

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
 Data File : BF134090.D
 Acq On : 06 Jul 2023 04:30
 Operator : CG\JU
 Sample : 03473-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CMC-EME-TOPSOIL-017

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-BF062623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF134090.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.128	3	7	18	rVB	744482	984156	57.58%	6.296%
2	2.246	22	27	37	rVB4	23291	46797	2.74%	0.299%
3	3.946	312	316	323	rVB	14072	20291	1.19%	0.130%
4	5.075	505	508	519	rVB	158470	180660	10.57%	1.156%
5	5.475	572	576	592	rBV	1603493	1466760	85.82%	9.383%
6	6.481	742	747	764	rBV	1532653	1570638	91.90%	10.048%
7	6.845	805	809	818	rBV	1001613	809371	47.35%	5.178%
8	7.404	900	904	915	rBV	1213541	1012986	59.27%	6.480%
9	8.122	1023	1026	1030	rBV	1169690	985647	57.67%	6.305%
10	9.198	1206	1209	1213	rBV	2038928	1709163	100.00%	10.934%
11	9.880	1322	1325	1329	rBV	1157590	971865	56.86%	6.217%
12	10.598	1443	1447	1453	rBV	211189	172823	10.11%	1.106%
13	10.675	1456	1460	1470	rBV	824862	751347	43.96%	4.807%
14	10.910	1496	1500	1503	rBV	32520	28078	1.64%	0.180%
15	11.374	1576	1579	1583	rBV	959814	860298	50.33%	5.504%
16	11.427	1586	1588	1591	rVB	45464	37725	2.21%	0.241%
17	12.968	1846	1850	1854	rBV	1332613	1299764	76.05%	8.315%
18	13.204	1887	1890	1894	rBV	50034	48575	2.84%	0.311%
19	13.551	1945	1949	1952	rBV	60148	61272	3.58%	0.392%
20	13.880	1991	2005	2017	rBV2	130827	381714	22.33%	2.442%
21	14.033	2027	2031	2038	rBV	742552	747418	43.73%	4.781%
22	14.198	2056	2059	2063	rBV	101380	98360	5.75%	0.629%
23	14.510	2108	2112	2115	rBV	118200	120367	7.04%	0.770%
24	14.821	2162	2165	2171	rVB	113215	127944	7.49%	0.818%
25	14.898	2174	2178	2183	rVB	126839	129843	7.60%	0.831%
26	15.174	2221	2225	2230	rBV	129884	156414	9.15%	1.001%
27	15.539	2282	2287	2291	rBV2	425962	586427	34.31%	3.752%
28	16.033	2367	2371	2375	rBV2	56886	86213	5.04%	0.552%
29	16.480	2444	2447	2454	rVB9	42615	71562	4.19%	0.458%
30	16.574	2458	2463	2469	rBV2	48864	107142	6.27%	0.685%

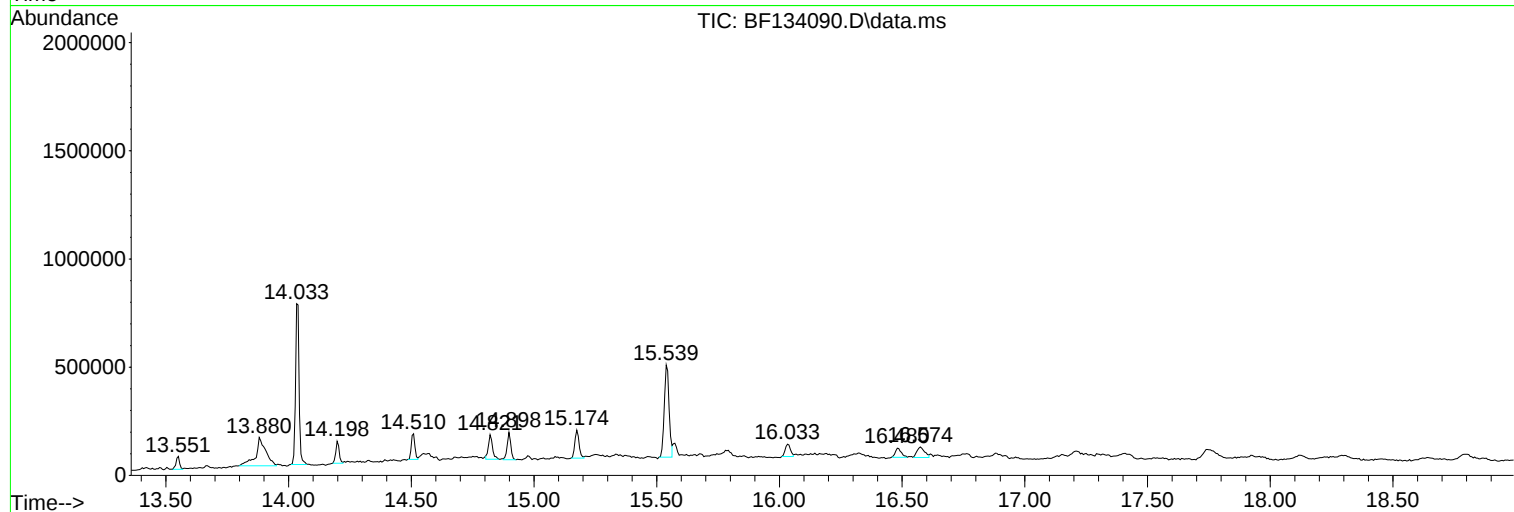
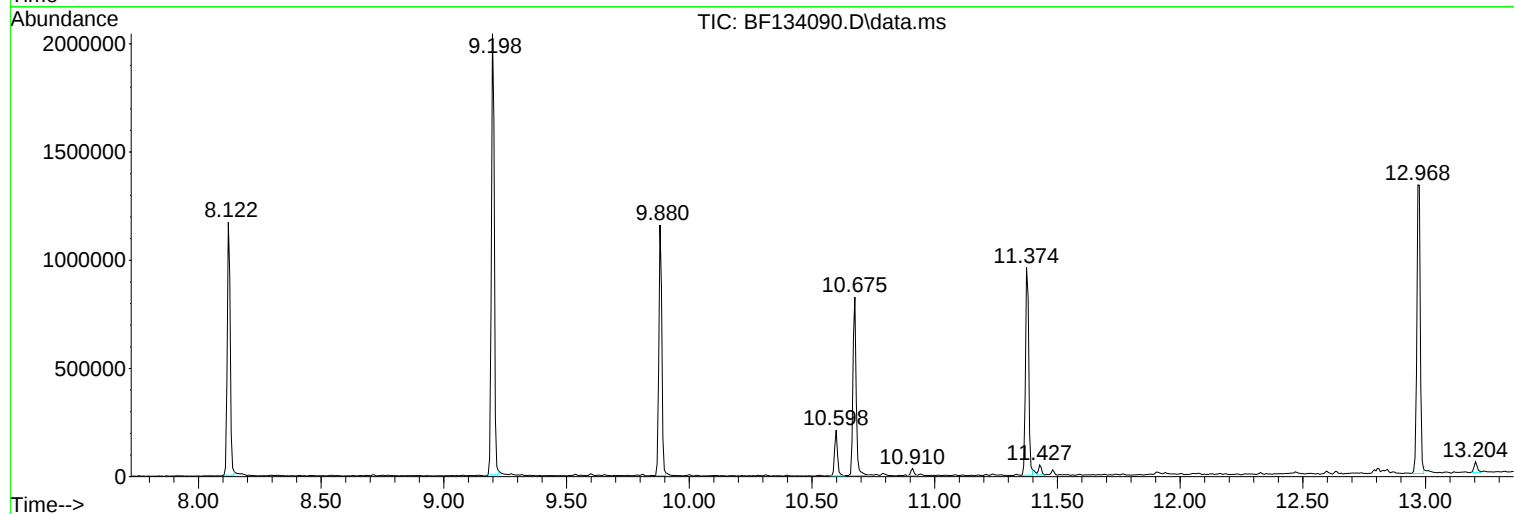
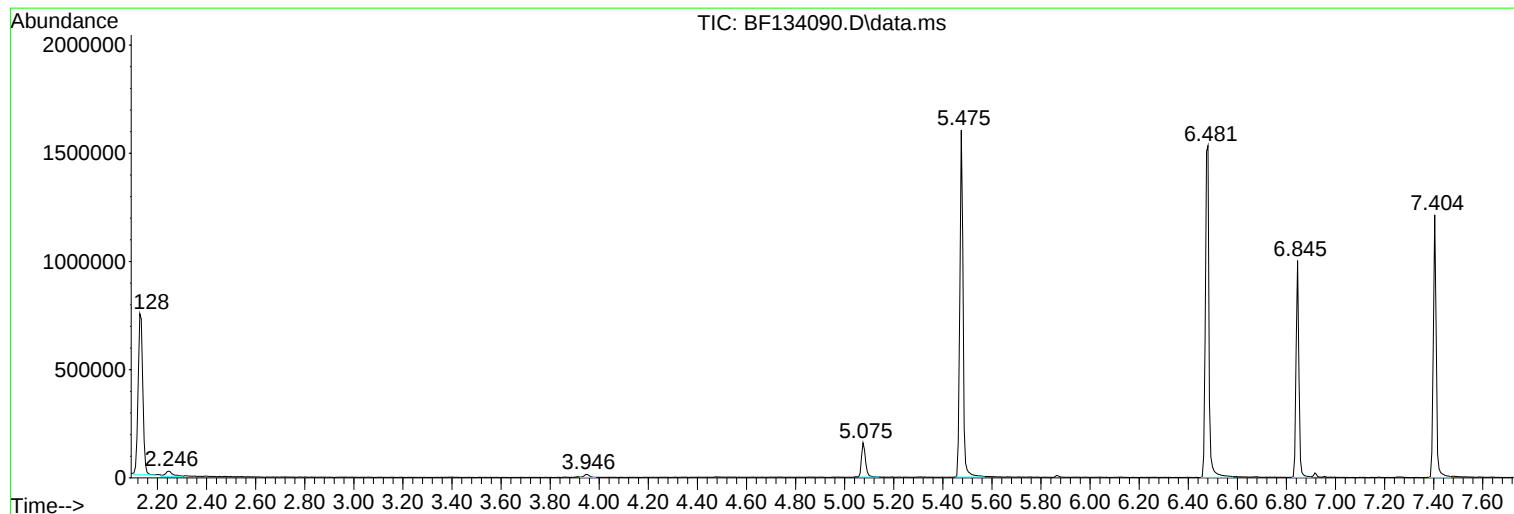
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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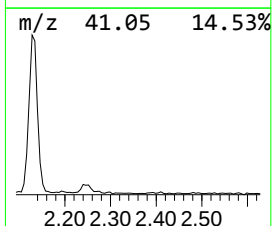
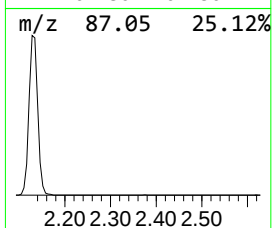
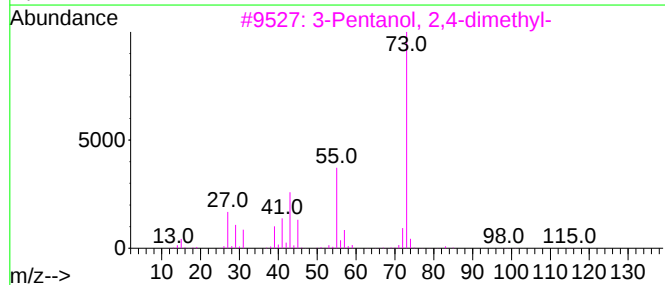
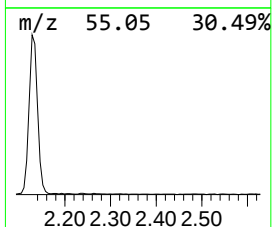
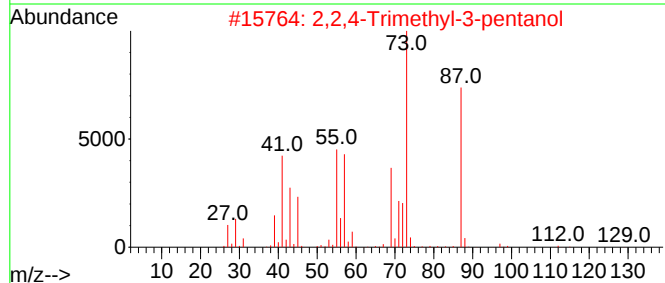
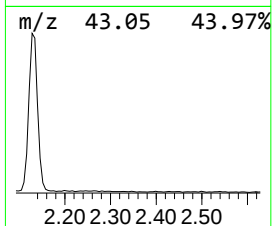
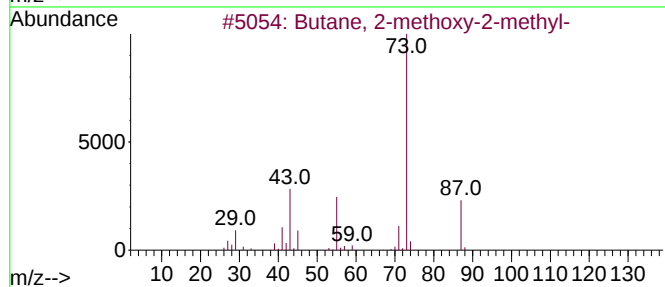
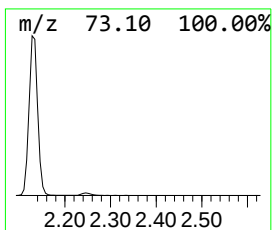
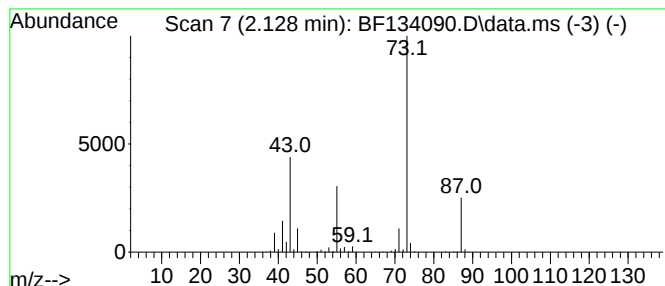
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.128	24.32 ng	984156	1,4-Dichlorobenzene-d4	6.845

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			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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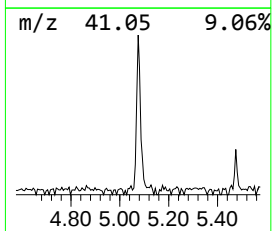
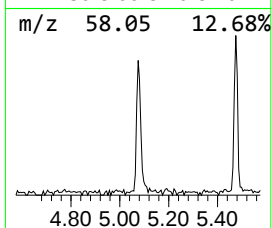
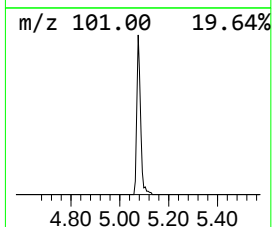
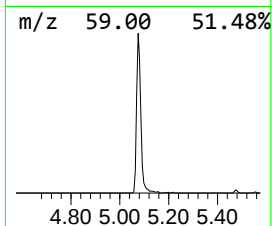
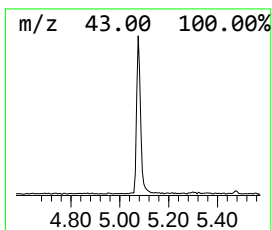
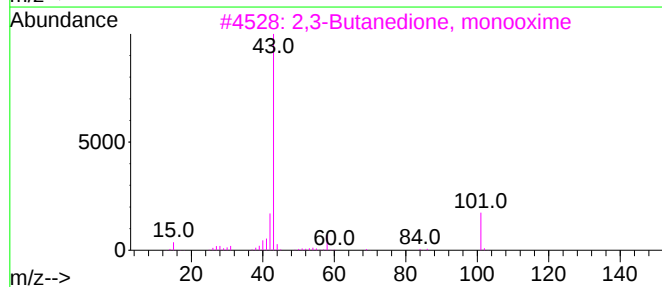
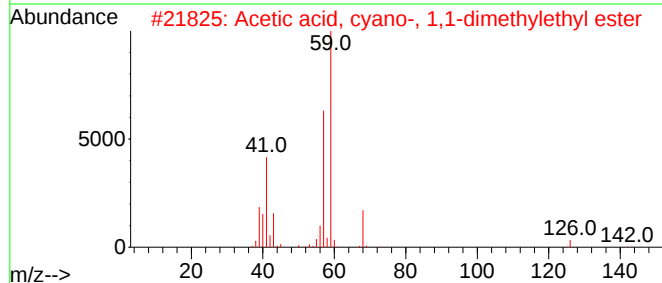
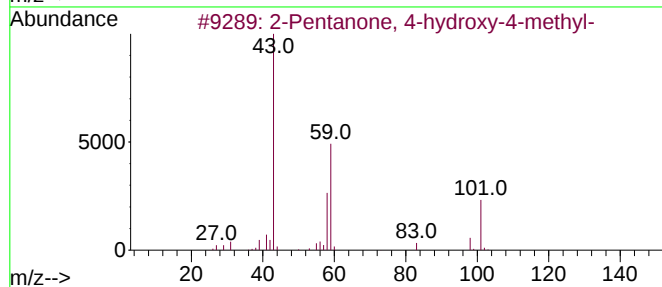
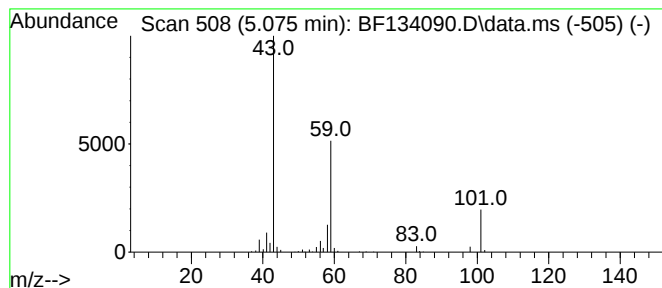
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.075	4.46 ng	180660	1,4-Dichlorobenzene-d4	6.845

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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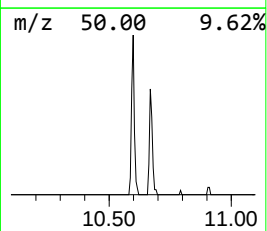
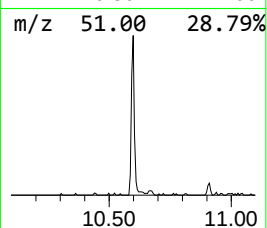
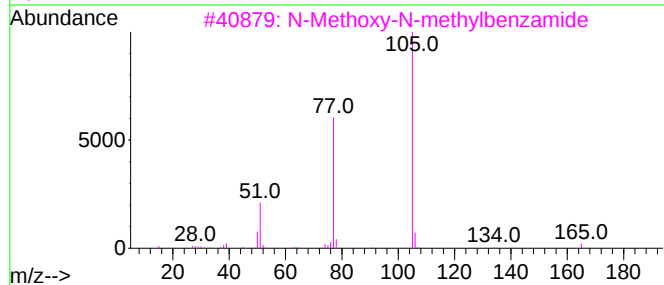
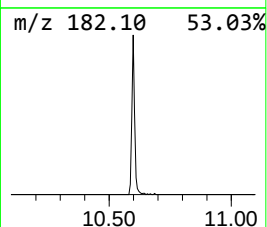
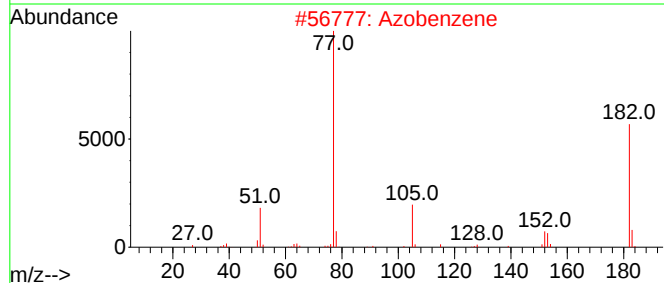
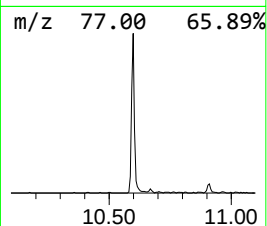
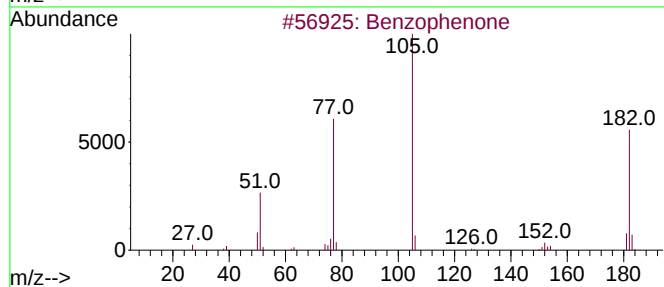
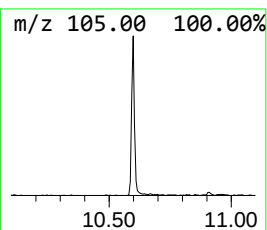
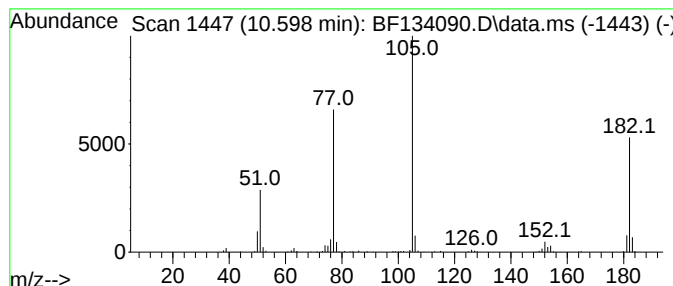
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.598	3.56 ng	172823	Acenaphthene-d10	9.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Azobenzene	182	C12H10N2	000103-33-3	58
3			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	50
4			acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	50
5			2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9NO2	037791-41-6	50



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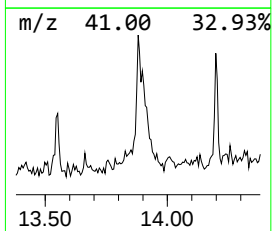
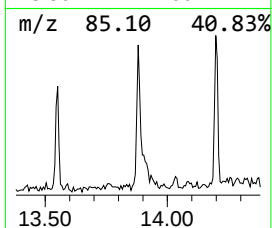
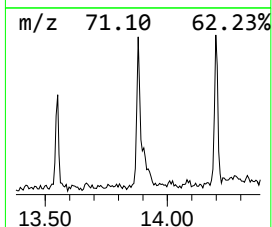
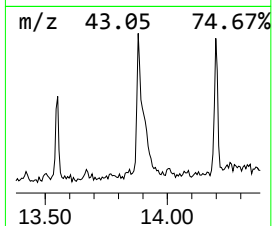
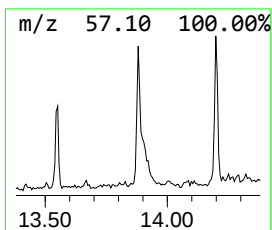
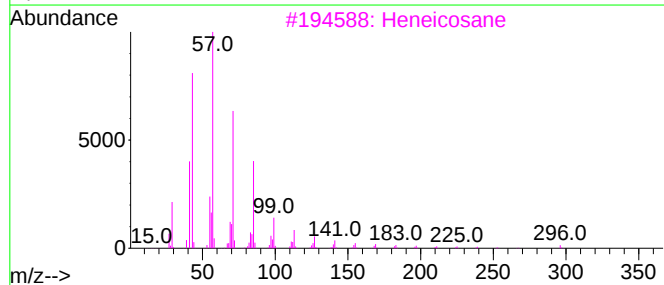
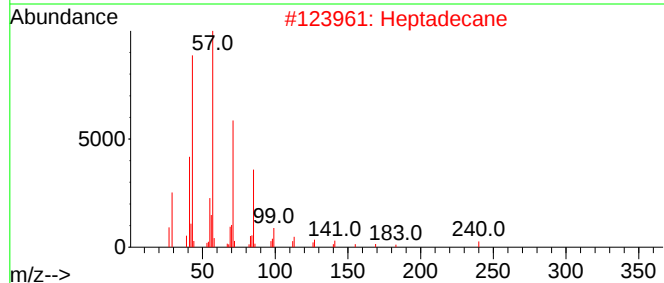
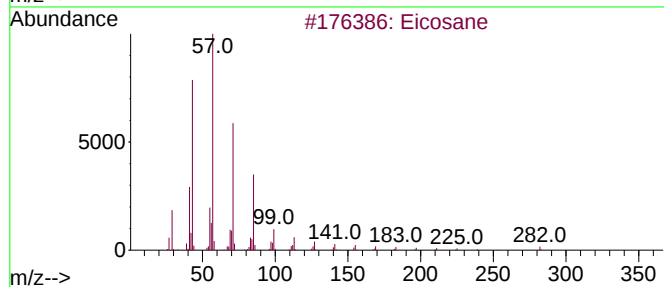
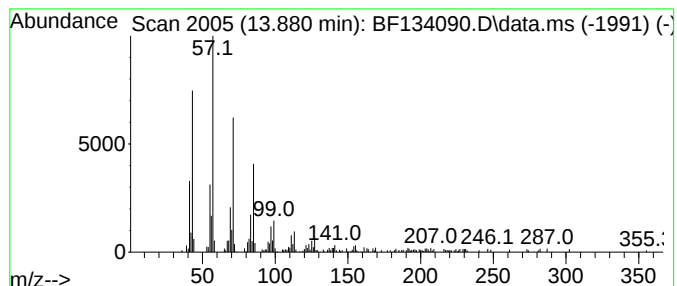
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Peak Number 4 Eicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.880	10.21 ng	381714	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Heptadecane	240	C17H36	000629-78-7	94
3			Heneicosane	296	C21H44	000629-94-7	90
4			Tetratetracontane	619	C44H90	007098-22-8	90
5			Hexatriacontane	507	C36H74	000630-06-8	87



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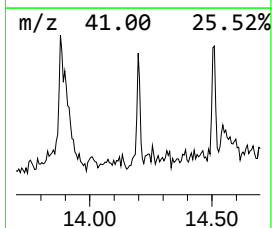
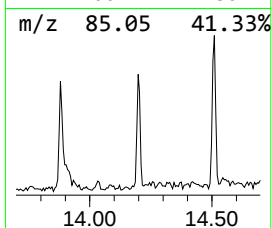
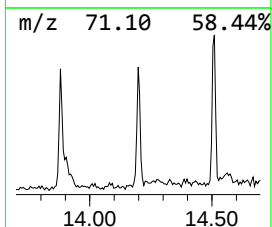
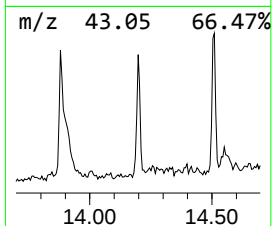
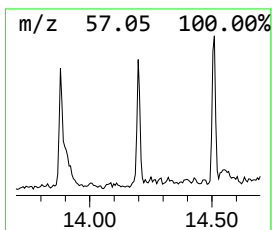
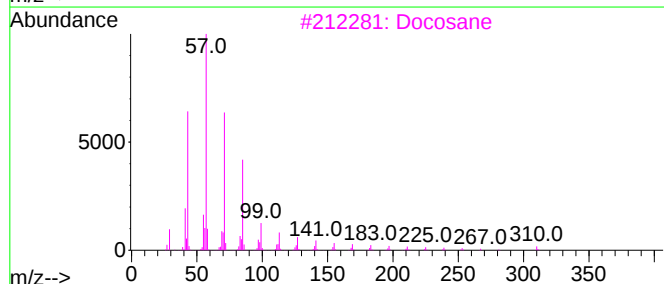
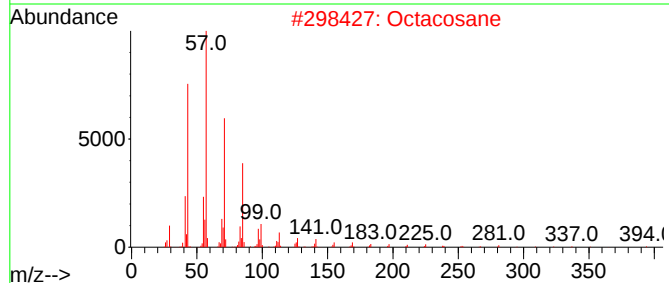
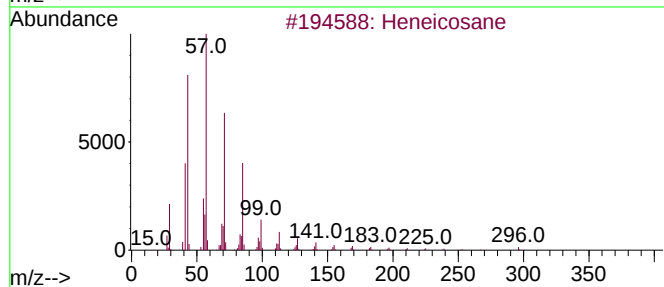
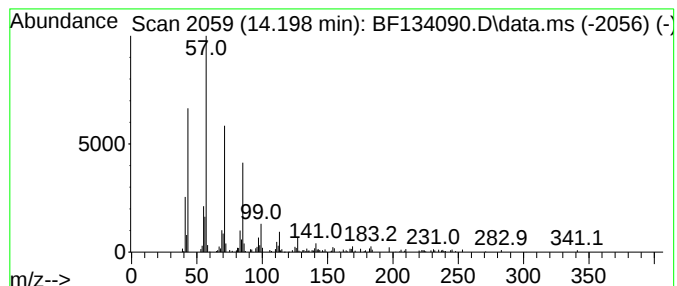
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Peak Number 5 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.198	2.63 ng	98360	Chrysene-d12	14.039	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Octacosane	394	C28H58	000630-02-4	90
3	Docosane	310	C22H46	000629-97-0	87
4	Eicosane	282	C20H42	000112-95-8	87
5	Heptacosane	380	C27H56	000593-49-7	87



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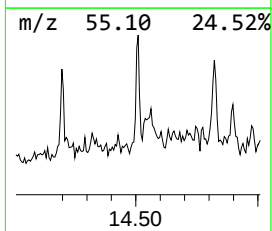
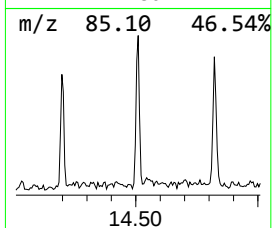
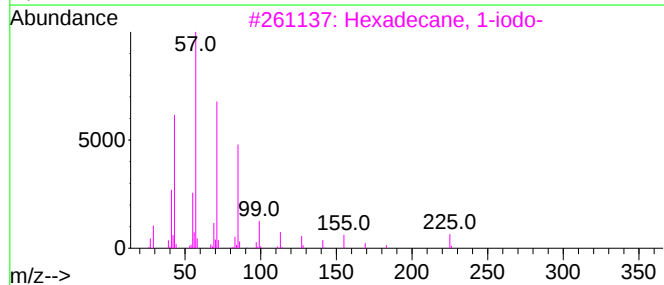
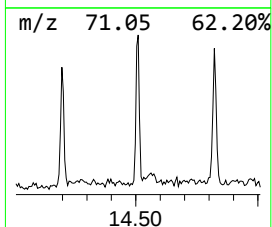
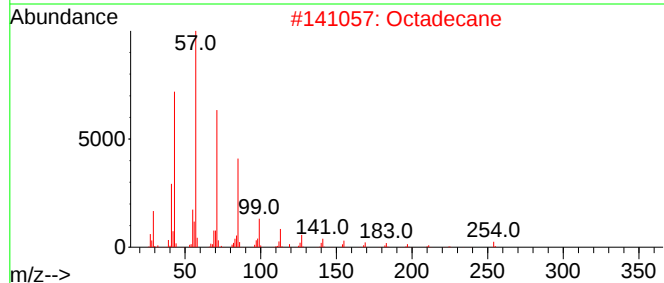
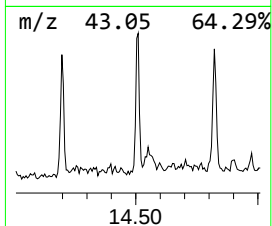
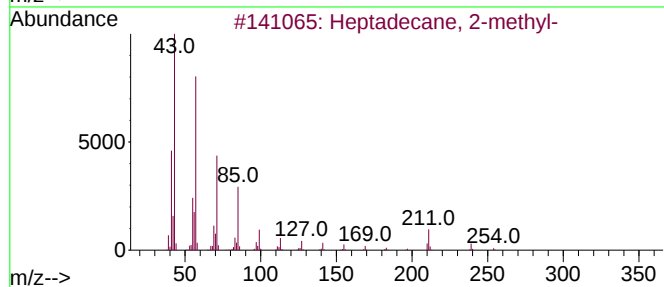
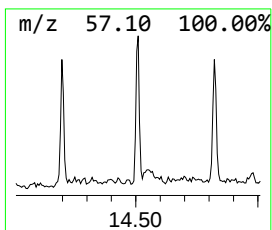
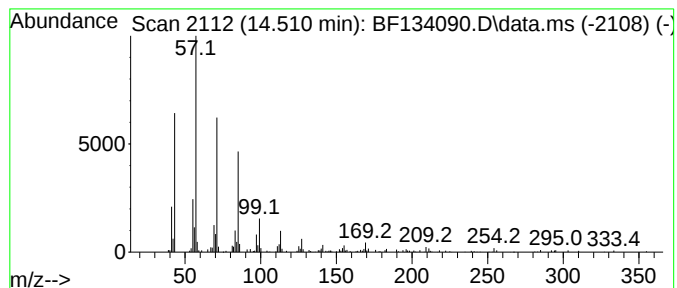
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.510	3.22 ng	120367	Chrysene-d12	14.039

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	94
2		Octadecane	254	C18H38	000593-45-3	91
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	91
4		Docosane	310	C22H46	000629-97-0	91
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134090.D
Acq On : 06 Jul 2023 04:30
Operator : CG\JU
Sample : 03473-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CMC-EME-TOPSOIL-017

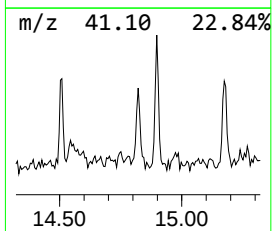
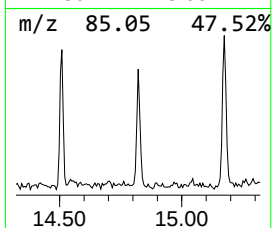
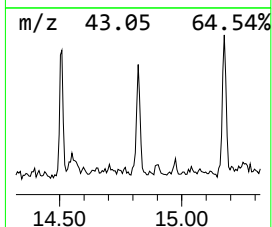
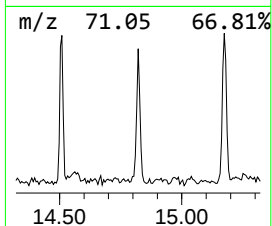
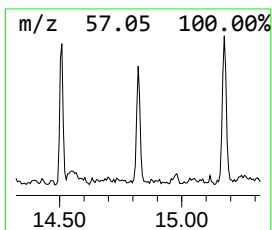
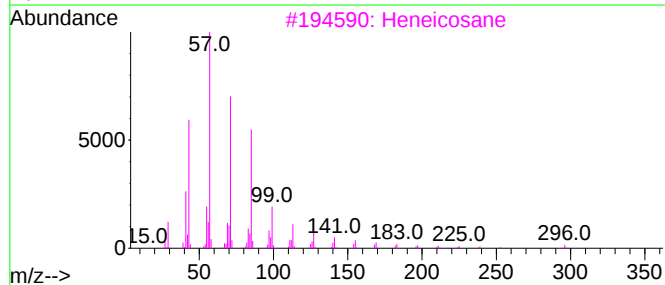
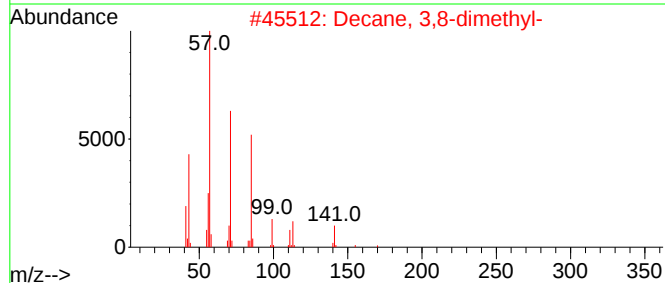
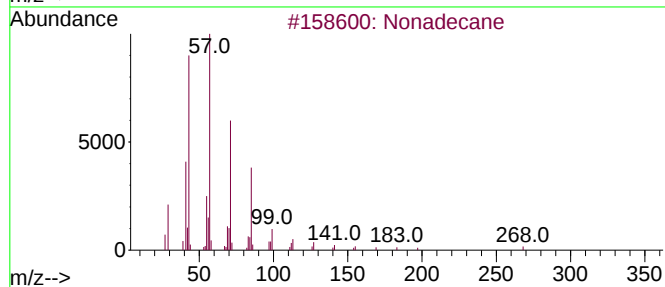
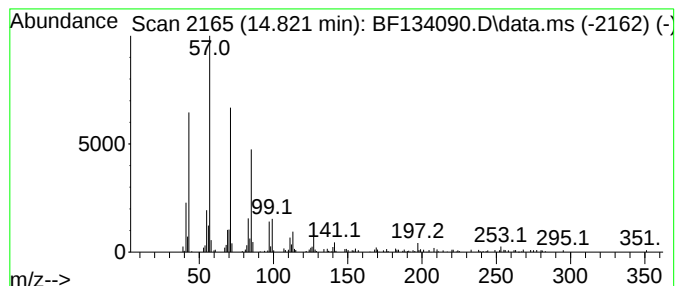
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.821	4.36 ng	127944	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			Hentriacontane	437	C31H64	000630-04-6	87
5			Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134090.D
Acq On : 06 Jul 2023 04:30
Operator : CG\JU
Sample : 03473-12
Misc :
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Instrument :
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ClientSampleId :
CMC-EME-TOPSOIL-017

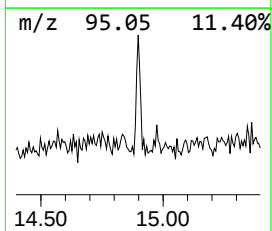
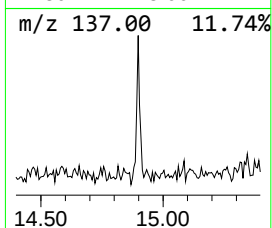
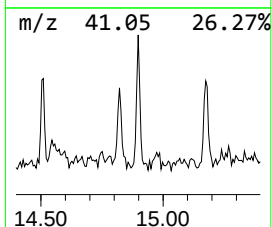
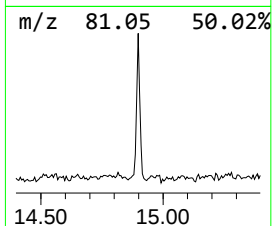
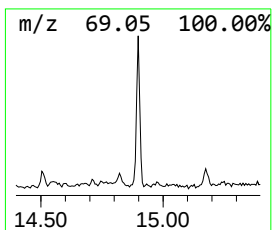
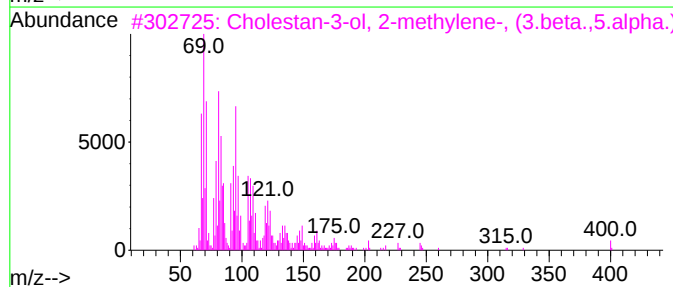
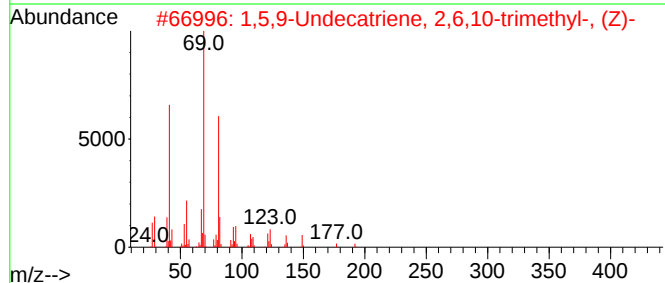
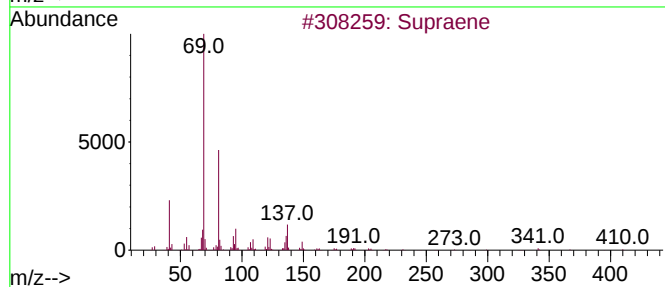
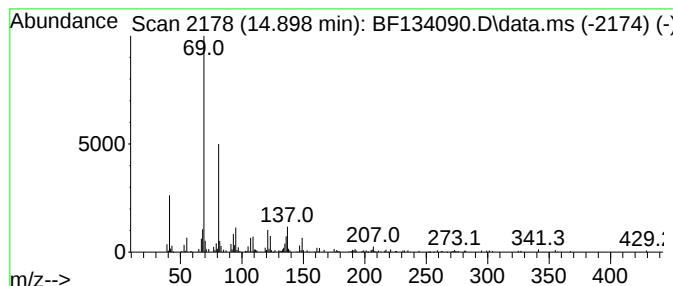
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Supraene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.898	4.43 ng	129843	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	90
2			1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	83
3			Cholestan-3-ol, 2-methylene-, (3...	400	C28H48O	022599-96-8	76
4			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5			1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134090.D
Acq On : 06 Jul 2023 04:30
Operator : CG\JU
Sample : 03473-12
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CMC-EME-TOPSOIL-017

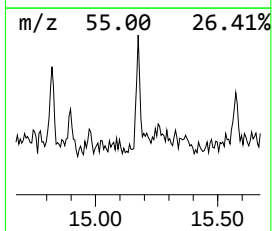
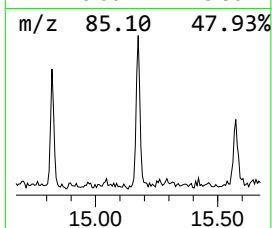
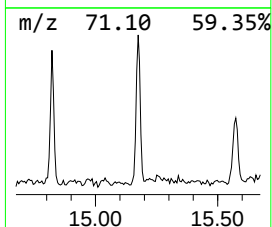
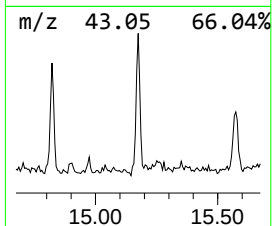
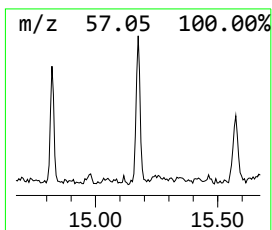
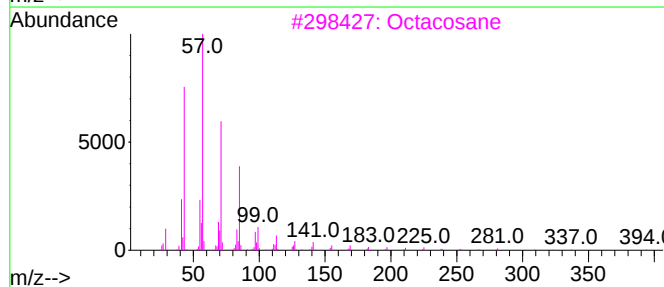
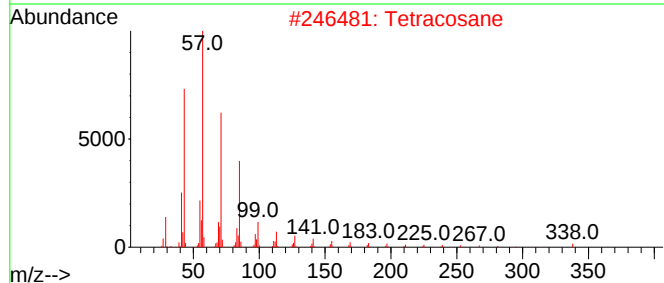
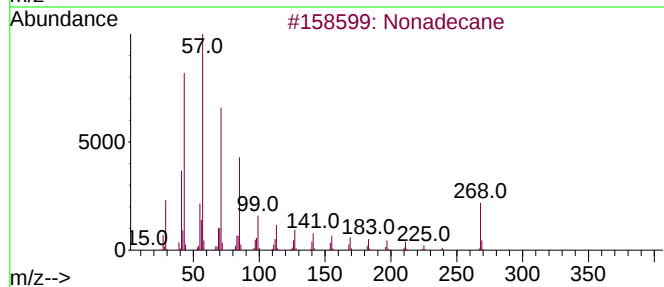
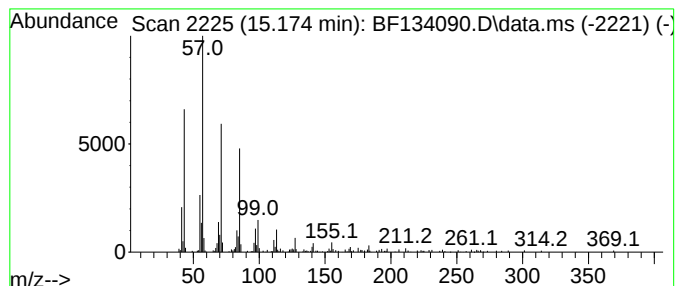
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.174	5.33 ng	156414	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Tetracosane	338	C24H50	000646-31-1	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Docosane	310	C22H46	000629-97-0	87
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134090.D
Acq On : 06 Jul 2023 04:30
Operator : CG\JU
Sample : 03473-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CMC-EME-TOPSOIL-017

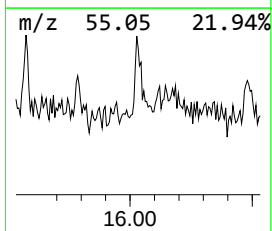
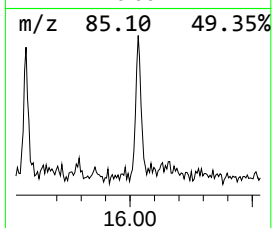
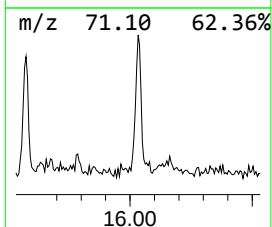
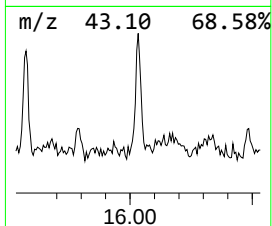
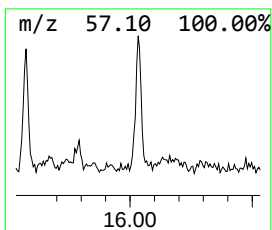
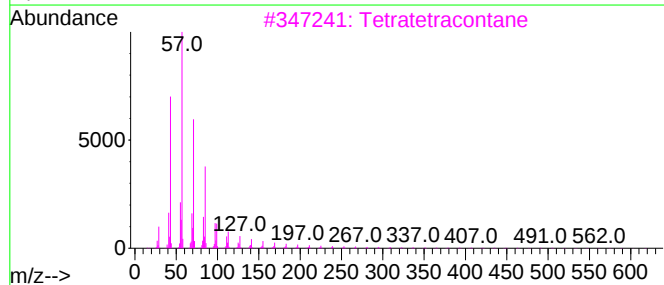
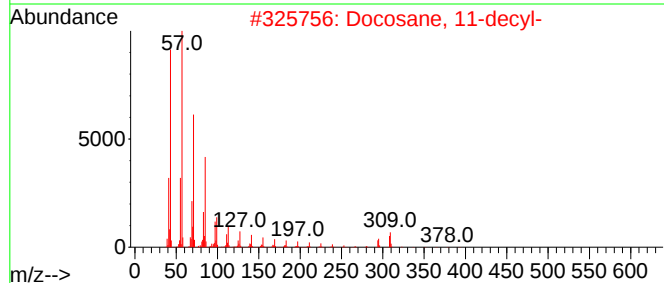
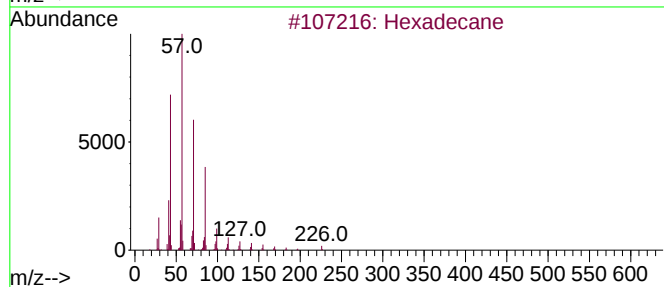
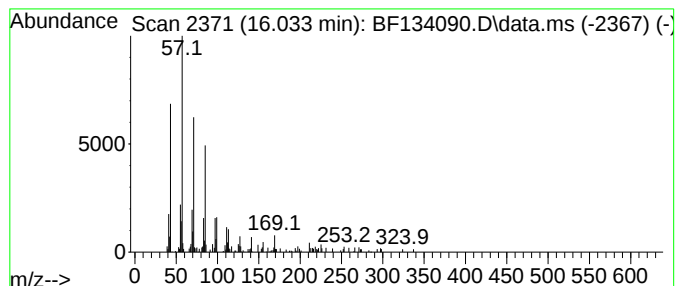
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.033	2.94 ng	86213	Perylene-d12	15.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	95
2		Docosane, 11-decyl-	451	C32H66	055401-55-3	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5		2-Methylhexacosane	380	C27H56	001561-02-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134090.D
Acq On : 06 Jul 2023 04:30
Operator : CG\JU
Sample : 03473-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CMC-EME-TOPSOIL-017

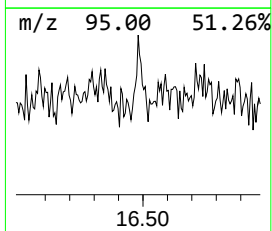
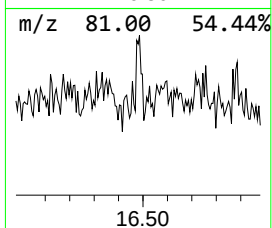
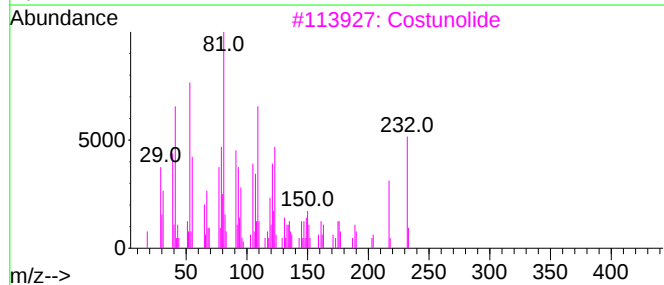
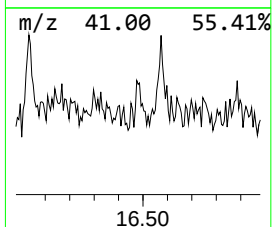
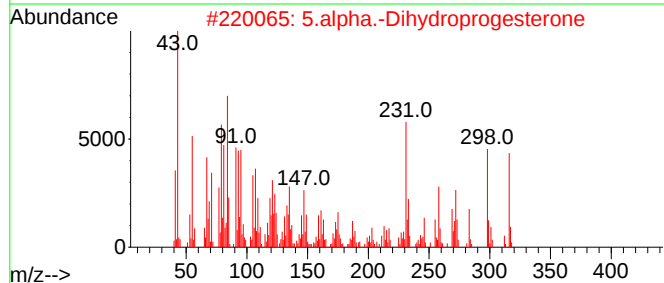
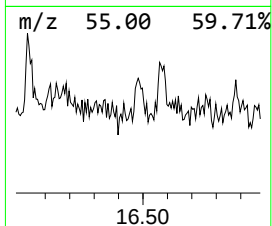
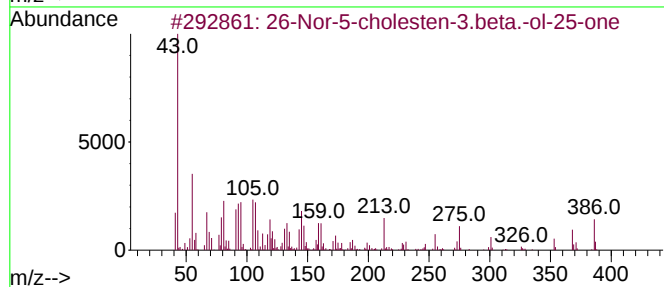
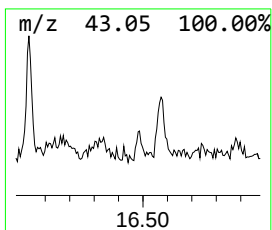
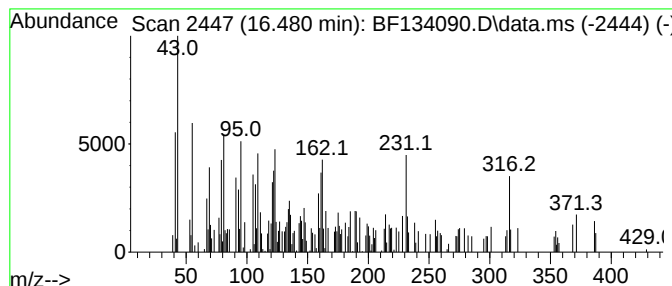
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 26-Nor-5-cholesten-3.beta.-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.480	2.44 ng	71562	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	53		
2	5.alpha.-Dihydroprogesterone	316	C21H32O2	000566-65-4	38		
3	Costunolide	232	C15H20O2	000553-21-9	30		
4	3-n-Heptyl-7-methyl-9-(2,6,6-tri...	368	C26H40O	1000216-09-3	30		
5	5.alpha.-Pregnane-12,20-dione	316	C21H32O2	006022-48-6	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
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Operator : CG\JU
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ALS Vial : 19 Sample Multiplier: 1

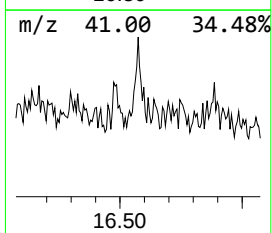
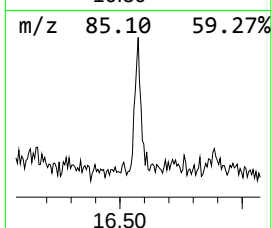
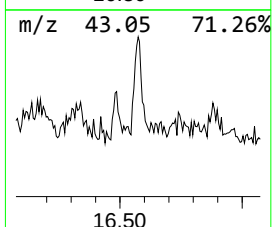
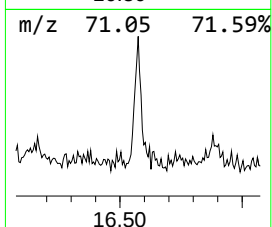
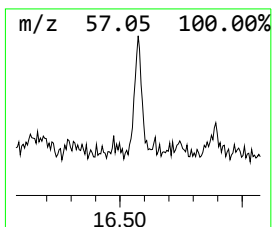
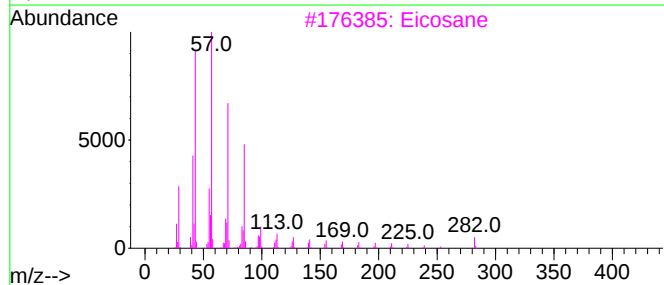
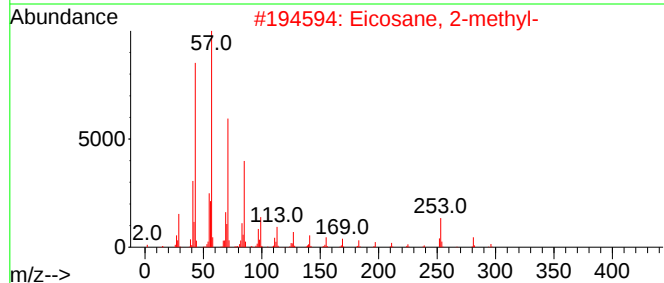
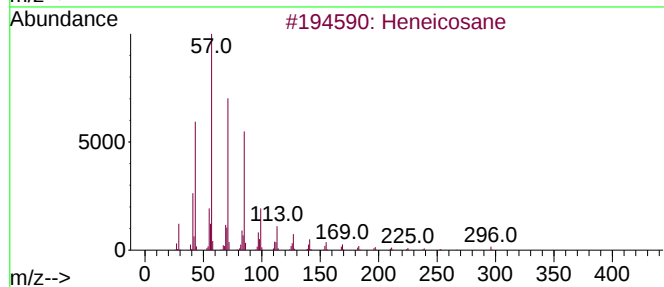
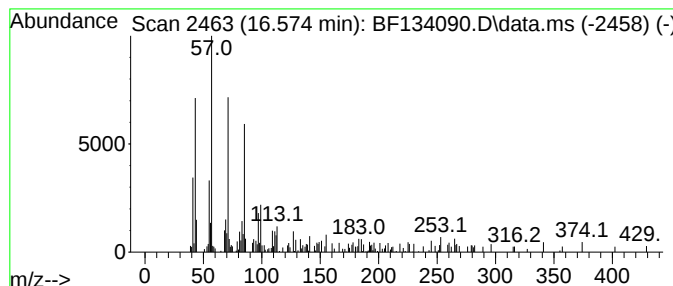
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CMC-EME-TOPSOIL-017

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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, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.574	3.65 ng	107142	Perylene-d12	15.539	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Eicosane, 2-methyl-	296	C21H44	001560-84-5	90
3	Eicosane	282	C20H42	000112-95-8	86
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	81
5	Pentacosane	352	C25H52	000629-99-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134090.D
Acq On : 06 Jul 2023 04:30
Operator : CG\JU
Sample : 03473-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CMC-EME-TOPSOIL-017

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.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...	2.128	24.3	ng	984156	1	6.845	809371	20.0
2-Pentanone, 4-...	5.075	4.5	ng	180660	1	6.845	809371	20.0
Benzophenone	10.598	3.6	ng	172823	3	9.880	971865	20.0
Eicosane	13.880	10.2	ng	381714	5	14.039	747418	20.0
Heneicosane	14.198	2.6	ng	98360	5	14.039	747418	20.0
Heptadecane, 2-...	14.510	3.2	ng	120367	5	14.039	747418	20.0
Nonadecane	14.821	4.4	ng	127944	6	15.539	586427	20.0
Supraene	14.898	4.4	ng	129843	6	15.539	586427	20.0
Tetracosane	15.174	5.3	ng	156414	6	15.539	586427	20.0
Hexadecane	16.033	2.9	ng	86213	6	15.539	586427	20.0
26-Nor-5-choles...	16.480	2.4	ng	71562	6	15.539	586427	20.0
Eicosane, 2-met...	16.574	3.6	ng	107142	6	15.539	586427	20.0