

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
 Data File : BF132307.D
 Acq On : 02 Feb 2023 23:24
 Operator : CG\JU
 Sample : 01371-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1-BERKSHIRE-NW

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-BF013123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF132307.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.313	415	421	430	rVB	1271458	1629326	34.39%	4.566%
2	5.695	480	486	489	rBV	3515509	3781292	79.80%	10.597%
3	6.678	647	653	656	rBV	3205254	3770087	79.57%	10.566%
4	7.066	715	719	722	rVB	1369751	1081593	22.83%	3.031%
5	7.625	809	814	818	rBV	2515873	2559805	54.02%	7.174%
6	8.025	877	882	890	rVB	58434	80884	1.71%	0.227%
7	8.348	934	937	941	rBV	1697715	1440622	30.40%	4.037%
8	9.430	1116	1121	1124	rBV	5261511	4738240	100.00%	13.279%
9	9.748	1171	1175	1177	rBV	231046	189844	4.01%	0.532%
10	10.113	1231	1237	1241	rBV	1804460	1607332	33.92%	4.504%
11	10.795	1349	1353	1355	rBV	94234	88366	1.86%	0.248%
12	10.830	1355	1359	1362	rVB	316900	271840	5.74%	0.762%
13	10.907	1368	1372	1375	rBV	3056208	2654937	56.03%	7.440%
14	11.613	1488	1492	1495	rBV	1751467	1439871	30.39%	4.035%
15	11.636	1495	1496	1500	rVB	122461	72146	1.52%	0.202%
16	12.124	1575	1579	1584	rBV	744938	718173	15.16%	2.013%
17	12.836	1695	1700	1709	rVB	233835	270788	5.71%	0.759%
18	12.913	1710	1713	1721	rBV	227073	211397	4.46%	0.592%
19	13.065	1733	1739	1745	rBV	166639	198398	4.19%	0.556%
20	13.207	1759	1763	1766	rBV	3699794	3270930	69.03%	9.167%
21	13.642	1833	1837	1840	rBV2	207633	211159	4.46%	0.592%
22	14.101	1912	1915	1929	rBV2	341440	616822	13.02%	1.729%
23	14.277	1940	1945	1947	rBV	912064	916415	19.34%	2.568%
24	14.301	1947	1949	1957	rVB2	71666	100287	2.12%	0.281%
25	14.730	2020	2022	2024	rBV	65619	56359	1.19%	0.158%
26	14.760	2024	2027	2032	rVB4	68932	106865	2.26%	0.299%
27	15.154	2091	2094	2098	rVB	283419	294129	6.21%	0.824%
28	15.230	2104	2107	2111	rVB	187157	201987	4.26%	0.566%
29	15.383	2129	2133	2136	rBV	60443	96123	2.03%	0.269%
30	15.448	2140	2144	2147	rVV	117922	140514	2.97%	0.394%
31	15.495	2147	2152	2166	rVV3	178882	505239	10.66%	1.416%
32	15.789	2199	2202	2208	rVB	52787	68695	1.45%	0.193%
33	15.865	2209	2215	2226	rVB	782316	1107637	23.38%	3.104%
34	16.101	2251	2255	2262	rVB	280505	389451	8.22%	1.091%
35	16.365	2295	2300	2307	rVB	141070	218035	4.60%	0.611%
36	16.448	2308	2314	2319	rBV3	98829	269151	5.68%	0.754%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.M
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37	16.983	2401	2405	2413	rVB	68667	127514	2.69%	0.357%
38	17.271	2451	2454	2464	rVB5	55018	102427	2.16%	0.287%
39	17.712	2525	2529	2533	rBV7	40440	78149	1.65%	0.219%

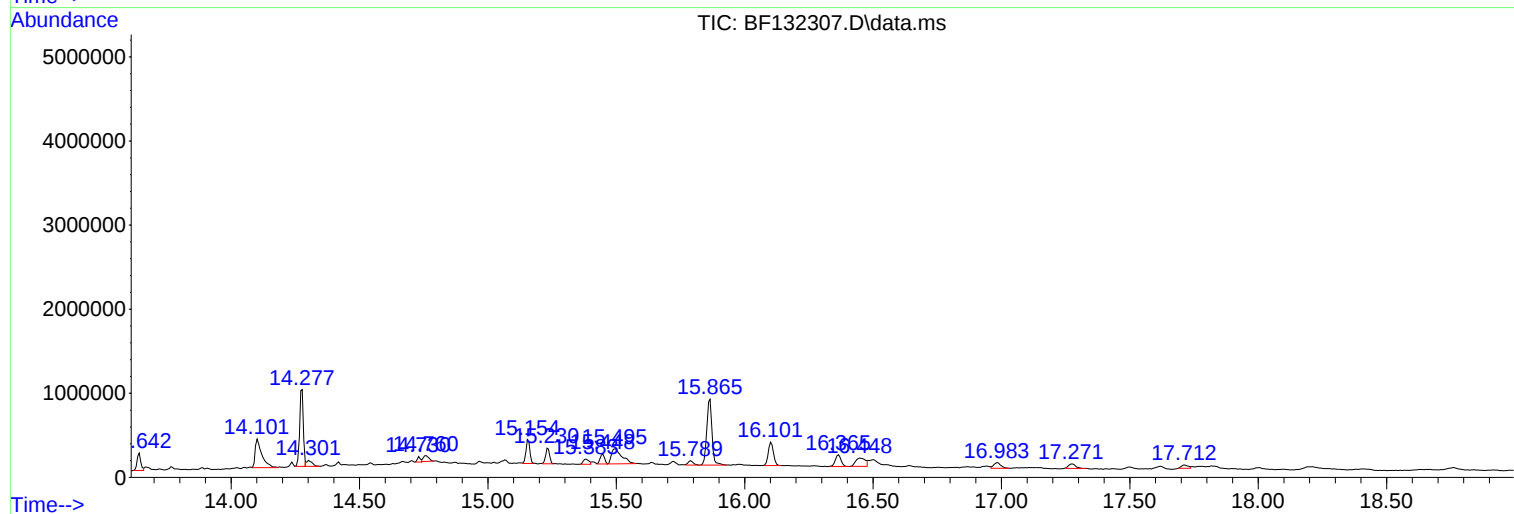
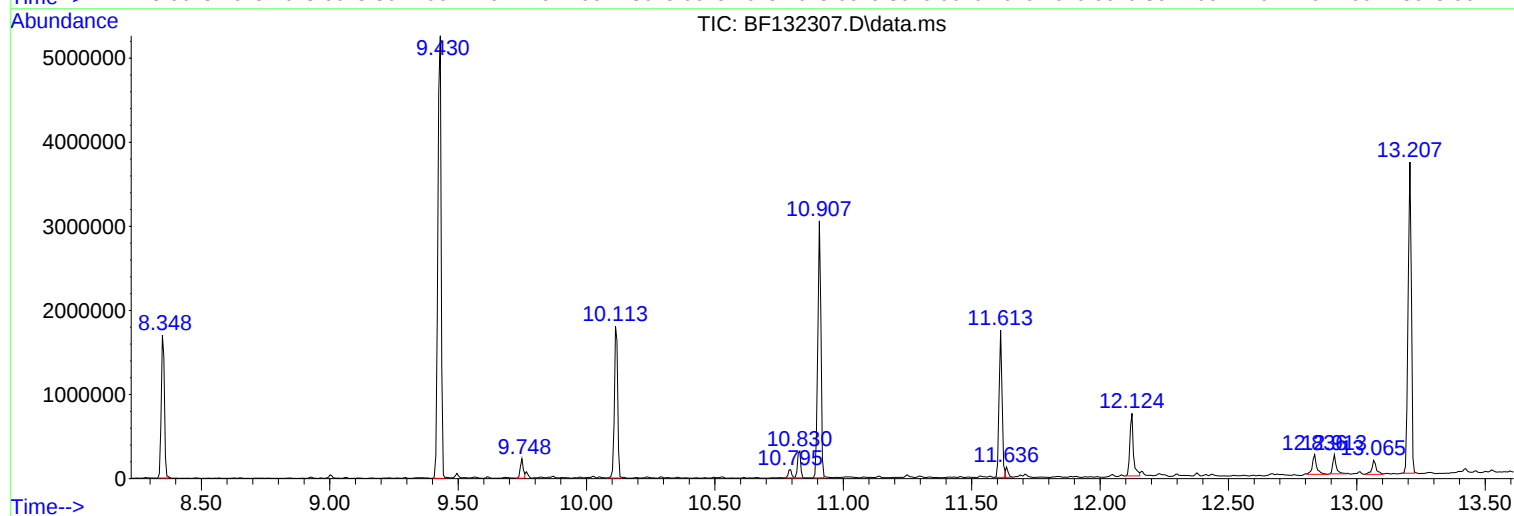
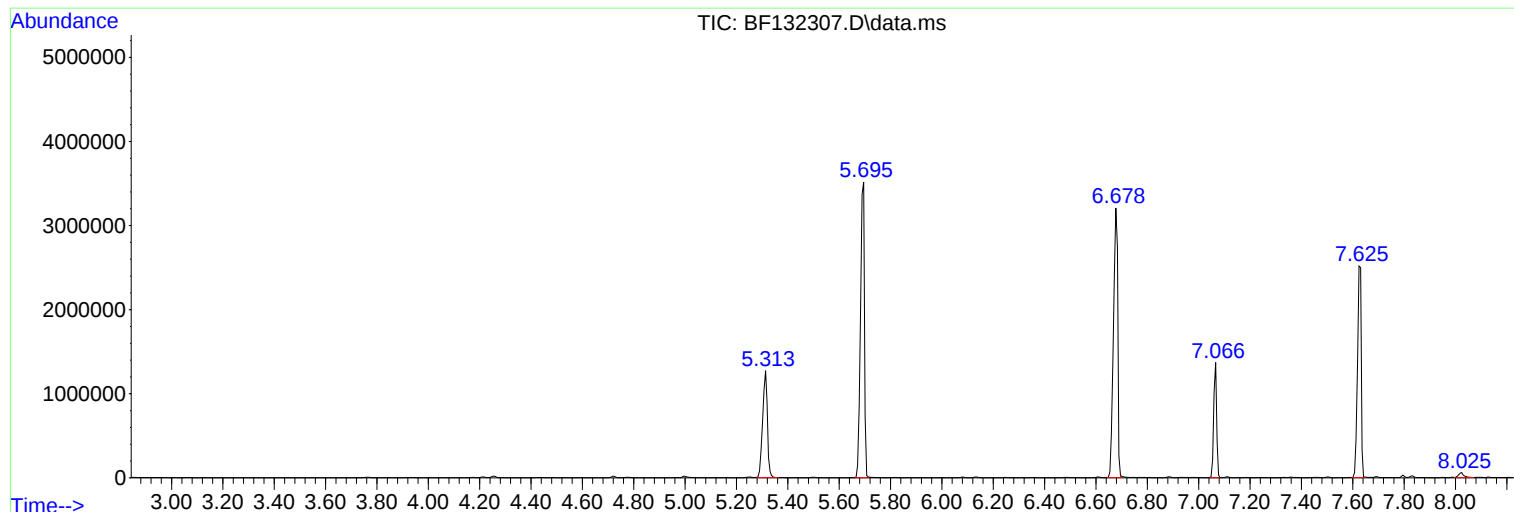
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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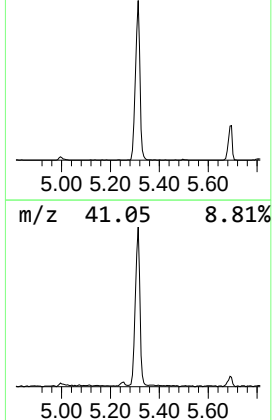
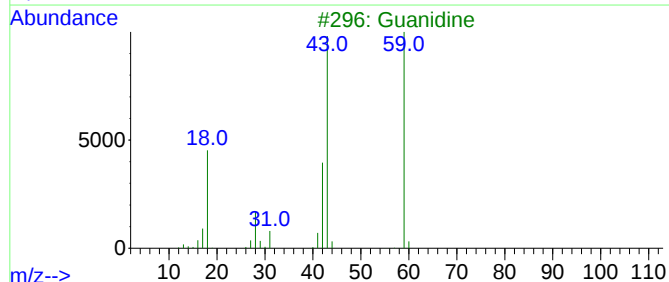
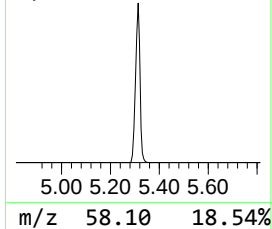
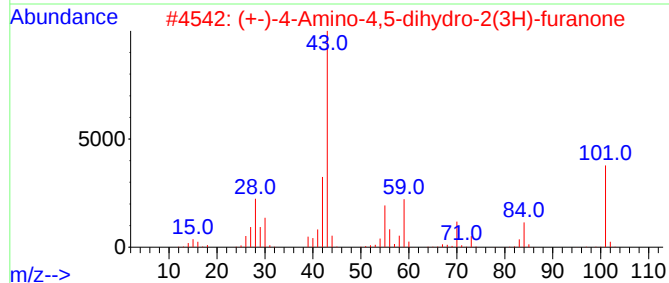
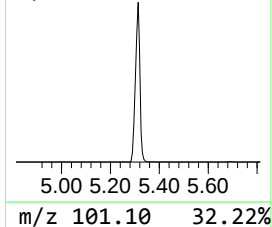
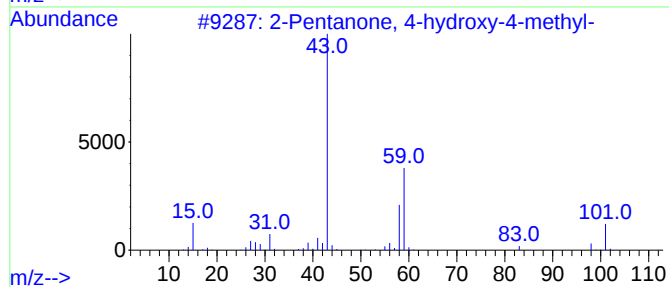
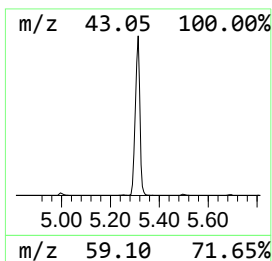
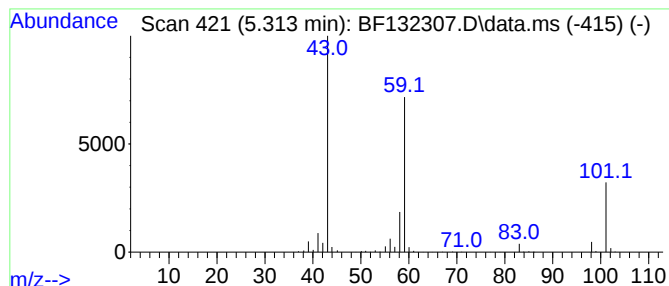
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.313	30.13 ng	1629330	1,4-Dichlorobenzene-d4	7.066	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3	Guanidine	59	CH5N3	000113-00-8	43
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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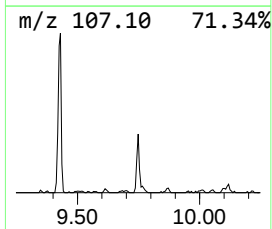
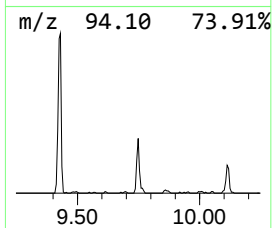
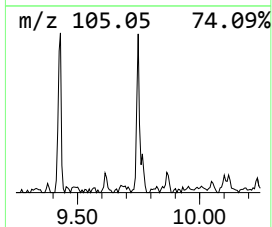
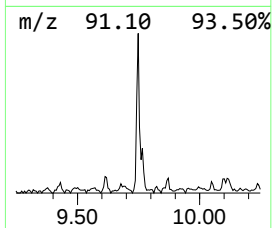
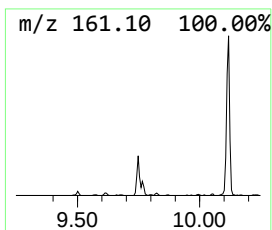
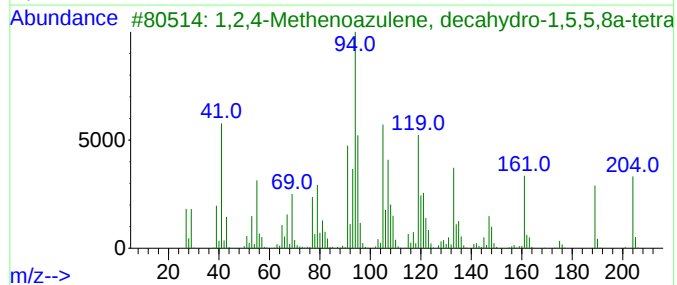
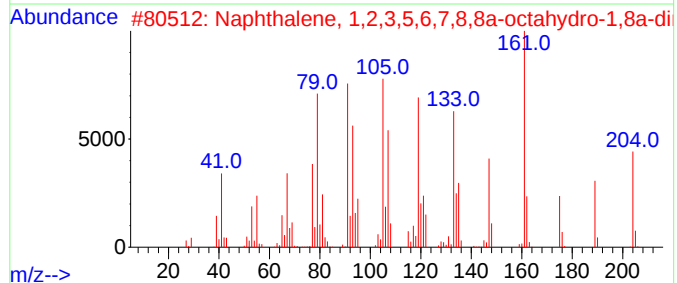
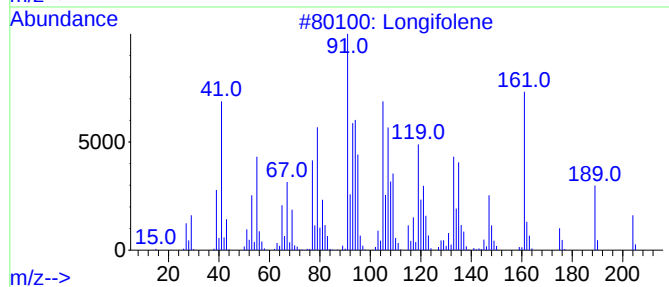
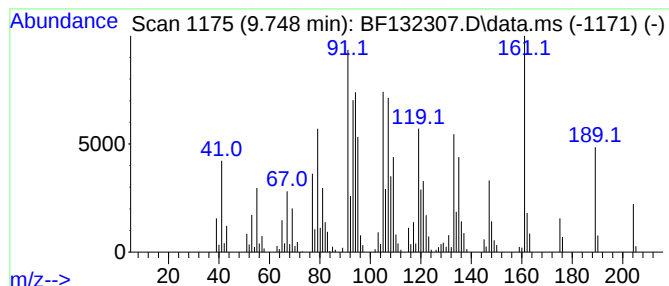
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Peak Number 2 Longifolene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.748	2.36 ng	189844	Acenaphthene-d10	10.113

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Longifolene	204	C15H24	000475-20-7	99
2			Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	91
3			1,2,4-Methenoazulene, decahydro-...	204	C15H24	001137-12-8	91
4			.beta.-Neoclovene	204	C15H24	056684-96-9	90
5			Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	010219-75-7	87



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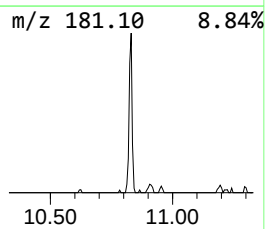
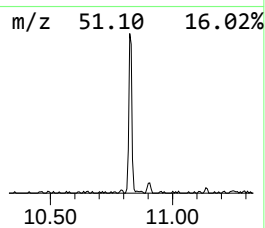
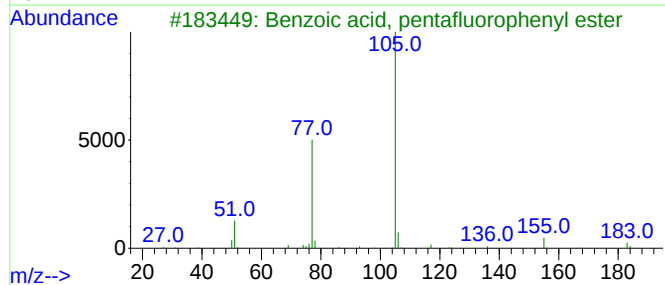
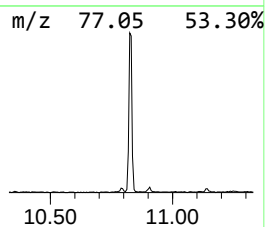
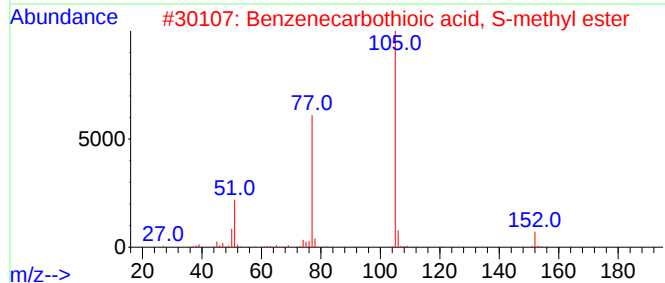
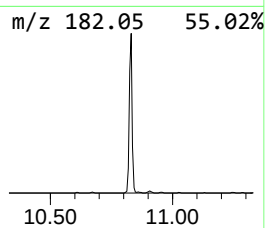
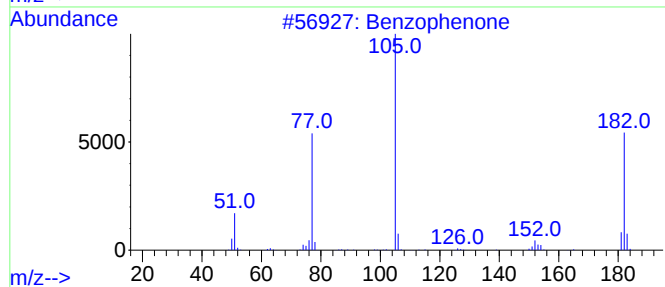
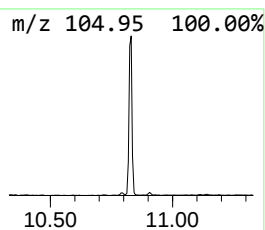
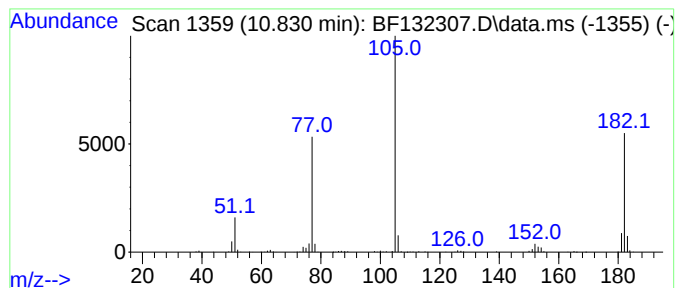
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzophenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.830	3.38 ng	271840	Acenaphthene-d10	10.113

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	47
3			Benzoic acid, pentafluorophenyl ...	288	C13H5F5O2	1000360-67-9	47
4			2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47
5			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	43



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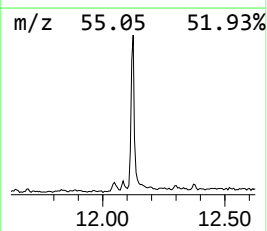
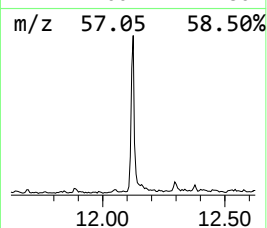
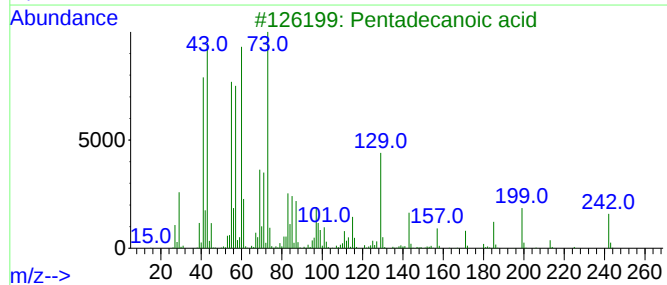
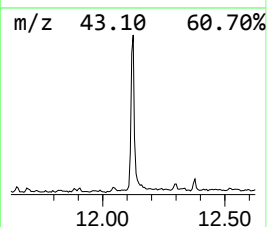
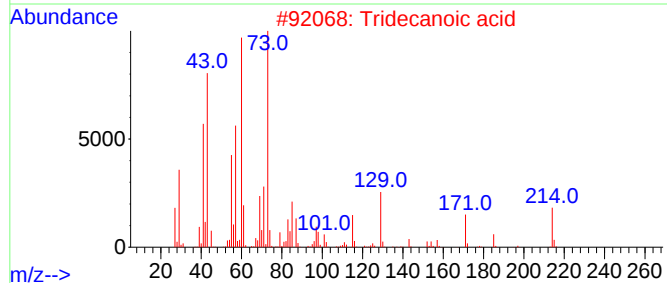
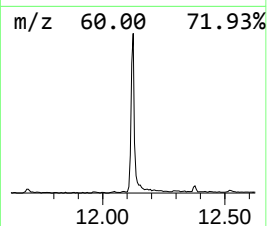
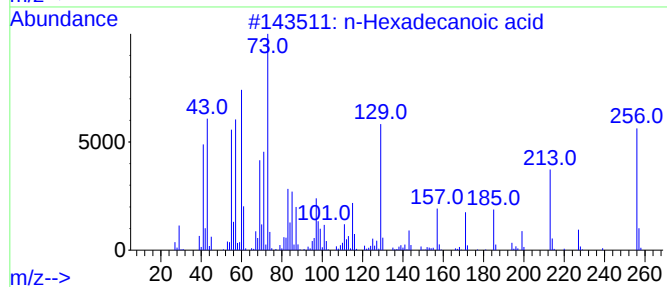
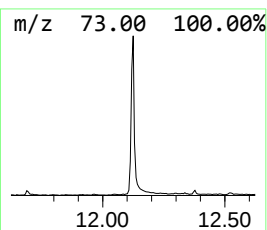
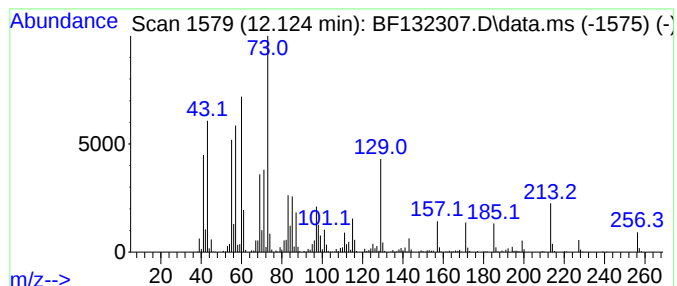
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.124	9.98 ng	718173	Phenanthrene-d10	11.613

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	90
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4			n-Decanoic acid	172	C10H20O2	000334-48-5	58
5			Dodecanoic acid	200	C12H24O2	000143-07-7	45



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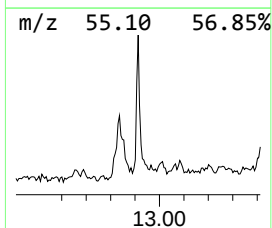
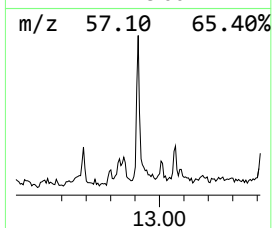
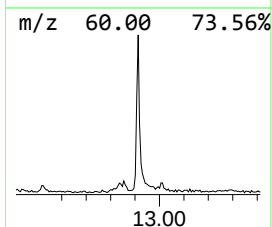
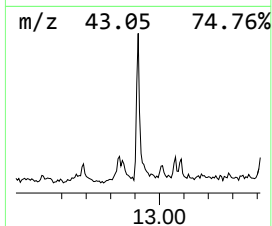
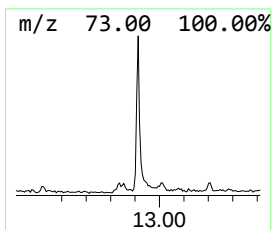
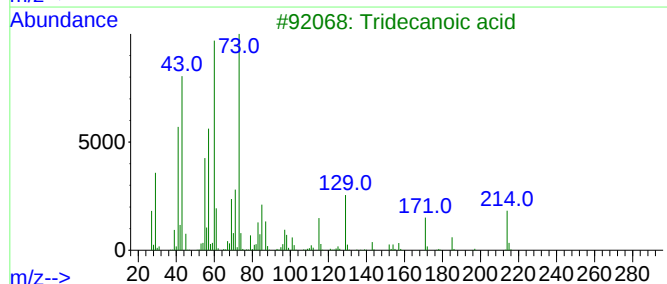
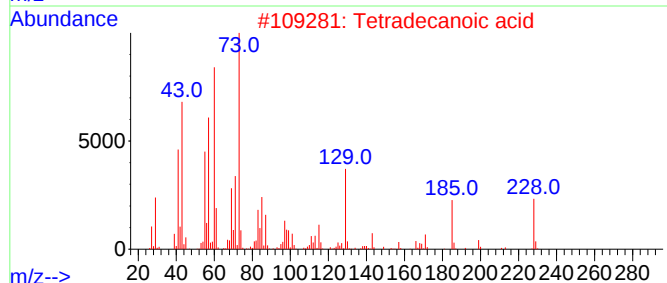
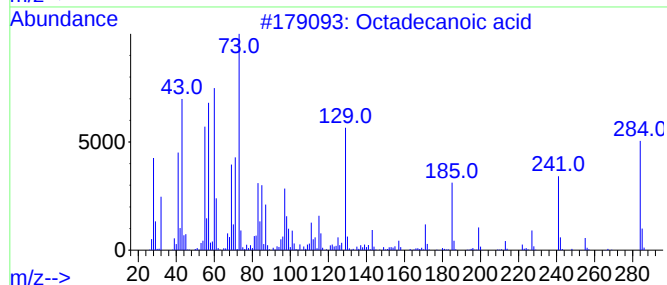
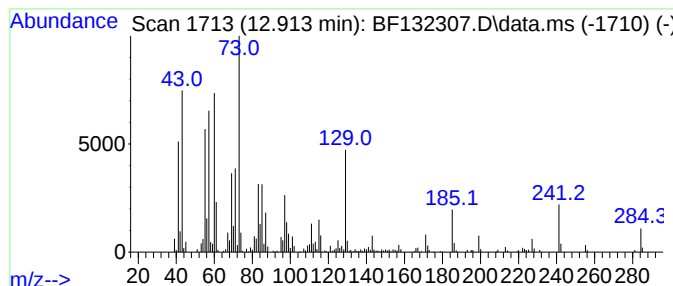
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.913	2.94 ng	211397	Phenanthrene-d10	11.613

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3			Tridecanoic acid	214	C13H26O2	000638-53-9	70
4			n-Decanoic acid	172	C10H20O2	000334-48-5	64
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
Data File : BF132307.D
Acq On : 02 Feb 2023 23:24
Operator : CG\JU
Sample : 01371-01
Misc :
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ClientSampleId :
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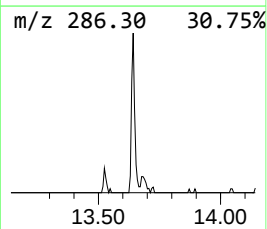
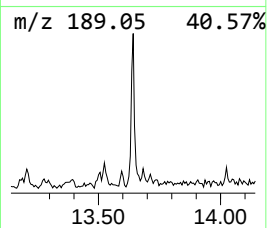
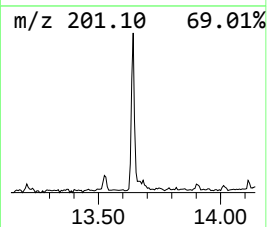
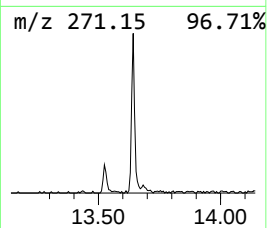
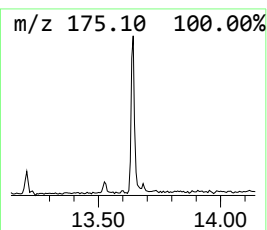
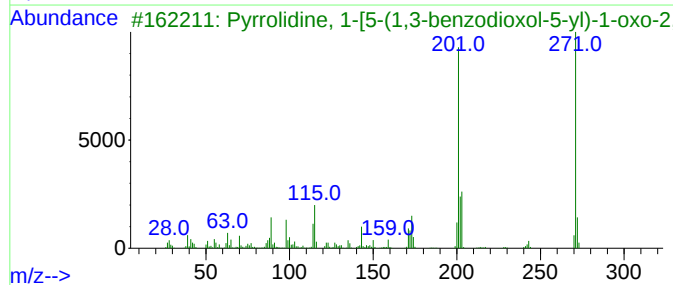
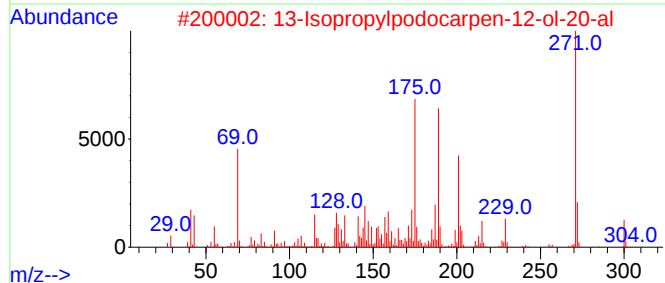
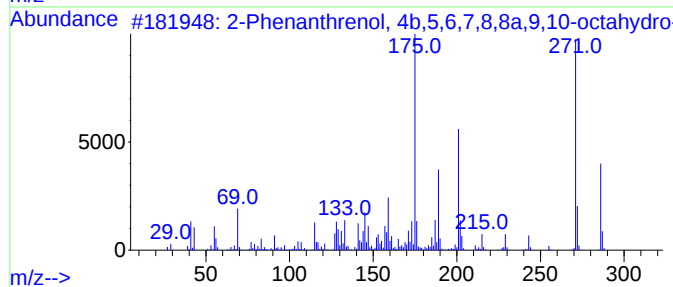
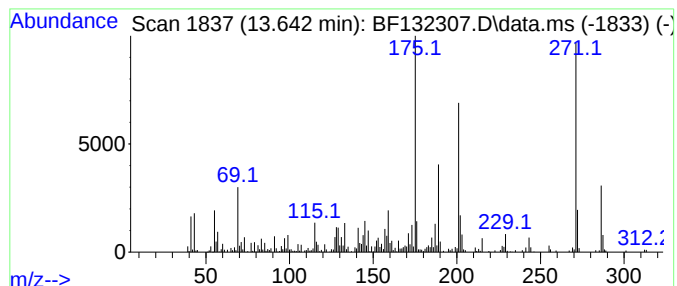
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.642	4.61 ng	211159	Chrysene-d12	14.277

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	99		
2	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	41		
3	Pyrrolidine, 1-[5-(1,3-benzodiox...	271	C16H17NO3	025924-78-1	30		
4	3-Cyano-6-trifluoromethylphenant...	271	C16H8F3N	1000253-90-3	22		
5	2,2-Bis(4'-methoxyphenyl)-2-etho...	286	C18H22O3	1000283-53-7	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
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ClientSampleId :
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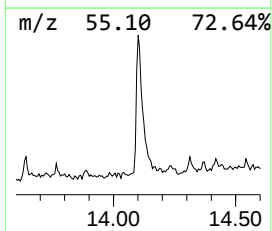
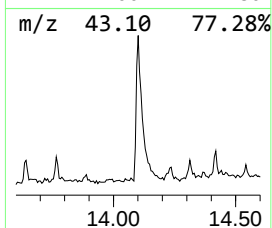
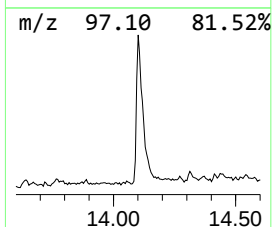
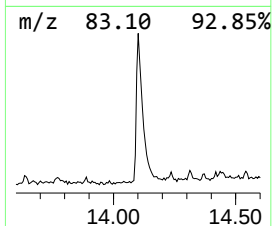
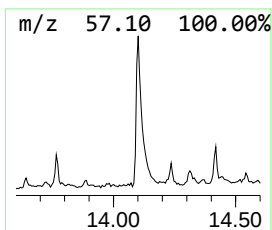
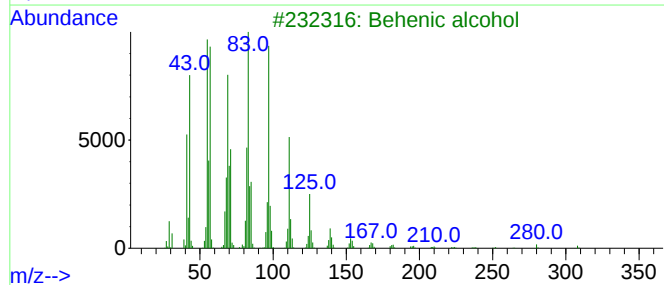
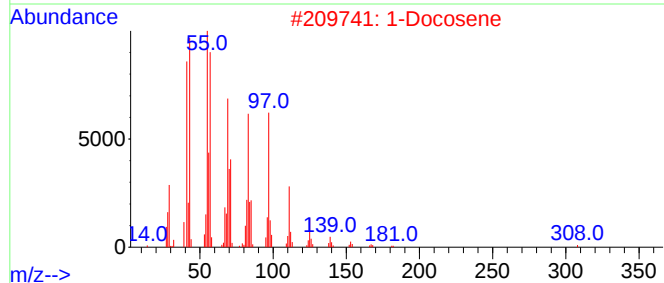
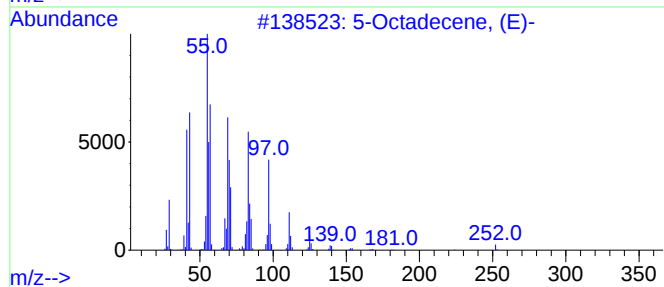
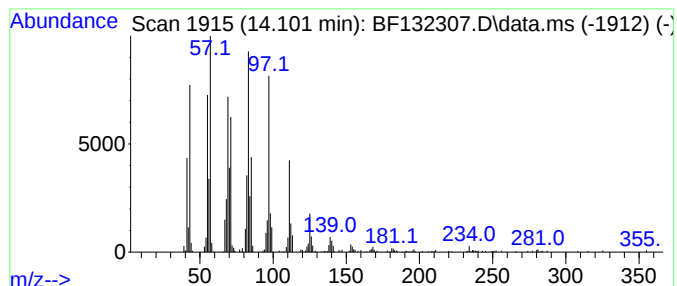
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 5-Octadecene, (E)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.101	13.46 ng	616822	Chrysene-d12	14.277

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	93
2			1-Docosene	308	C22H44	001599-67-3	91
3			Behenic alcohol	326	C22H46O	000661-19-8	91
4			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	90
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



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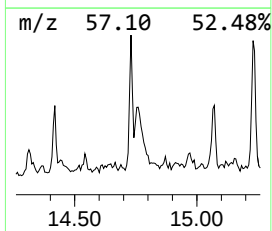
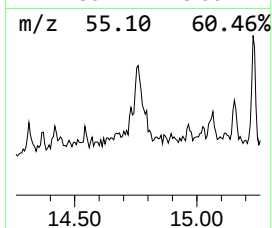
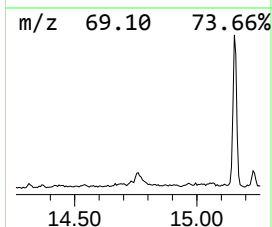
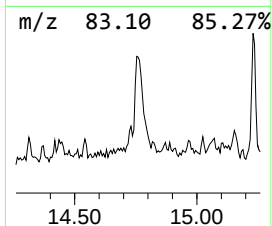
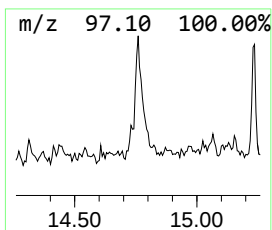
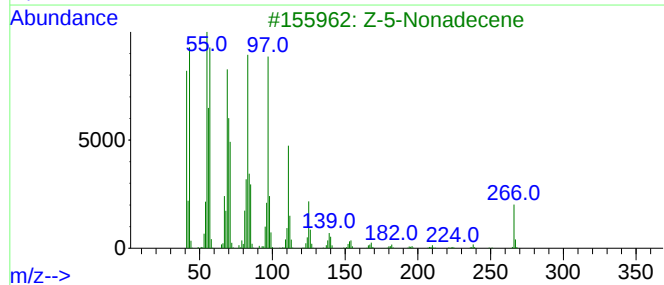
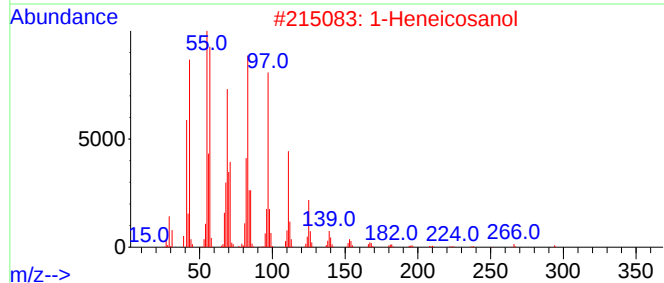
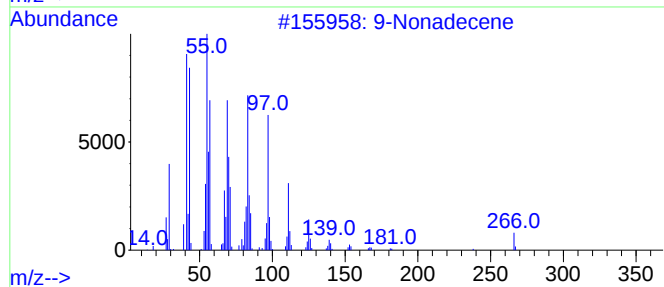
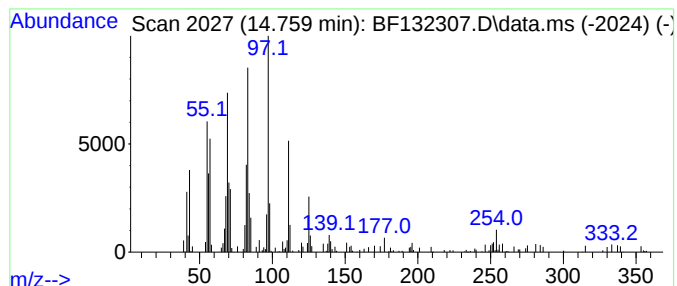
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Nonadecene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.759	2.33 ng	106865	Chrysene-d12		14.277	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9-Nonadecene		266	C19H38	031035-07-1	91
2	1-Heneicosanol		312	C21H44O	015594-90-8	91
3	Z-5-Nonadecene		266	C19H38	1000131-11-8	86
4	n-Nonadecanol-1		284	C19H40O	001454-84-8	76
5	Cyclohexadecane		224	C16H32	000295-65-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
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Instrument :
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ClientSampleId :
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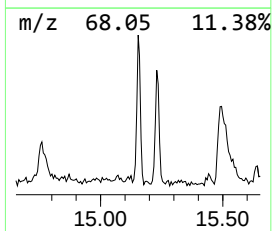
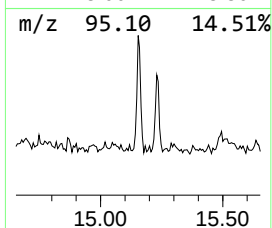
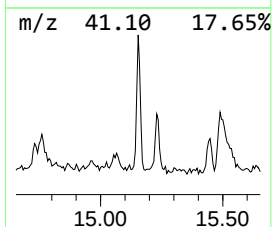
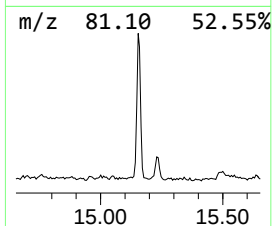
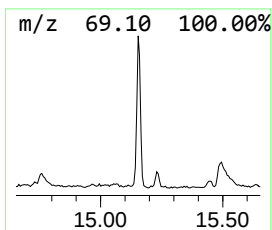
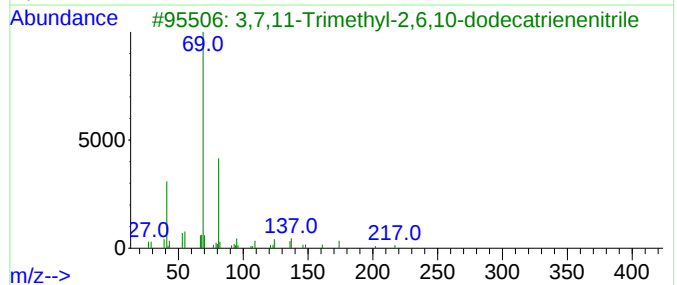
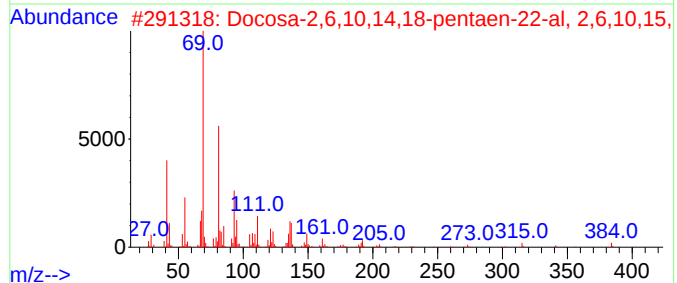
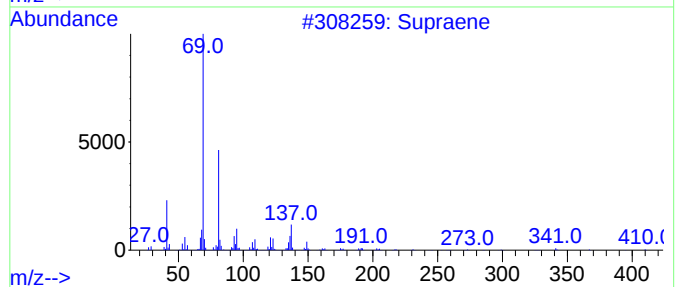
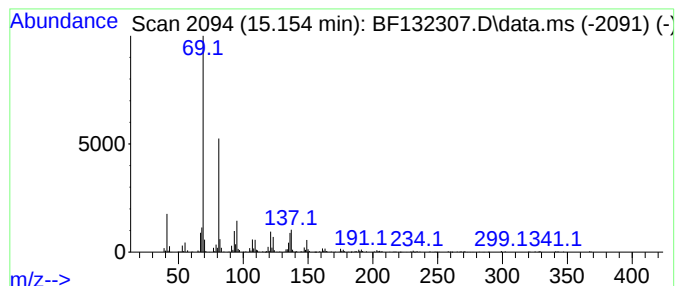
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.154	5.31 ng	294129	Perylene-d12	15.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	90
3			3,7,11-Trimethyl-2,6,10-dodecatr...	217	C15H23N	029789-67-1	59
4			Bis(3-methylbut-3-enyl) phthalate	302	C18H22O4	022711-53-1	53
5			1,6-Octadiene, 3,5-dimethyl-, tr...	138	C10H18	074630-87-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
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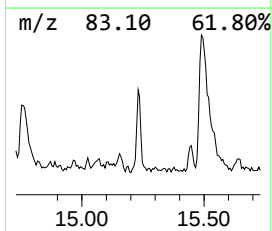
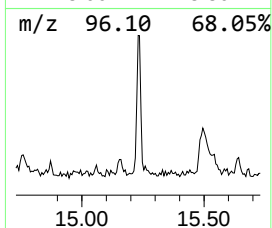
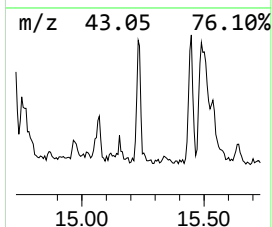
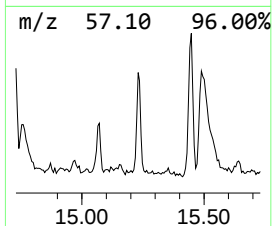
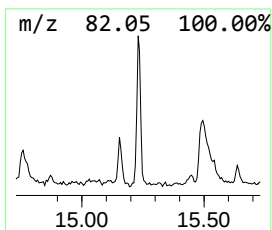
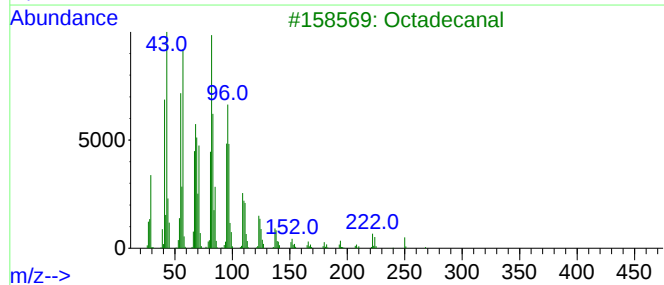
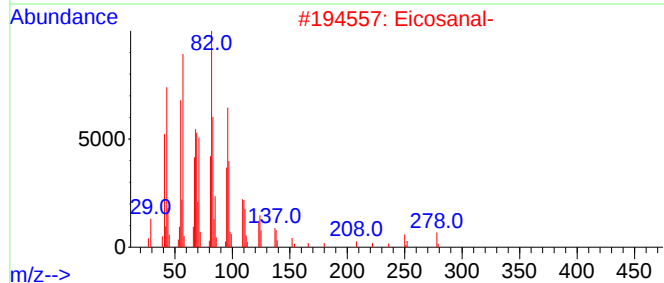
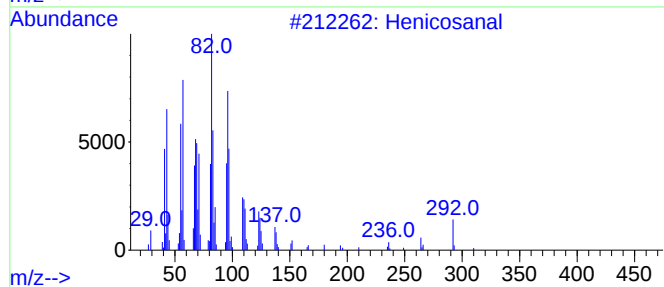
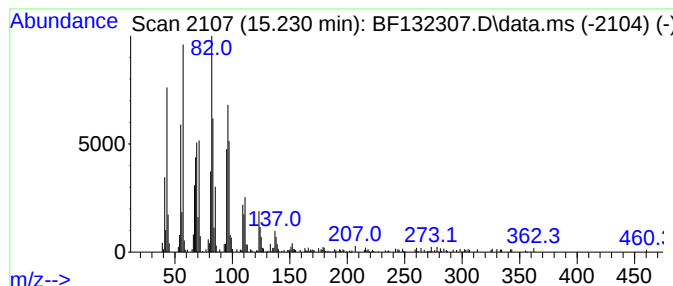
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Henicosanal Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.230	3.65 ng	201987	Perylene-d12	15.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Henicosanal	310	C21H42O	051227-32-8	87
2			Eicosanal-	296	C20H40O	002400-66-0	87
3			Octadecanal	268	C18H36O	000638-66-4	83
4			Hexacosanal	380	C26H52O	026627-85-0	83
5			Undec-10-ynoic acid, undecyl ester	336	C22H40O2	1000406-16-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
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ClientSampleId :
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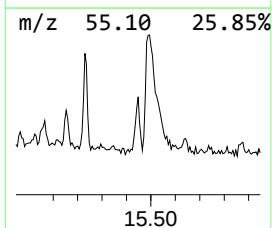
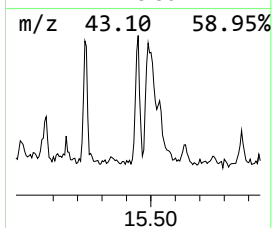
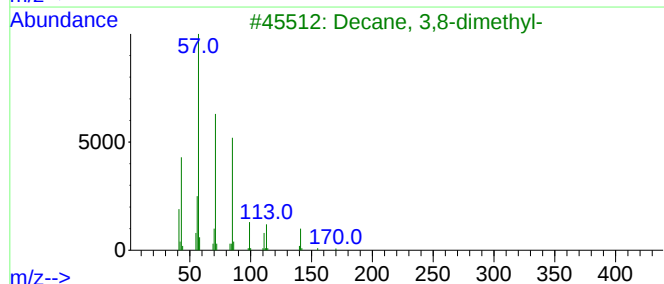
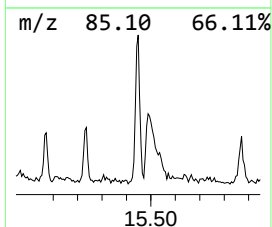
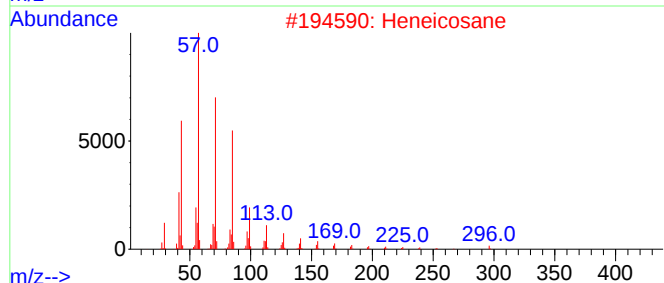
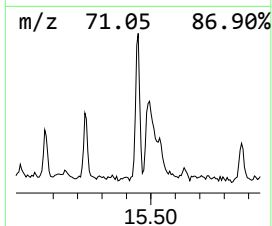
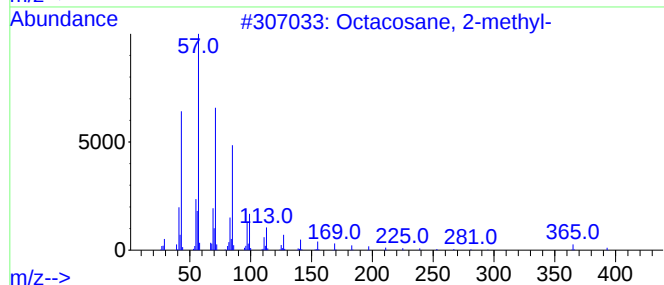
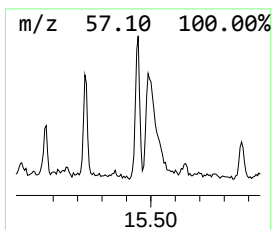
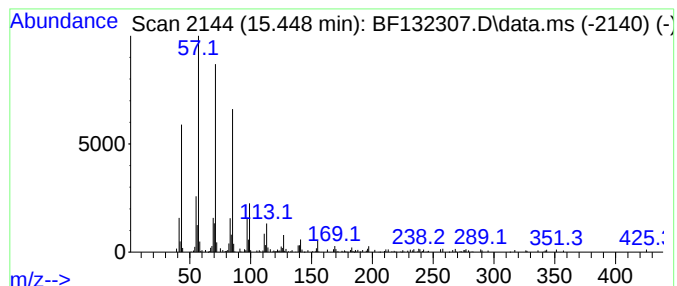
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.448	2.54 ng	140514	Perylene-d12	15.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87
4			Hexacosane	366	C26H54	000630-01-3	87
5			Hexadecane	226	C16H34	000544-76-3	86



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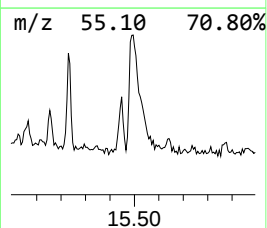
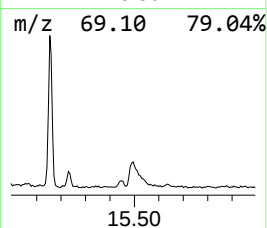
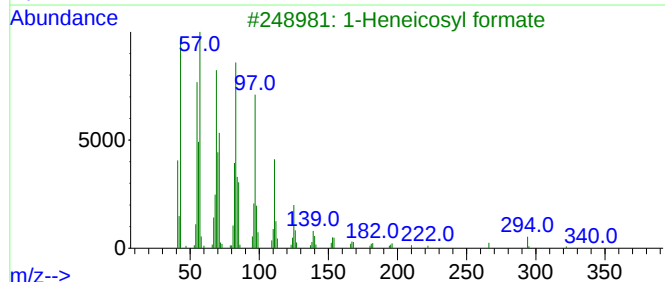
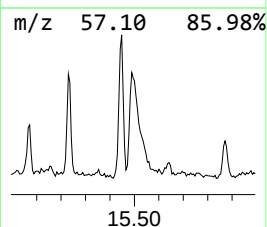
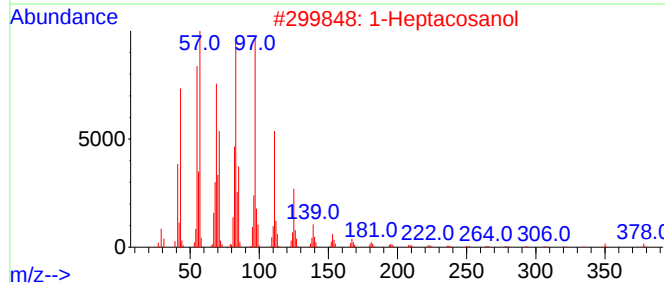
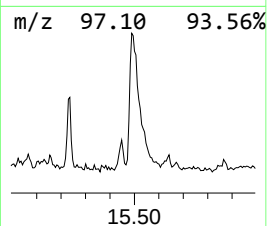
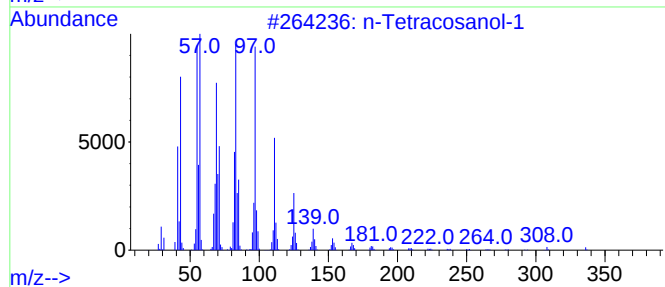
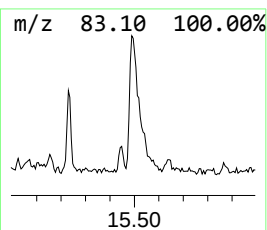
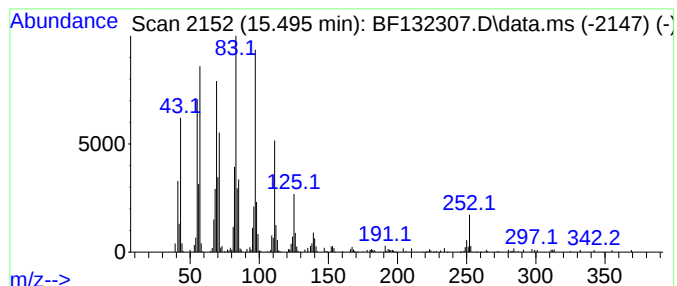
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 n-Tetracosanol-1 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.495	9.12 ng	505239	Perylene-d12	15.865

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2		1-Heptacosanol	396	C27H56O	002004-39-9	94
3		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
4		1-Tetracosene	336	C24H48	010192-32-2	91
5		1-Heneicosanol	312	C21H44O	015594-90-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
Data File : BF132307.D
Acq On : 02 Feb 2023 23:24
Operator : CG\JU
Sample : 01371-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1-BERKSHIRE-NW

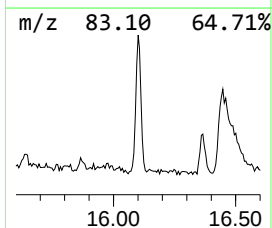
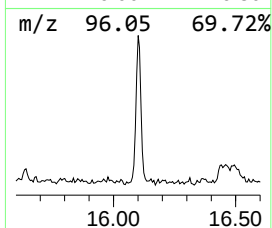
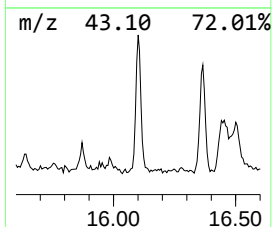
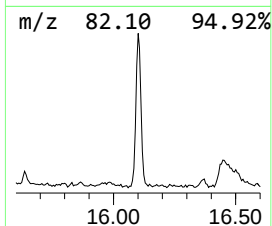
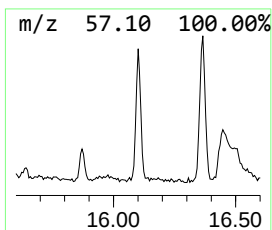
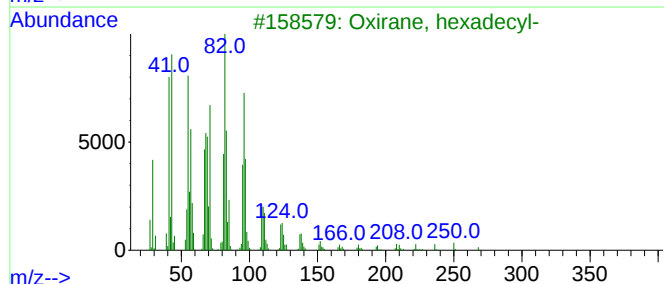
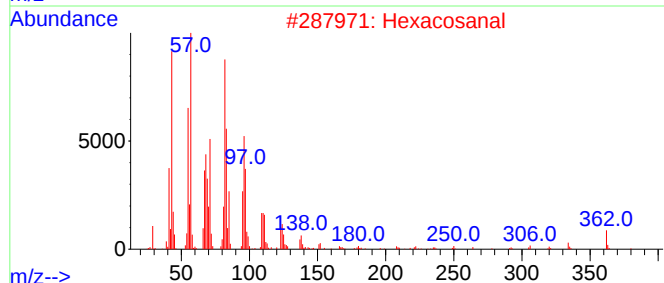
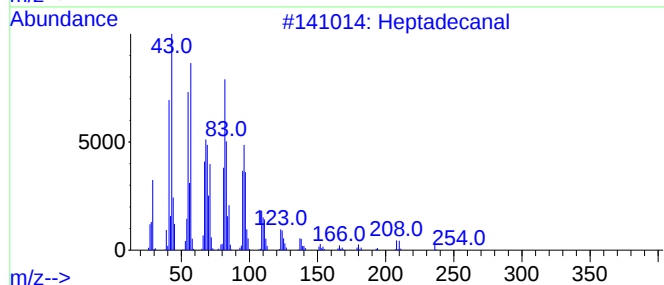
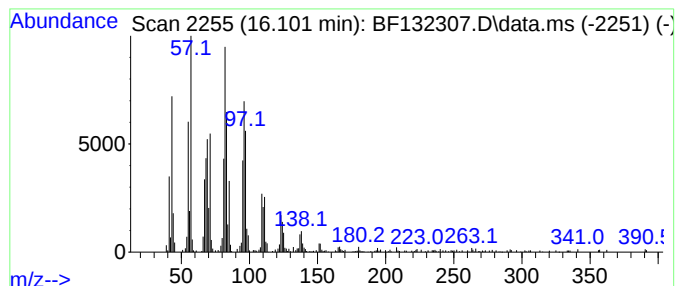
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecanal Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.101	7.03 ng	389451	Perylene-d12	15.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecanal	254	C17H34O	1000376-70-0	94
2			Hexacosanal	380	C26H52O	026627-85-0	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4			Octadecanal	268	C18H36O	000638-66-4	91
5			Octacosanal	408	C28H56O	022725-64-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
Data File : BF132307.D
Acq On : 02 Feb 2023 23:24
Operator : CG\JU
Sample : 01371-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
1-BERKSHIRE-NW

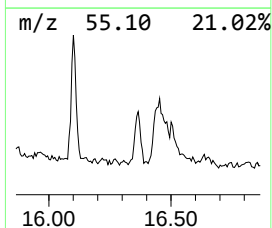
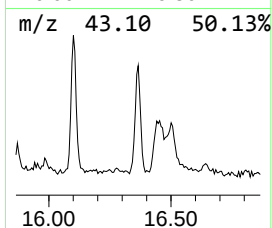
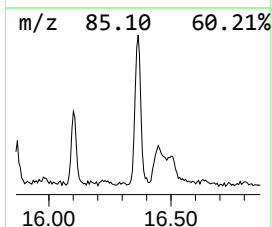
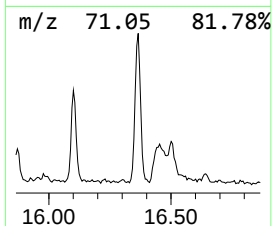
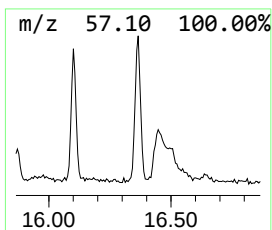
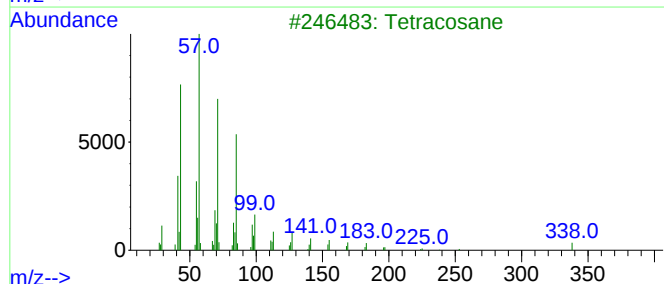
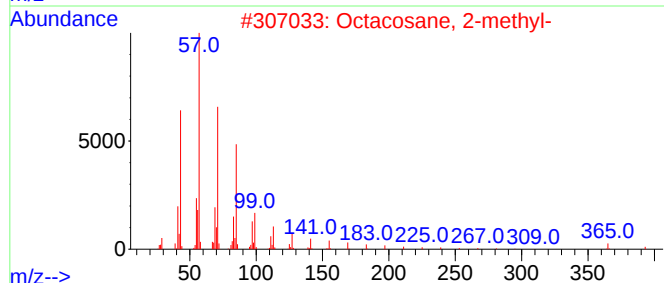
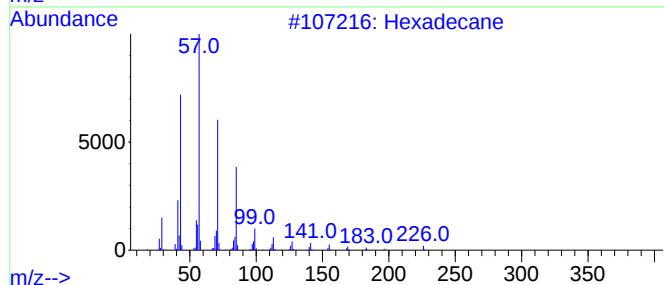
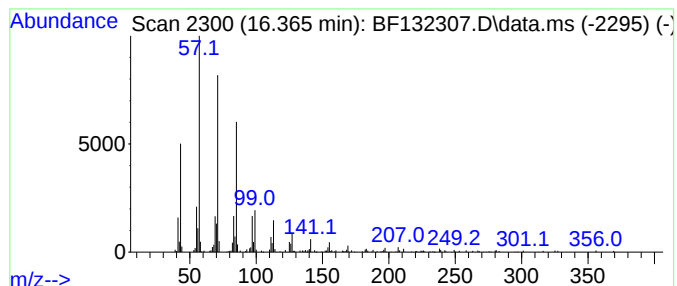
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.365	3.94 ng	218035	Perylene-d12	15.865

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	93
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
Data File : BF132307.D
Acq On : 02 Feb 2023 23:24
Operator : CG\JU
Sample : 01371-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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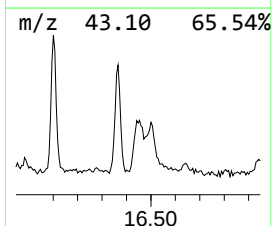
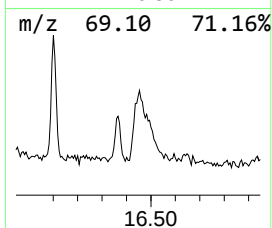
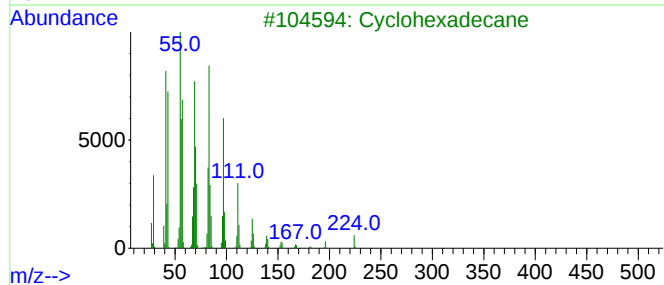
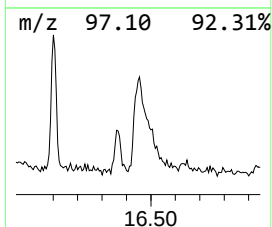
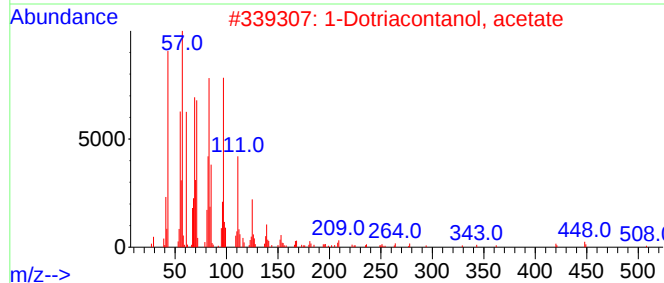
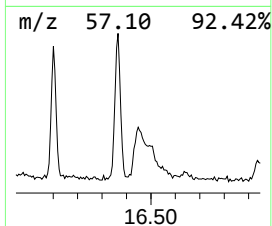
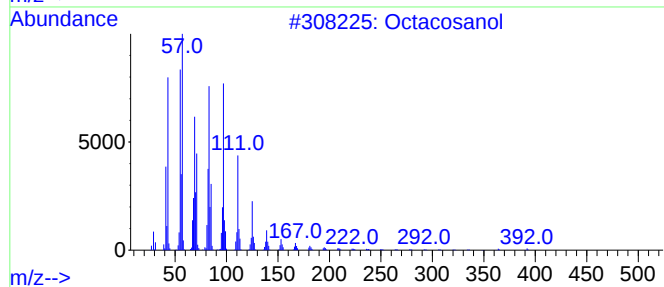
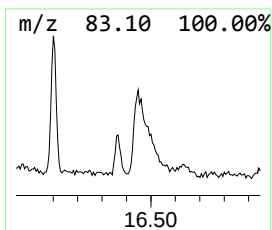
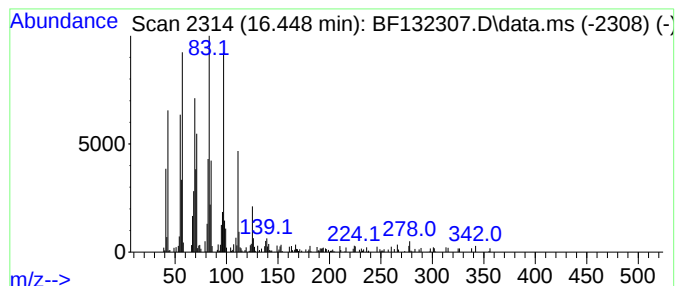
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octacosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.448	4.86 ng	269151	Perylene-d12	15.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosanol	410	C28H58O	000557-61-9	86
2			1-Dotriacontanol, acetate	509	C34H68O2	023811-94-1	80
3			Cyclohexadecane	224	C16H32	000295-65-8	74
4			1-Heneicosanol	312	C21H44O	015594-90-8	72
5			2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020223\
Data File : BF132307.D
Acq On : 02 Feb 2023 23:24
Operator : CG\JU
Sample : 01371-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.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
2-Pentanone, 4-...	5.313	30.1	ng	1629330	1	7.066	1081590	20.0
Longifolene	9.748	2.4	ng	189844	3	10.113	1607330	20.0
Benzophenone	10.830	3.4	ng	271840	3	10.113	1607330	20.0
n-Hexadecanoic ...	12.124	10.0	ng	718173	4	11.613	1439870	20.0
Octadecanoic acid	12.913	2.9	ng	211397	4	11.613	1439870	20.0
2-Phenanthrenol...	13.642	4.6	ng	211159	5	14.277	916415	20.0
5-Octadecene, (E)-	14.101	13.5	ng	616822	5	14.277	916415	20.0
9-Nonadecene	14.759	2.3	ng	106865	5	14.277	916415	20.0
Supraene	15.154	5.3	ng	294129	6	15.865	1107640	20.0
Henicosanal	15.230	3.6	ng	201987	6	15.865	1107640	20.0
Octacosane, 2-m...	15.448	2.5	ng	140514	6	15.865	1107640	20.0
n-Tetracosanol-1	15.495	9.1	ng	505239	6	15.865	1107640	20.0
Heptadecanal	16.101	7.0	ng	389451	6	15.865	1107640	20.0
Hexadecane	16.365	3.9	ng	218035	6	15.865	1107640	20.0
Octacosanol	16.448	4.9	ng	269151	6	15.865	1107640	20.0
5-(3-Hydroxy-4,...	16.983	2.3	ng	127514	6	15.865	1107640	20.0