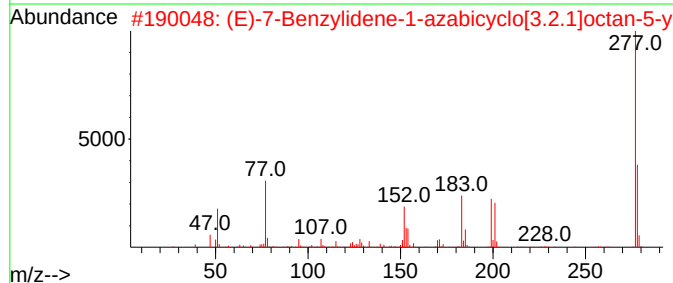
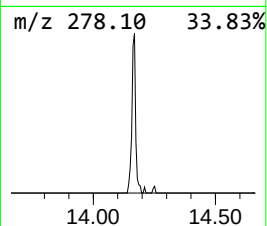
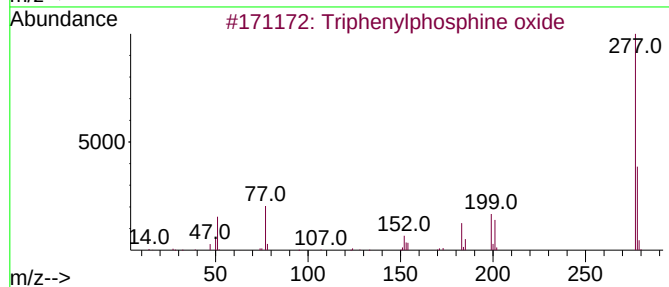
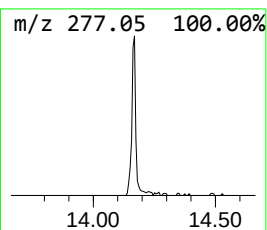
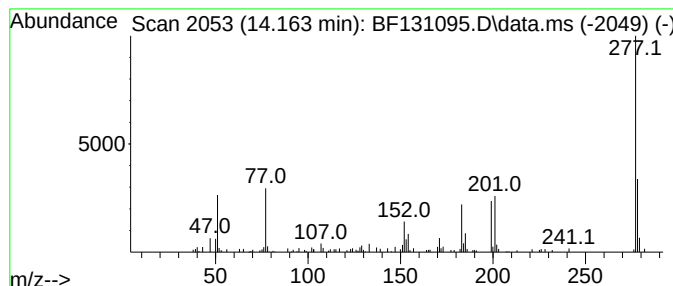
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Triphenylphosphine oxide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.163	2.58 ng	67663	Chrysene-d12	14.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	95
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	80
3			Vanadium, cycloheptatrienyl-(pen...	277	C17H22V	1000155-20-5	38
4			Cobalt, (.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	38
5			4-Nitro-4'-methylidiphenylsulfone	277	C13H11N04S	004094-37-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
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Operator : CG\JU
Sample : N5437-01
Misc :
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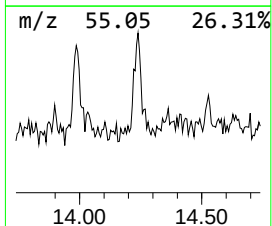
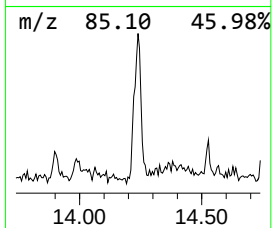
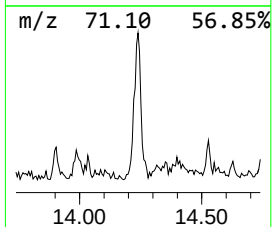
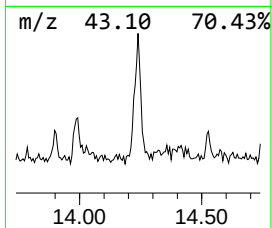
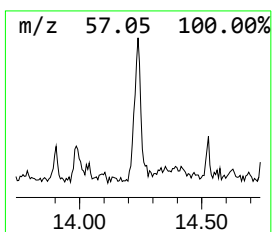
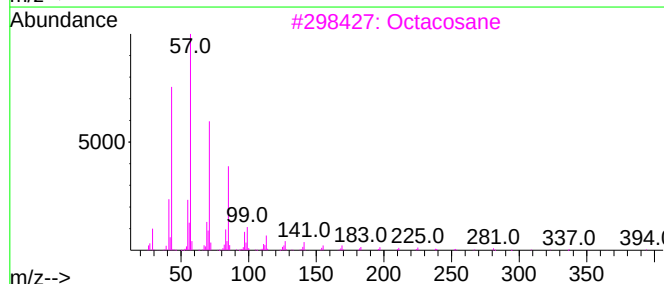
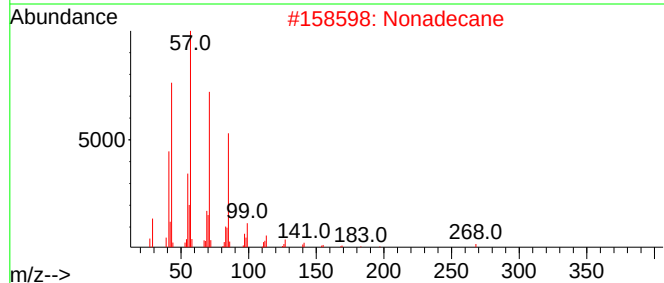
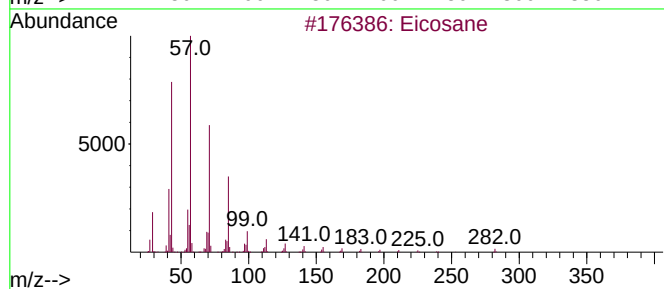
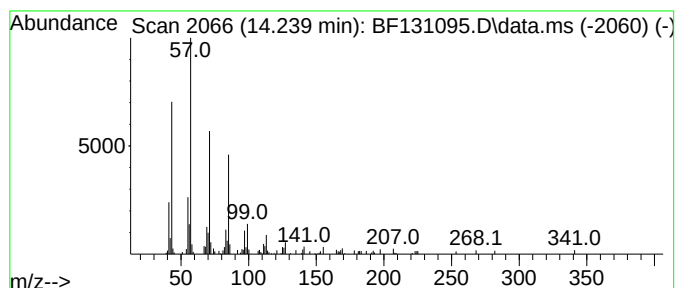
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Eicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.239	4.34 ng	113757	Chrysene-d12	14.069	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Nonadecane	268	C19H40	000629-92-5	95
3	Octacosane	394	C28H58	000630-02-4	90
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5	Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
Data File : BF131095.D
Acq On : 09 Nov 2022 19:32
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Sample : N5437-01
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Instrument :
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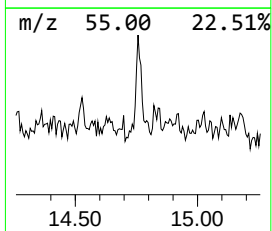
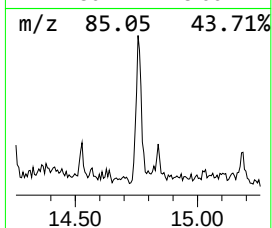
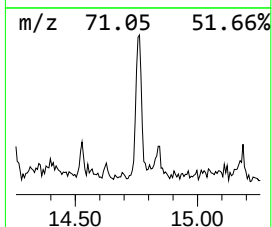
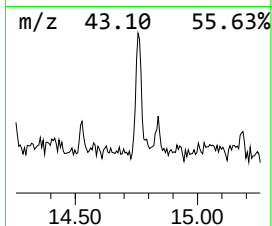
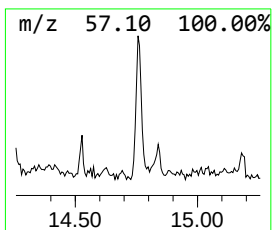
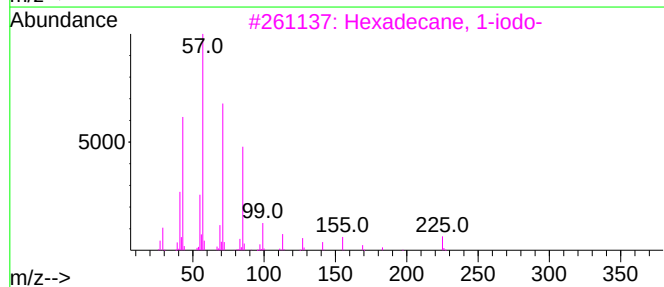
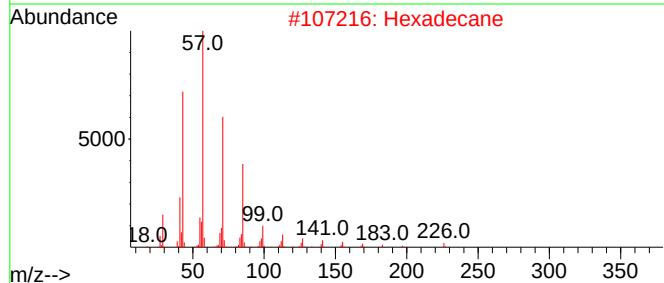
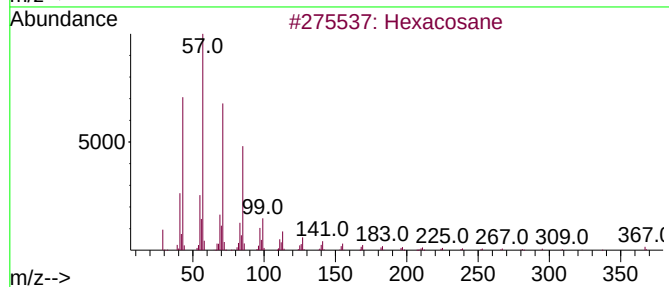
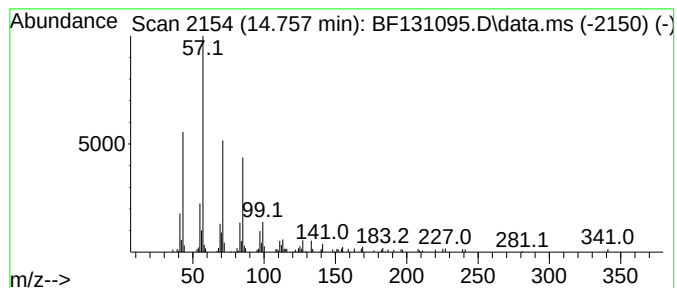
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.757	3.58 ng	93802	Chrysene-d12	14.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	87
2			Hexadecane	226	C16H34	000544-76-3	83
3			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	83
4			Heptadecane	240	C17H36	000629-78-7	83
5			Nonadecane	268	C19H40	000629-92-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
Data File : BF131095.D
Acq On : 09 Nov 2022 19:32
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Misc :
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Instrument :
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ClientSampleId :
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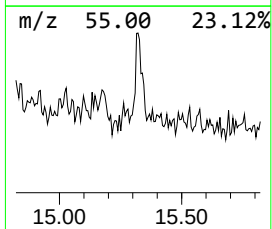
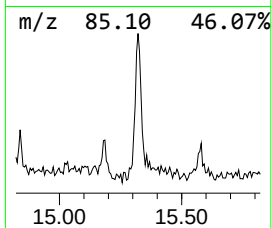
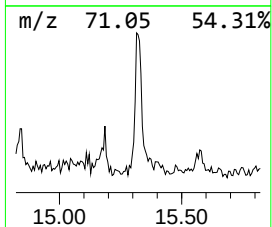
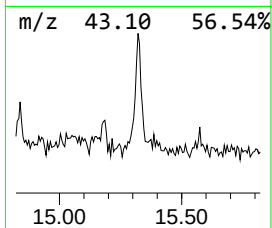
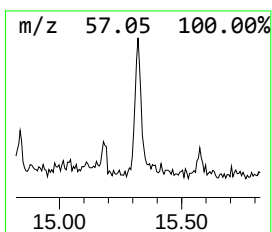
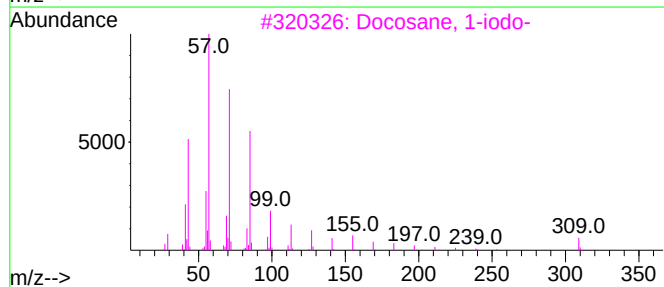
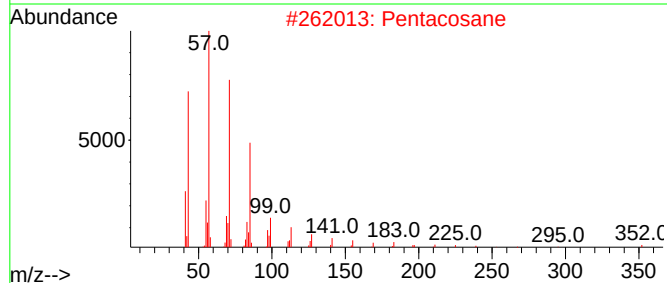
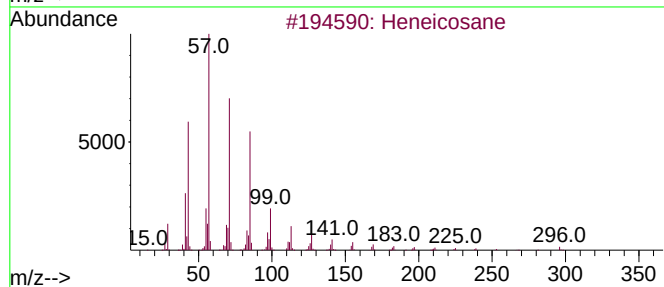
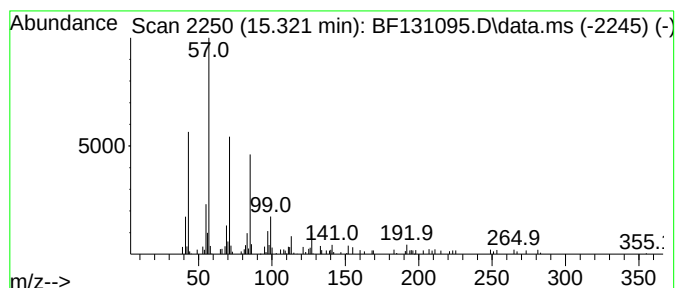
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.321	3.36 ng	79499	Perylene-d12	15.563

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	87
2		Pentacosane	352	C25H52	000629-99-2	87
3		Docosane, 1-iodo-	436	C22H45I	1000406-31-9	83
4		Hexacosane	366	C26H54	000630-01-3	80
5		Nonadecane	268	C19H40	000629-92-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
Data File : BF131095.D
Acq On : 09 Nov 2022 19:32
Operator : CG\JU
Sample : N5437-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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...	2.181	81.3	ng	2381070	1	6.893	586034	20.0
3-Penten-2-one,...	4.493	3.2	ng	93606	1	6.893	586034	20.0
2-Pentanone, 4-...	5.110	6.5	ng	188895	1	6.893	586034	20.0
n-Hexadecanoic ...	11.945	4.3	ng	183303	4	11.428	856692	20.0
1,3-Benzenedica...	12.345	2.2	ng	93687	4	11.428	856692	20.0
Tetratetracontane	12.839	4.2	ng	109206	5	14.069	524308	20.0
Tetracosane	13.633	4.6	ng	120207	5	14.069	524308	20.0
Trichloroacetic...	13.986	2.5	ng	64594	5	14.069	524308	20.0
Triphenylphosph...	14.163	2.6	ng	67663	5	14.069	524308	20.0
Eicosane	14.239	4.3	ng	113757	5	14.069	524308	20.0
Hexacosane	14.757	3.6	ng	93802	5	14.069	524308	20.0
Heneicosane	15.321	3.4	ng	79499	6	15.563	473140	20.0