

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
 Data File : BF134082.D
 Acq On : 05 Jul 2023 23:56
 Operator : CG\JU
 Sample : 03473-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CMC-EME-TOPSOIL-008

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: BF134082.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.152	4	11	20	rVB	1117326	1468524	64.95%	7.642%
2	2.263	26	30	41	rVB	33695	56727	2.51%	0.295%
3	5.087	506	510	521	rVB	192151	235506	10.42%	1.226%
4	5.481	573	577	588	rBV	2149340	2111574	93.39%	10.989%
5	6.481	743	747	764	rBV	2122508	2167681	95.87%	11.281%
6	6.845	805	809	819	rBV	1023202	823529	36.42%	4.286%
7	7.410	900	905	908	rBV	1653313	1510937	66.82%	7.863%
8	8.122	1023	1026	1030	rBV	1276684	1117155	49.41%	5.814%
9	9.204	1206	1210	1213	rBV	2702863	2261071	100.00%	11.767%
10	9.881	1322	1325	1329	rBV	1251633	1125824	49.79%	5.859%
11	10.598	1443	1447	1456	rBV	196971	160001	7.08%	0.833%
12	10.675	1456	1460	1468	rBV	1208097	1029913	45.55%	5.360%
13	11.375	1576	1579	1583	rBV	925501	851585	37.66%	4.432%
14	11.428	1586	1588	1592	rVB2	29613	29240	1.29%	0.152%
15	11.904	1661	1669	1673	rBV3	23863	36781	1.63%	0.191%
16	12.698	1799	1804	1811	rBV7	21683	44903	1.99%	0.234%
17	12.810	1819	1823	1825	rBV2	34411	34421	1.52%	0.179%
18	12.974	1847	1851	1854	rBV	1725752	1531282	67.72%	7.969%
19	13.204	1887	1890	1897	rBV	55462	49895	2.21%	0.260%
20	13.551	1944	1949	1952	rBV	73935	75025	3.32%	0.390%
21	13.880	2002	2005	2017	rVB2	133584	291898	12.91%	1.519%
22	14.039	2028	2032	2042	rVB	742608	745736	32.98%	3.881%
23	14.198	2056	2059	2062	rBV	106388	101474	4.49%	0.528%
24	14.510	2108	2112	2115	rBV	141826	125842	5.57%	0.655%
25	14.821	2162	2165	2170	rVB	118165	121796	5.39%	0.634%
26	14.898	2175	2178	2183	rVB	40620	43484	1.92%	0.226%
27	15.174	2221	2225	2229	rBV	150682	170335	7.53%	0.886%
28	15.539	2282	2287	2291	rBV	403837	567194	25.09%	2.952%
29	16.033	2367	2371	2379	rBV2	70840	113521	5.02%	0.591%
30	16.574	2460	2463	2471	rVB2	44062	74118	3.28%	0.386%
31	17.751	2658	2663	2677	rVB2	40522	139046	6.15%	0.724%

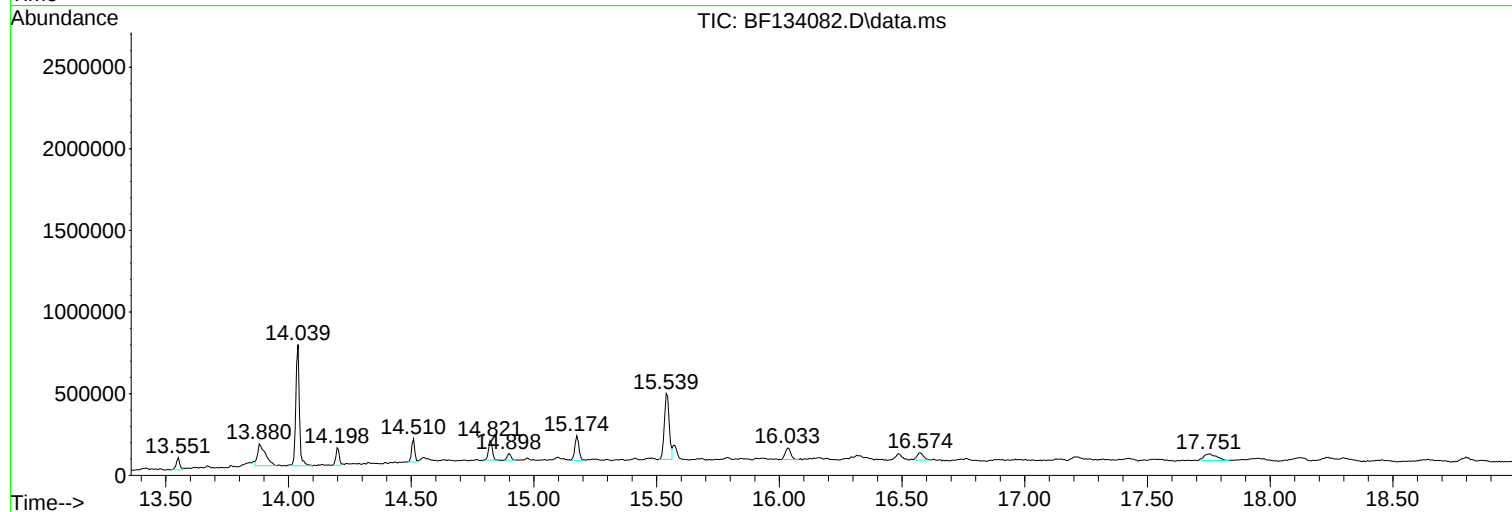
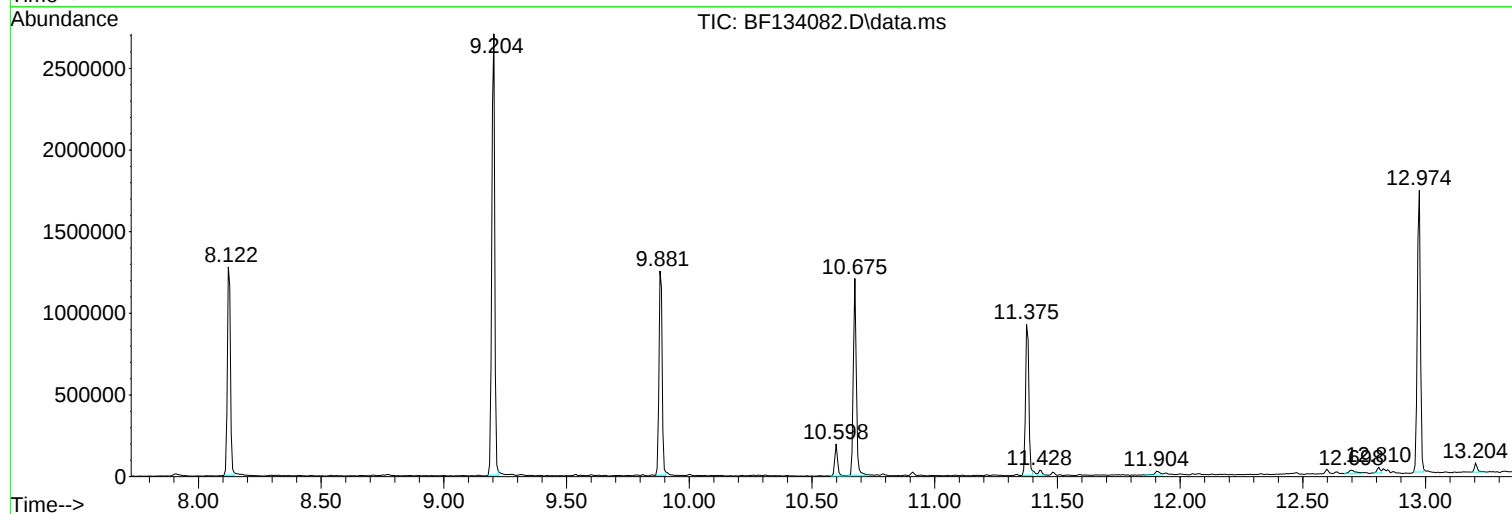
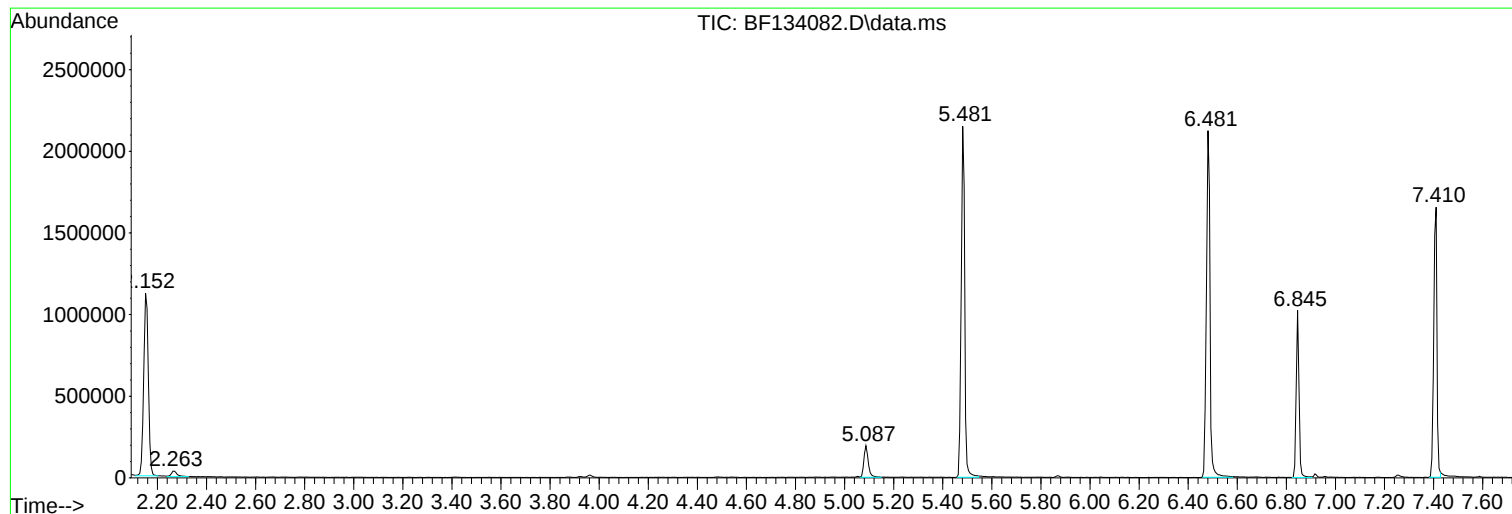
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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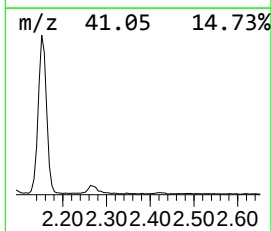
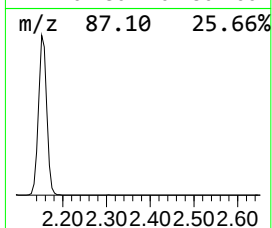
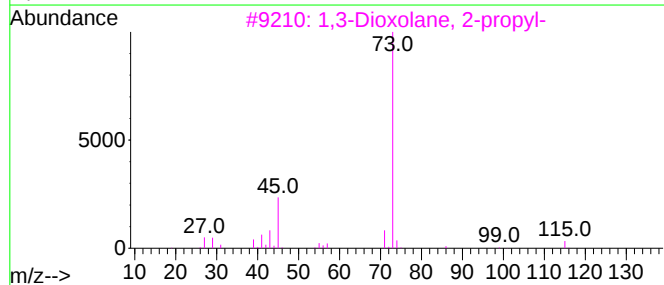
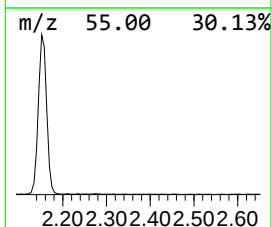
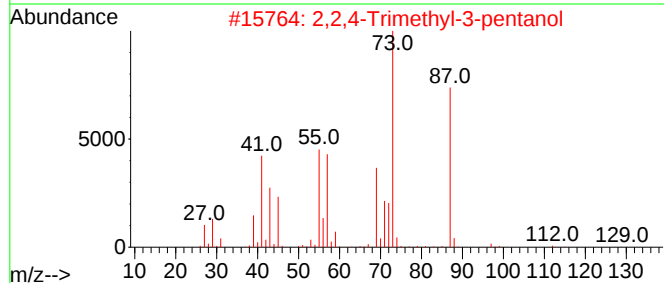
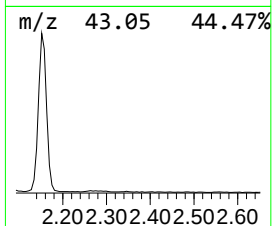
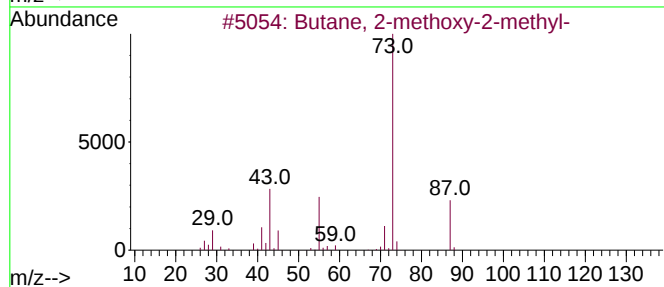
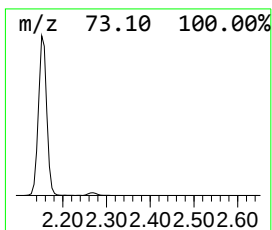
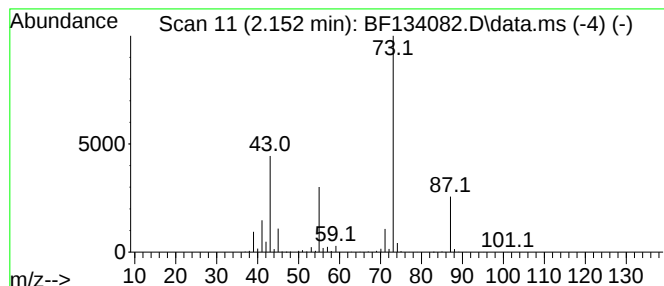
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.152	35.66 ng	1468520	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			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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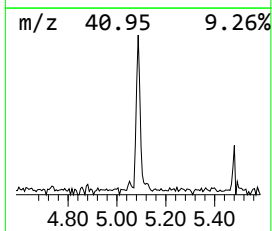
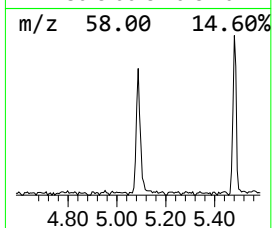
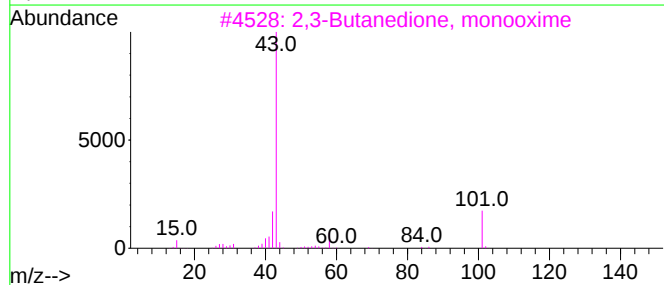
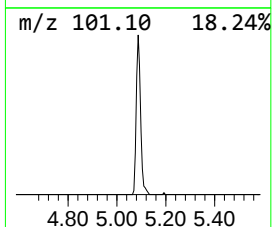
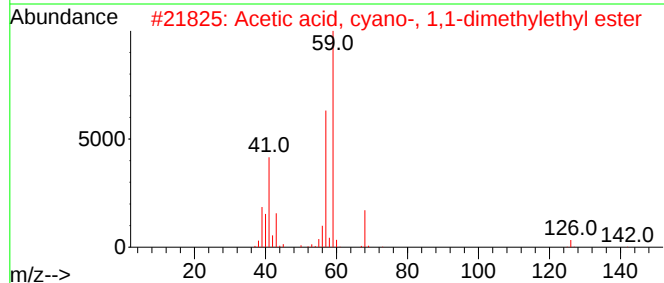
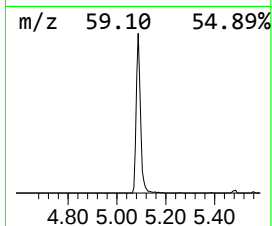
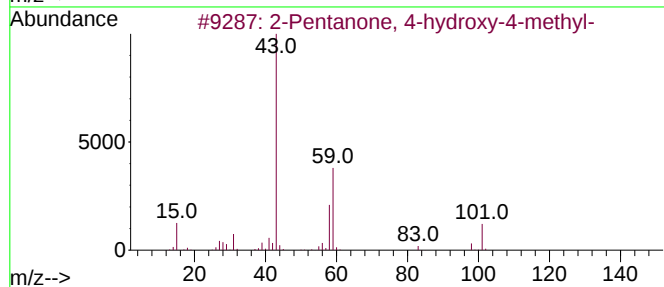
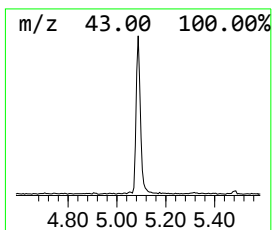
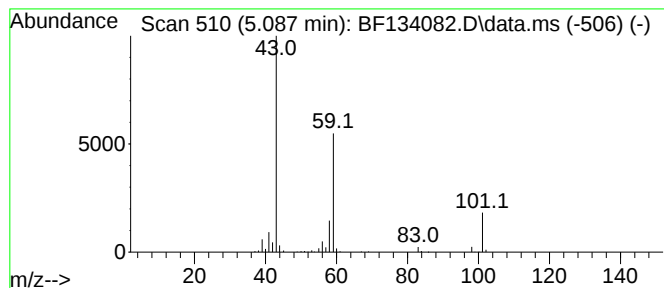
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.087	5.72 ng	235506	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	64		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
5	Oxirane, tetramethyl-	100	C6H12O	005076-20-0	9		



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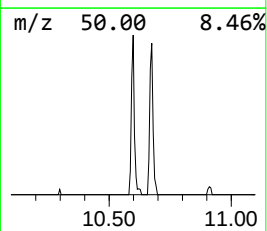
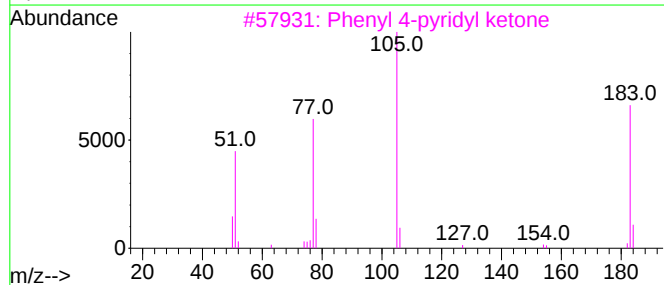
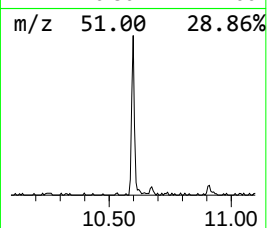
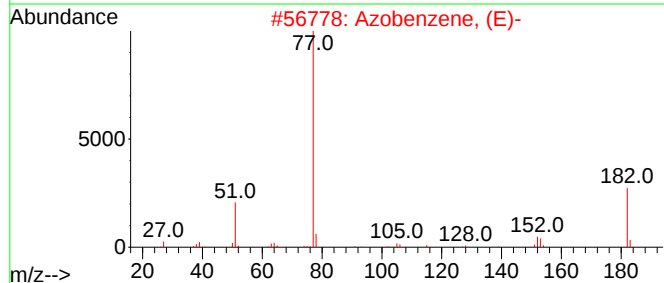
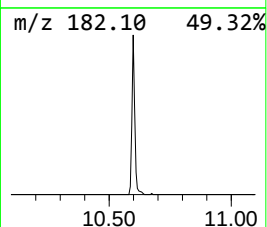
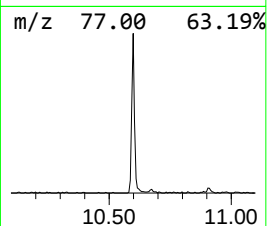
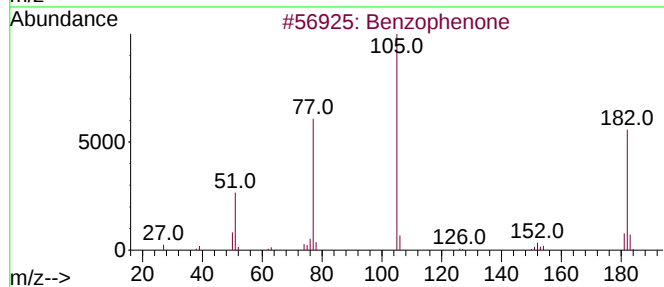
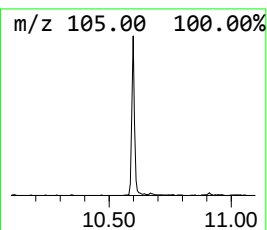
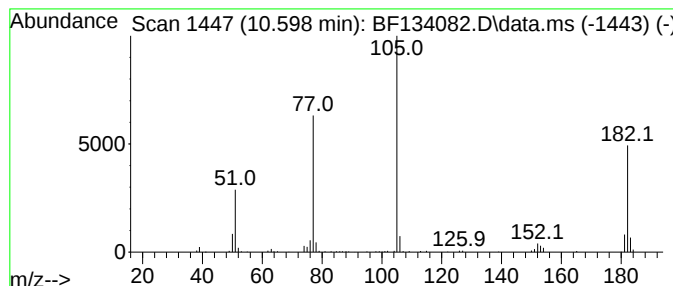
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Peak Number 3 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.598	2.84 ng	160001	Acenaphthene-d10	9.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Azobenzene, (E)-	182	C12H10N2	017082-12-1	64
3			Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	58
4			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	47
5			Vinyl benzoate	148	C9H8O2	000769-78-8	47



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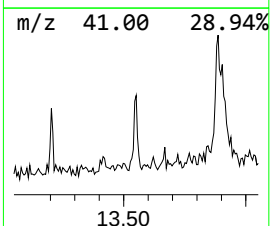
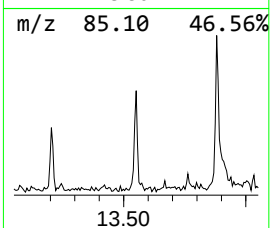
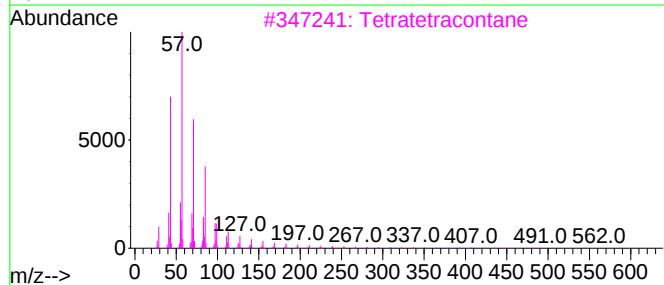
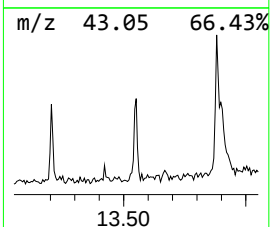
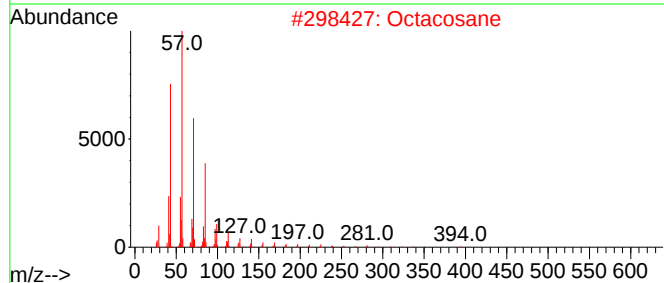
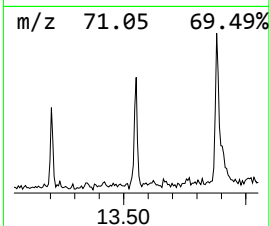
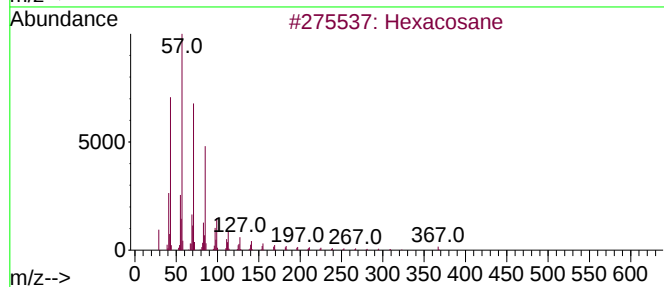
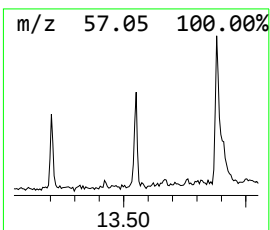
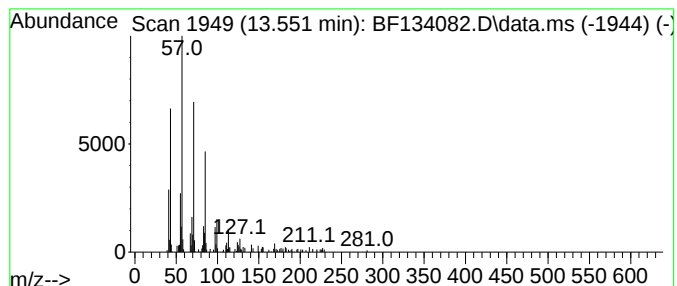
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Peak Number 4 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.551	2.01 ng	75025	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Octacosane	394	C28H58	000630-02-4	90
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Pentacosane	352	C25H52	000629-99-2	90
5			Nonadecane	268	C19H40	000629-92-5	87



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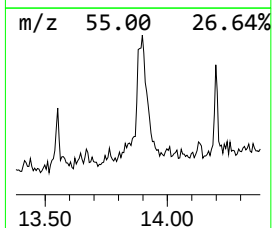
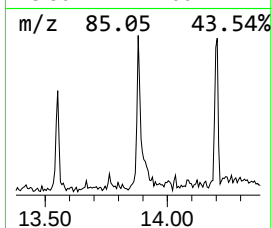
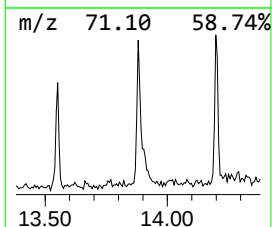
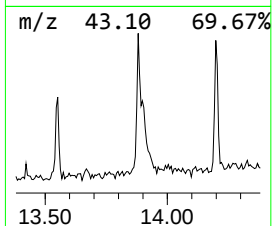
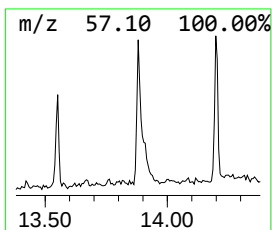
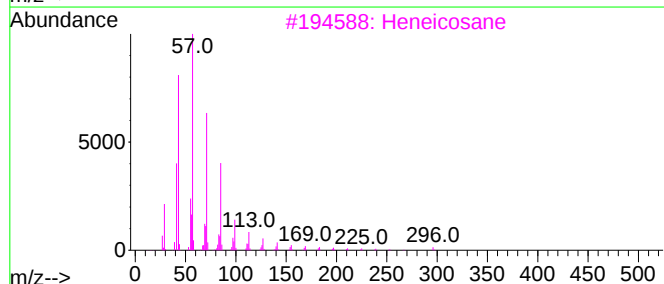
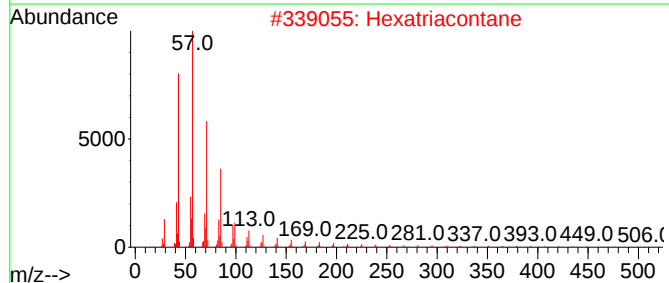
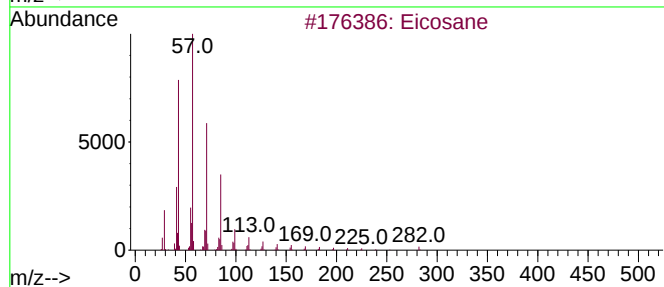
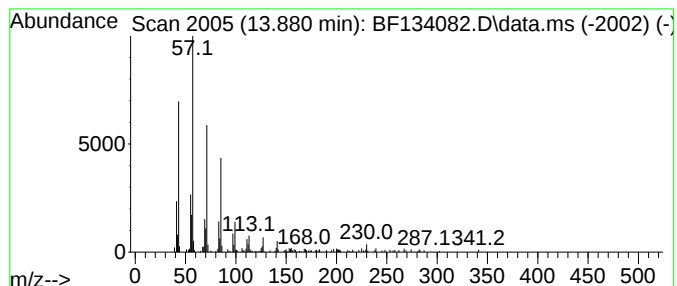
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Peak Number 5 Eicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.880	7.83 ng	291898	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Hexatriacontane	507	C36H74	000630-06-8	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			Hexacosane	366	C26H54	000630-01-3	87
5			Octadecane	254	C18H38	000593-45-3	83



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CMC-EME-TOPSOIL-008

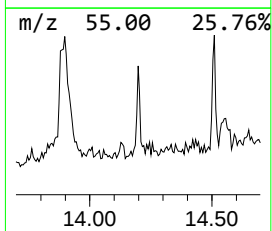
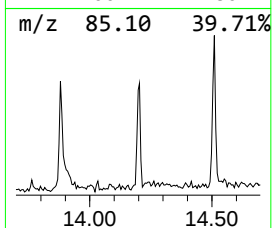
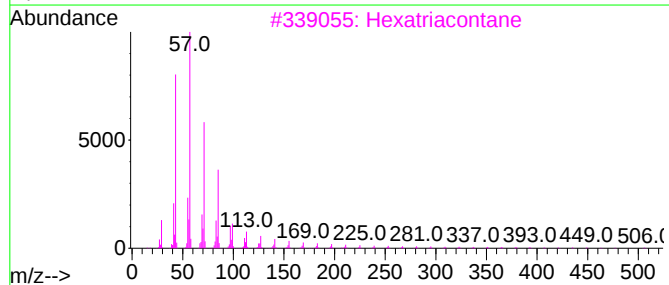
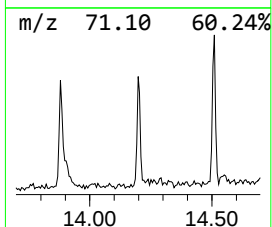
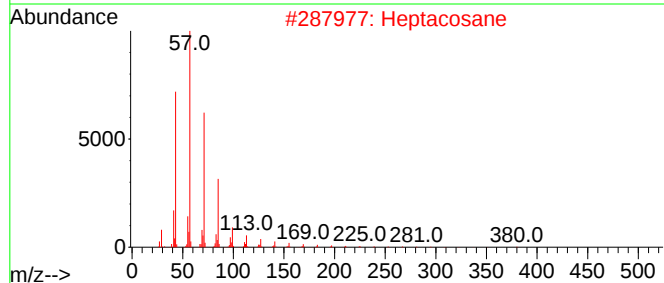
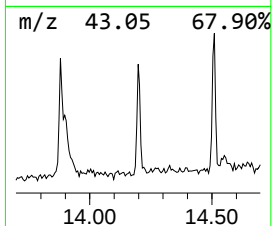
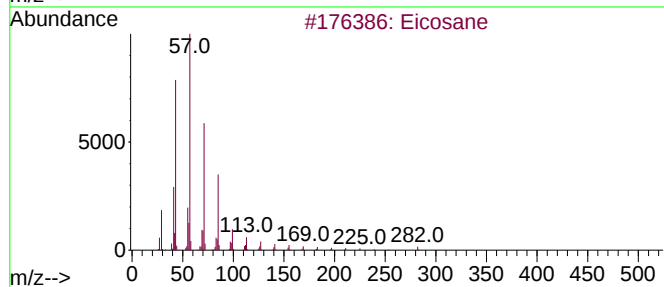
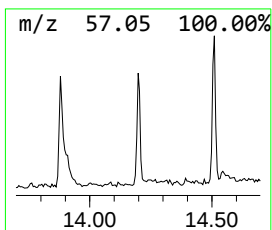
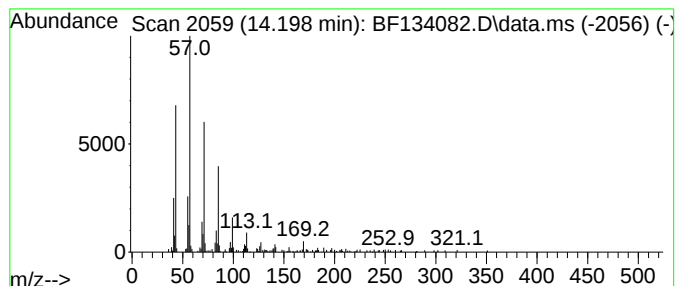
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.198	2.72 ng	101474	Chrysene-d12	14.039

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	87
2		Heptacosane	380	C27H56	000593-49-7	87
3		Hexatriacontane	507	C36H74	000630-06-8	87
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134082.D
Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CMC-EME-TOPSOIL-008

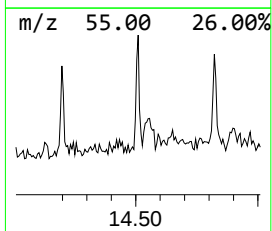
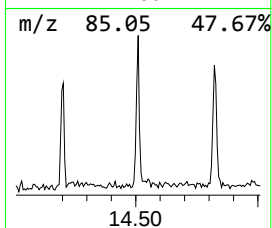
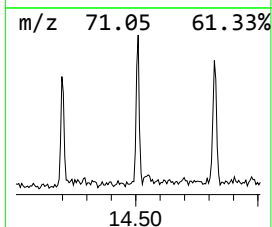
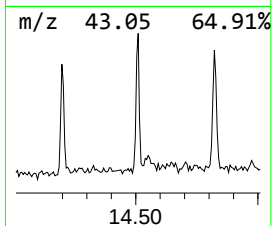
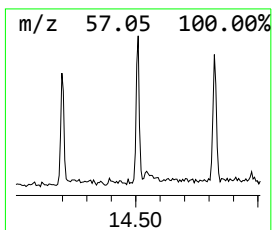
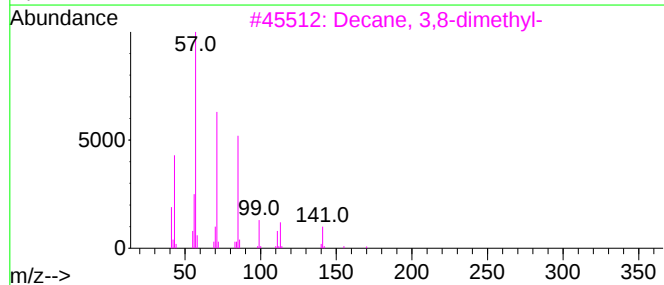
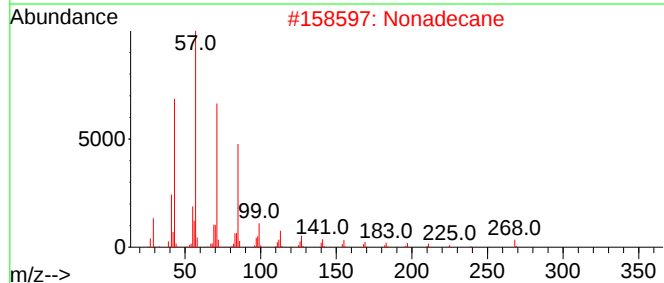
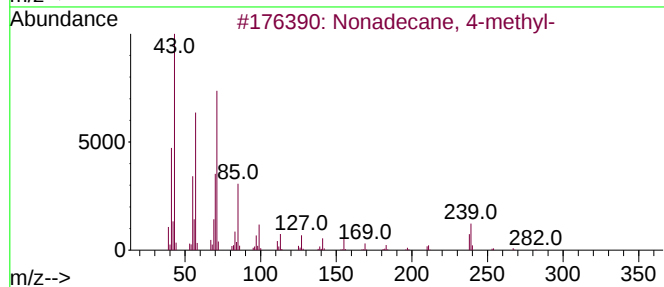
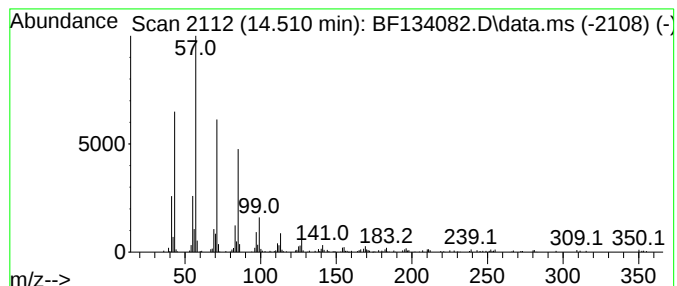
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.510	3.37 ng	125842	Chrysene-d12	14.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 4-methyl-	282	C20H42	025117-27-5	86
2			Nonadecane	268	C19H40	000629-92-5	83
3			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	83
4			Heneicosane	296	C21H44	000629-94-7	81
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134082.D
Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
Misc :
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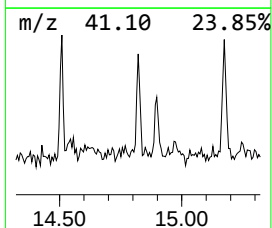
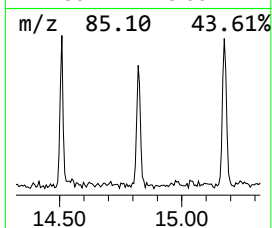
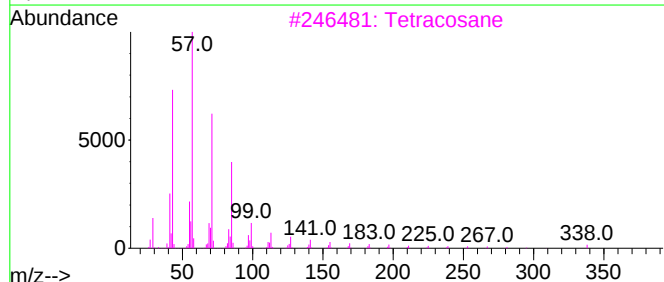
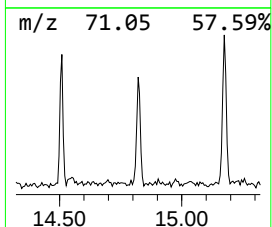
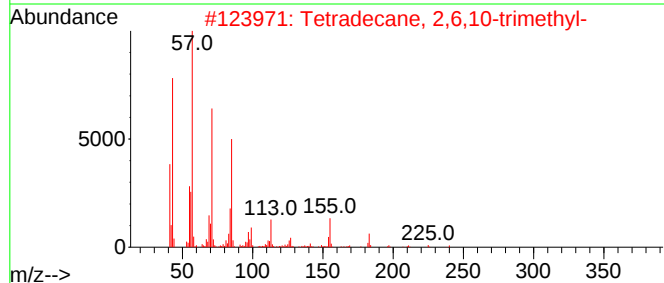
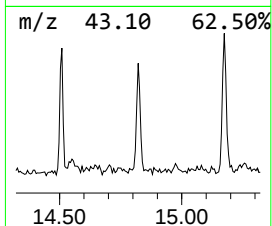
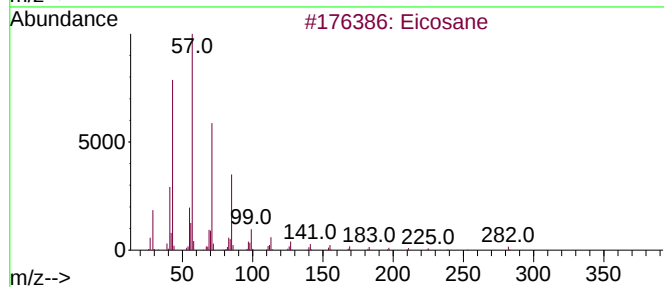
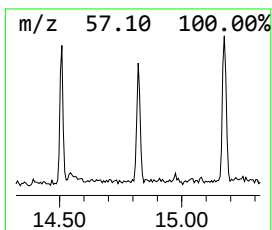
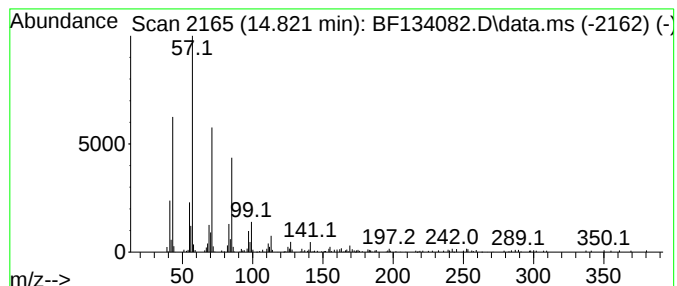
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetradecane, 2,6,10-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.821	4.29 ng	121796	Perylene-d12		15.539	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	90
3	Tetracosane		338	C24H50	000646-31-1	87
4	Octacosane		394	C28H58	000630-02-4	87
5	Heptadecane		240	C17H36	000629-78-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134082.D
Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CMC-EME-TOPSOIL-008

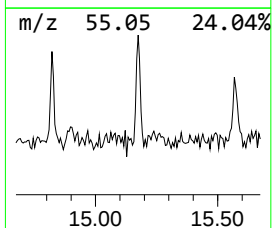
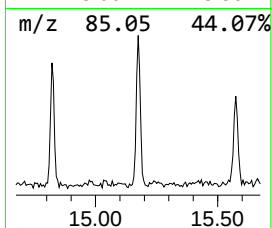
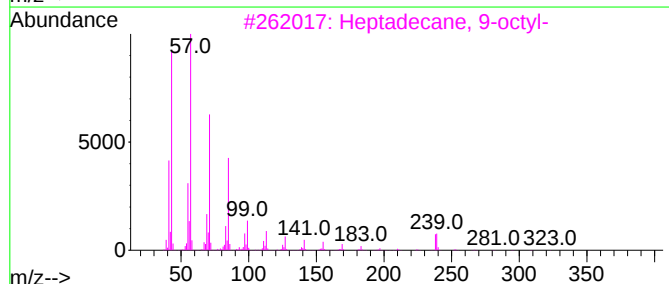
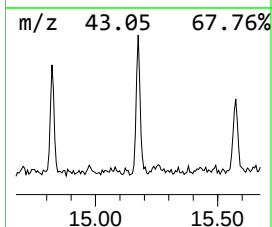
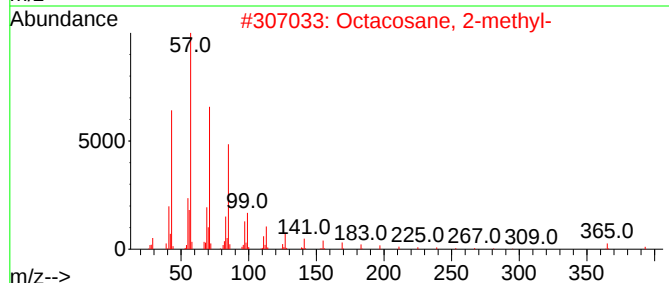
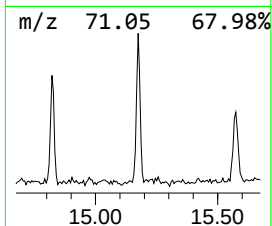
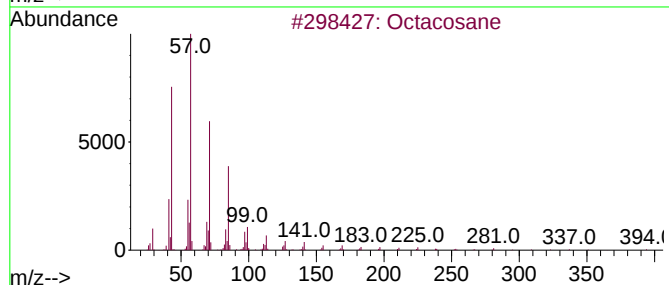
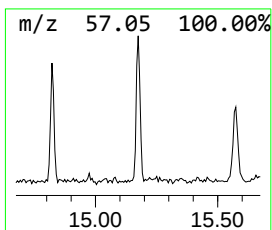
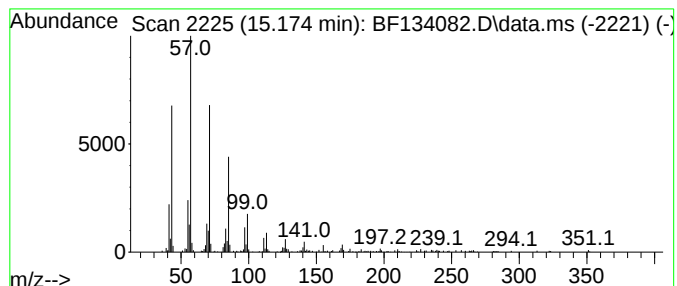
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.174	6.01 ng	170335	Perylene-d12	15.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	91
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Heptadecane	240	C17H36	000629-78-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134082.D
Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

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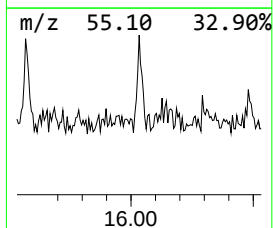
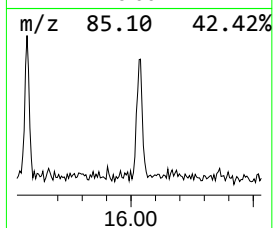
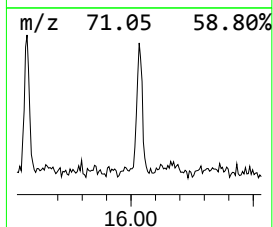
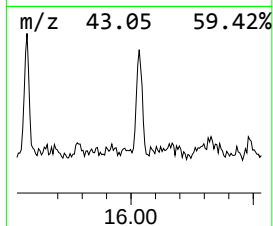
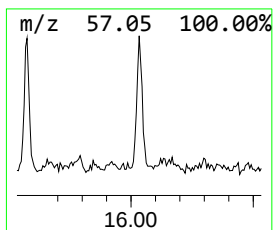
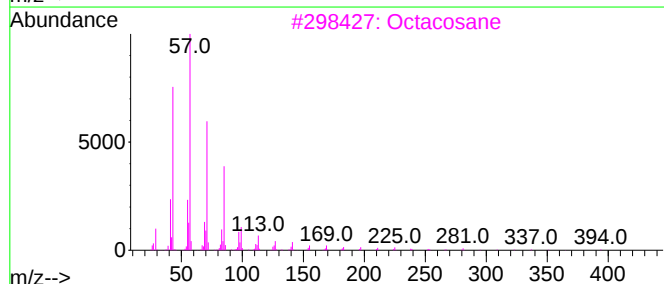
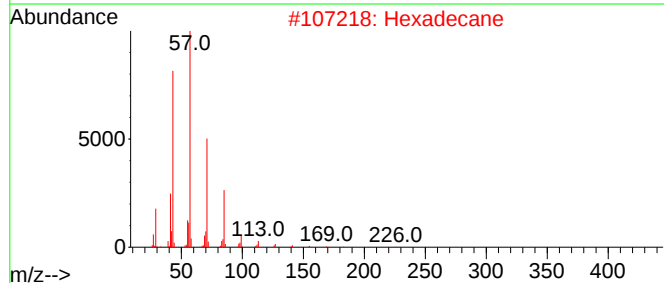
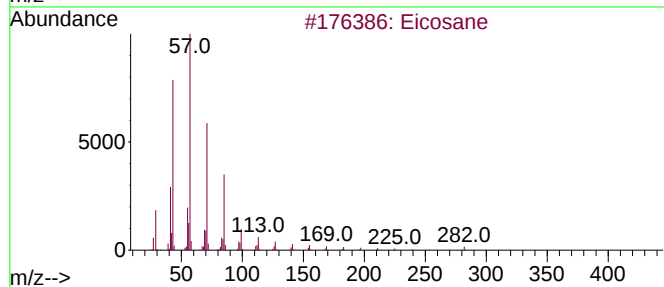
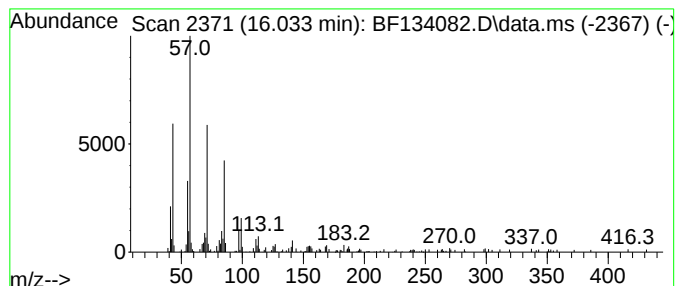
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.033	4.00 ng	113521	Perylene-d12	15.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	92
2		Hexadecane	226	C16H34	000544-76-3	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	86
5		Heptadecane	240	C17H36	000629-78-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
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Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
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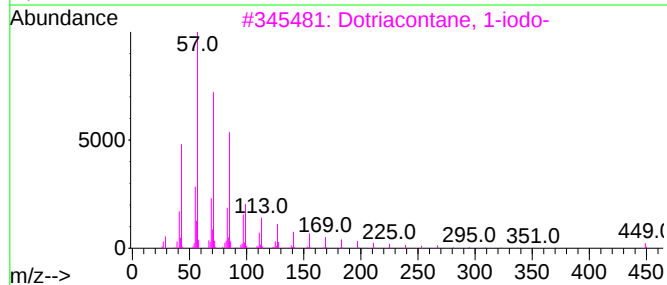
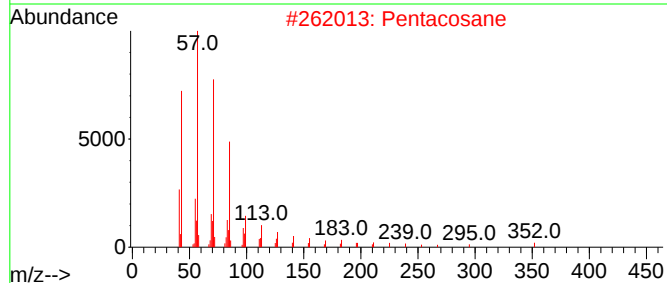
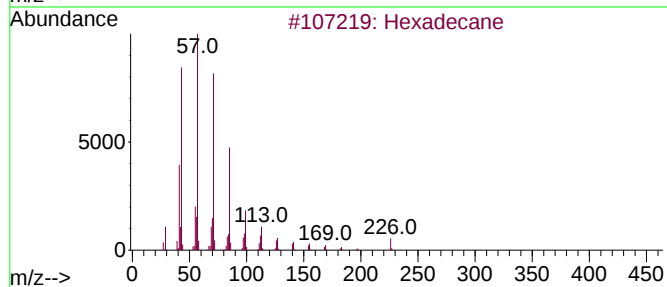
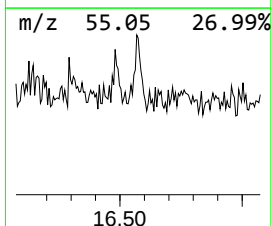
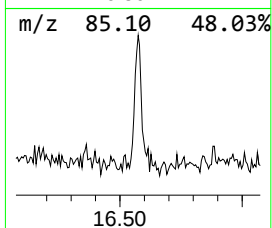
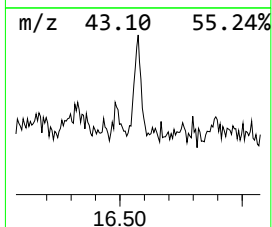
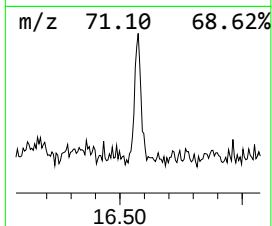
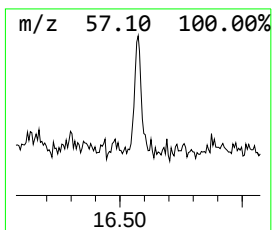
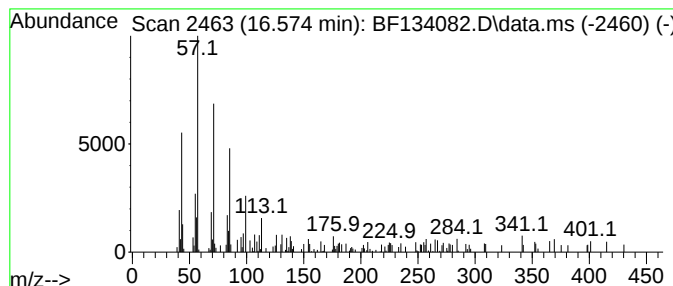
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pentacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.574	2.61 ng	74118	Perylene-d12	15.539	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	64
2	Pentacosane	352	C25H52	000629-99-2	64
3	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	59
4	Heneicosane	296	C21H44	000629-94-7	58
5	Hexacosane	366	C26H54	000630-01-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134082.D
Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CMC-EME-TOPSOIL-008

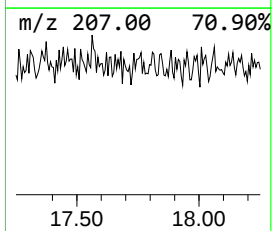
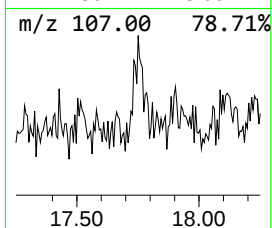
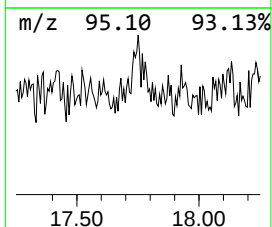
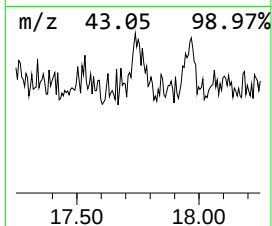
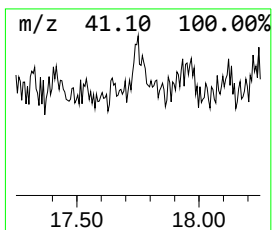
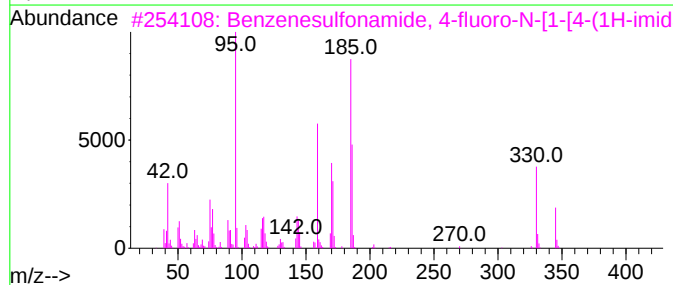
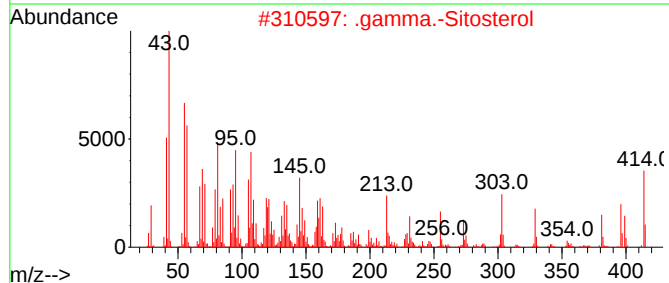
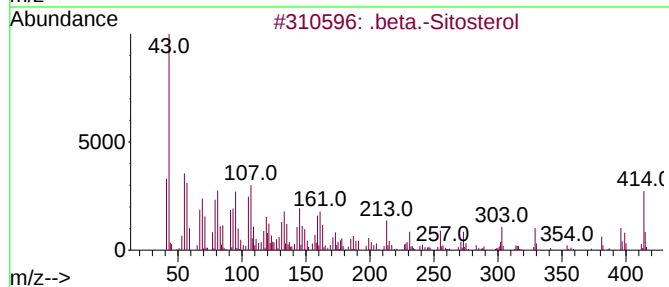
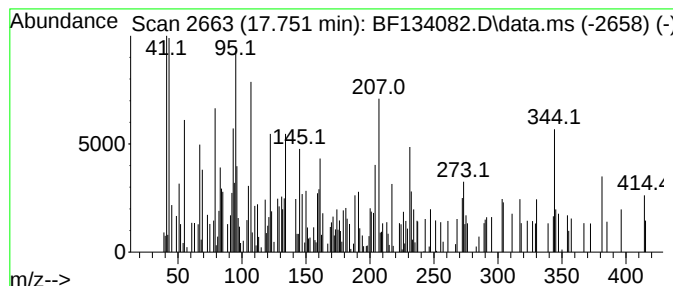
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown17.751 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.751	4.90 ng	139046	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-	Sitosterol	414	C29H50O	000083-46-5	38
2	.	gamma.-	Sitosterol	414	C29H50O	000083-47-6	38
3			Benzenesulfonamide, 4-fluoro-N-[...]	345	C17H16FN3O2S	1000318-81-2	25
4			5H,13H-Dibenzo[c,h]dipyrrolo[3,2...	344	C22H24N4	016739-54-1	22
5			Ethyl 5-(furan-2-yl)-1,2-oxazole...	207	C10H9N04	1000411-40-9	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070523\
Data File : BF134082.D
Acq On : 05 Jul 2023 23:56
Operator : CG\JU
Sample : 03473-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CMC-EME-TOPSOIL-008

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...	2.152	35.7	ng	1468520	1	6.845	823529	20.0
2-Pentanone, 4-...	5.087	5.7	ng	235506	1	6.845	823529	20.0
Benzophenone	10.598	2.8	ng	160001	3	9.881	1125820	20.0
Hexacosane	13.551	2.0	ng	75025	5	14.039	745736	20.0
Eicosane	13.880	7.8	ng	291898	5	14.039	745736	20.0
Heptacosane	14.198	2.7	ng	101474	5	14.039	745736	20.0
Nonadecane, 4-m...	14.510	3.4	ng	125842	5	14.039	745736	20.0
Tetradecane, 2,...	14.821	4.3	ng	121796	6	15.539	567194	20.0
Octacosane	15.174	6.0	ng	170335	6	15.539	567194	20.0
Hexadecane	16.033	4.0	ng	113521	6	15.539	567194	20.0
Pentacosane	16.574	2.6	ng	74118	6	15.539	567194	20.0
unknown17.751	17.751	4.9	ng	139046	6	15.539	567194	20.0