

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
 Data File : BF123717.D
 Acq On : 24 Apr 2021 15:14
 Operator : JU/CG
 Sample : M2149-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-40-9.5-10.0-DUP

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-BF041521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF123717.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.034	495	501	507	rVB	120892	162376	2.26%	0.291%
2	5.439	566	570	574	rBV	7105476	6670690	92.76%	11.949%
3	6.457	738	743	755	rBV	6916017	7191052	100.00%	12.882%
4	6.822	801	805	808	rBV	1583759	1345921	18.72%	2.411%
5	7.386	895	901	904	rBV	4722181	4692866	65.26%	8.406%
6	8.104	1019	1023	1026	rBV	2305135	1750618	24.34%	3.136%
7	9.180	1202	1206	1209	rBV	7749057	6706809	93.27%	12.014%
8	9.569	1269	1272	1276	rBV	208860	190208	2.65%	0.341%
9	9.857	1317	1321	1325	rBV	2253061	1869447	26.00%	3.349%
10	10.651	1451	1456	1459	rBV	5394810	5022226	69.84%	8.996%
11	11.345	1570	1574	1577	rBV	2415292	1970257	27.40%	3.529%
12	11.410	1580	1585	1594	rVB5	52811	78235	1.09%	0.140%
13	11.880	1661	1665	1672	rBV	304599	305920	4.25%	0.548%
14	12.939	1840	1845	1848	rBV	7414901	6741448	93.75%	12.076%
15	13.168	1880	1884	1901	rVB2	175772	404520	5.63%	0.725%
16	13.627	1959	1962	1966	rVB	259106	207300	2.88%	0.371%
17	13.839	1994	1998	2016	rBV	990111	2276378	31.66%	4.078%
18	13.986	2019	2023	2027	rVB	1965716	1749904	24.33%	3.135%
19	14.280	2070	2073	2078	rVB	611924	524636	7.30%	0.940%
20	14.486	2101	2108	2113	rBV	616516	1383966	19.25%	2.479%
21	14.586	2123	2125	2135	rVB2	82750	84499	1.18%	0.151%
22	14.915	2177	2181	2188	rVB	548454	577955	8.04%	1.035%
23	15.104	2209	2213	2218	rBV	233740	273346	3.80%	0.490%
24	15.162	2218	2223	2224	rBV4	98200	147589	2.05%	0.264%
25	15.451	2267	2272	2275	rBV	1075666	1361343	18.93%	2.439%
26	15.480	2275	2277	2283	rVB2	123532	174193	2.42%	0.312%
27	15.686	2308	2312	2317	rBV2	97050	127133	1.77%	0.228%
28	15.815	2330	2334	2339	rBV4	46790	85694	1.19%	0.154%
29	15.921	2347	2352	2357	rVB	167478	251770	3.50%	0.451%
30	16.009	2362	2367	2368	rBV4	66337	94279	1.31%	0.169%
31	16.186	2393	2397	2404	rVB5	52955	84769	1.18%	0.152%
32	16.421	2433	2437	2443	rVB3	80820	131331	1.83%	0.235%
33	16.950	2522	2527	2532	rBV6	75082	142695	1.98%	0.256%
34	17.009	2534	2537	2544	rVB3	65006	118963	1.65%	0.213%
35	17.221	2570	2573	2583	rVB6	113328	257368	3.58%	0.461%
36	17.562	2618	2631	2639	rBV10	206126	666910	9.27%	1.195%

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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-BF041521.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

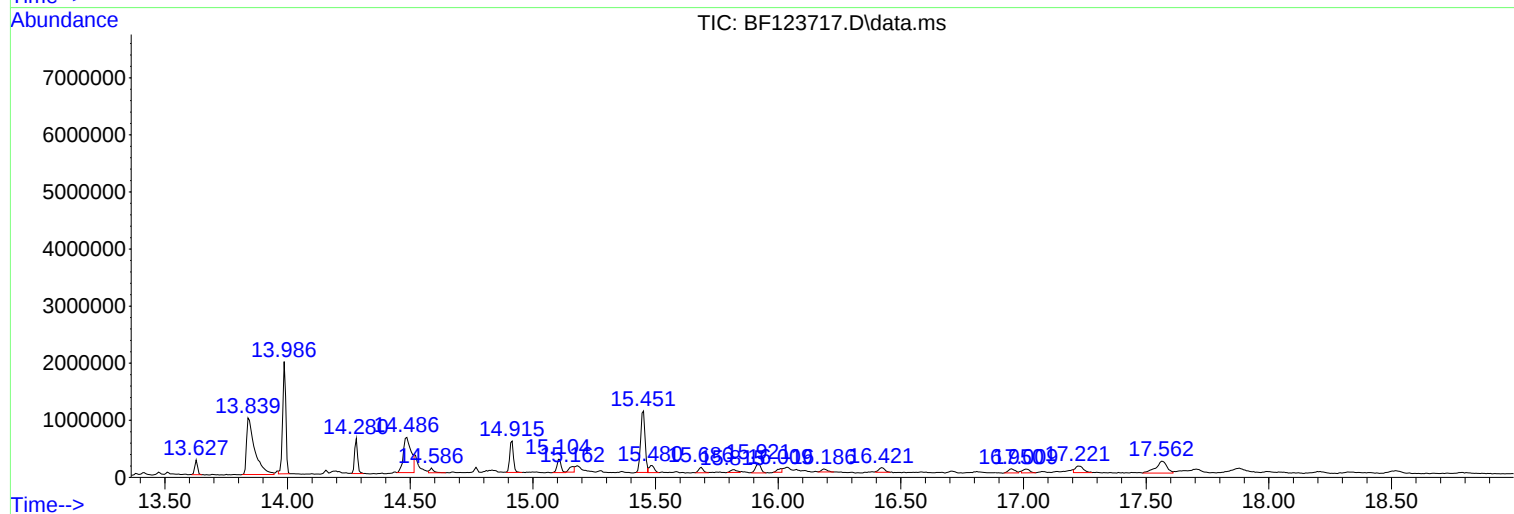
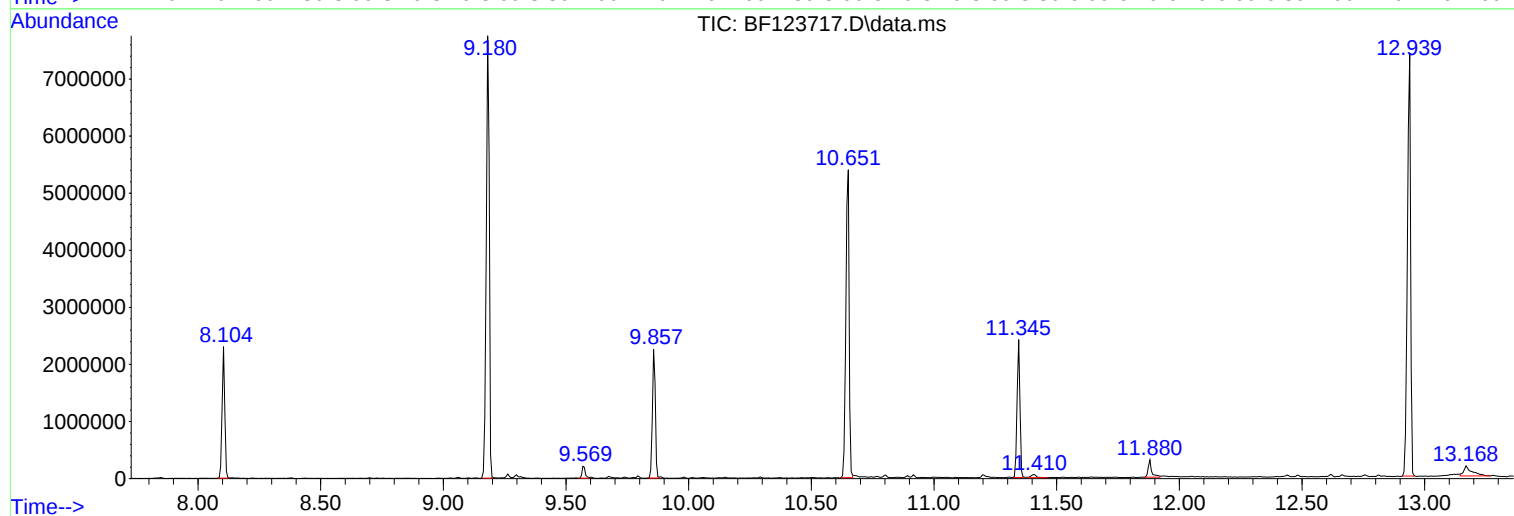
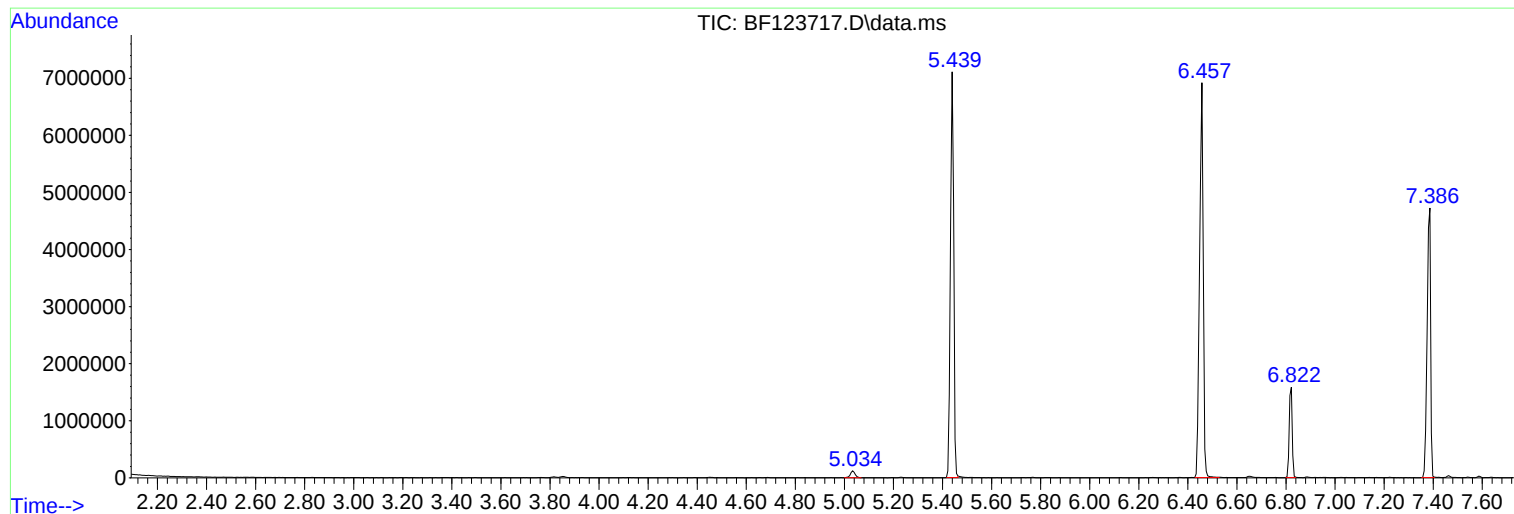
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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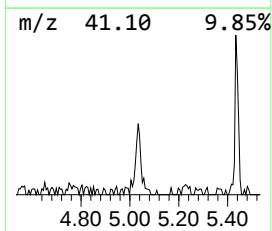
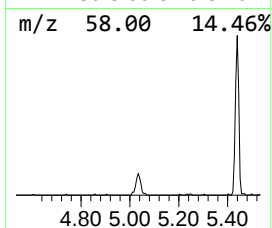
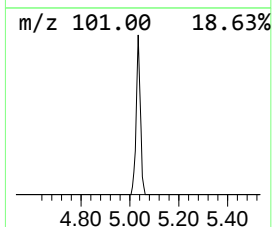
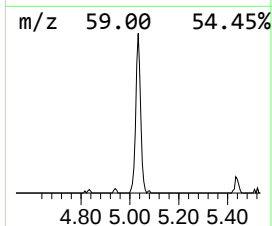
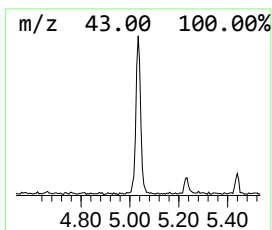
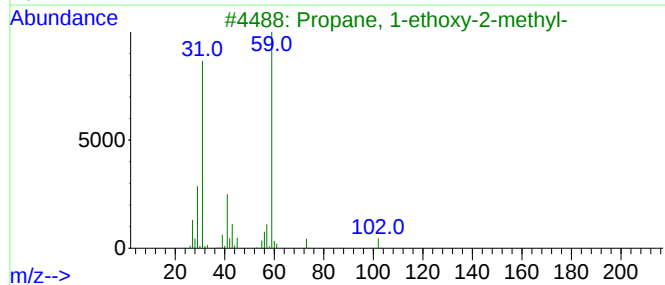
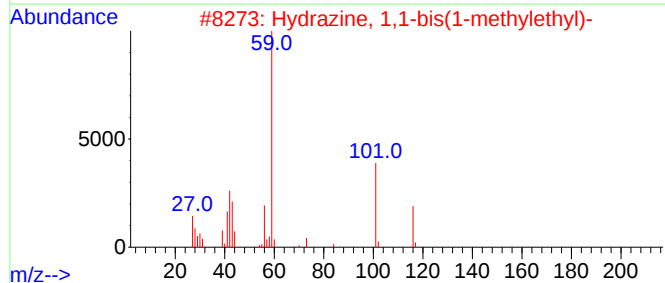
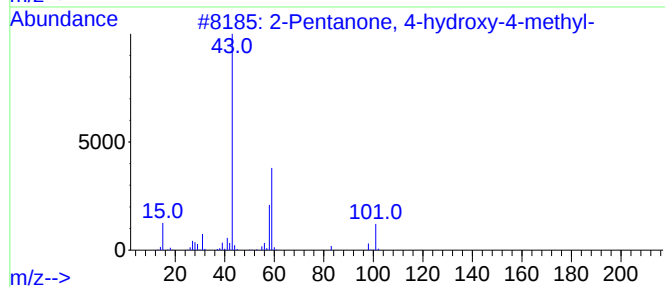
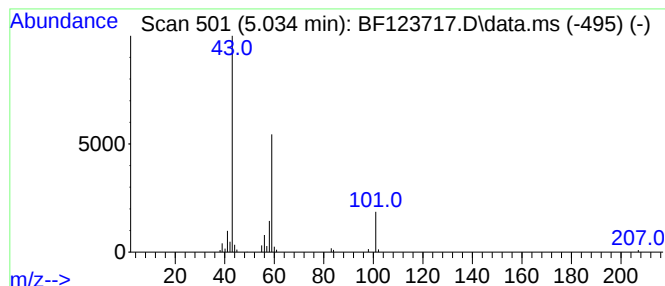
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.034	2.41 ng	162376	1,4-Dichlorobenzene-d4	6.822

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	40		
3	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38		
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27		



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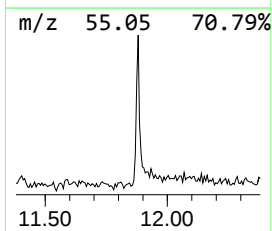
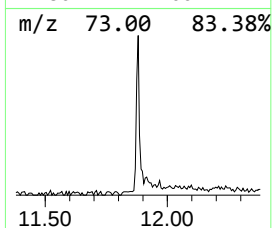
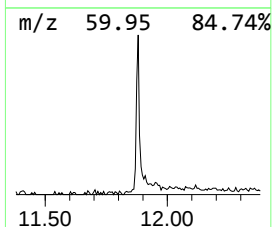
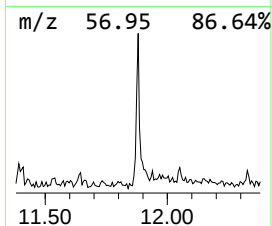
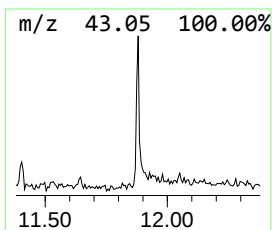
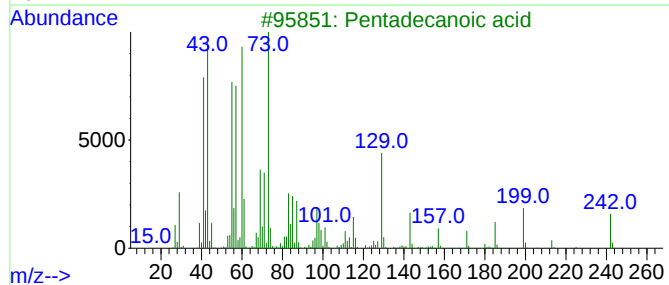
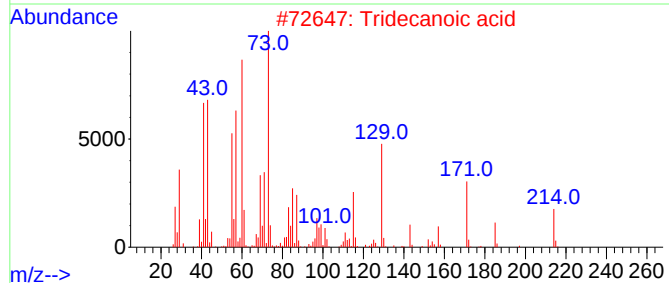
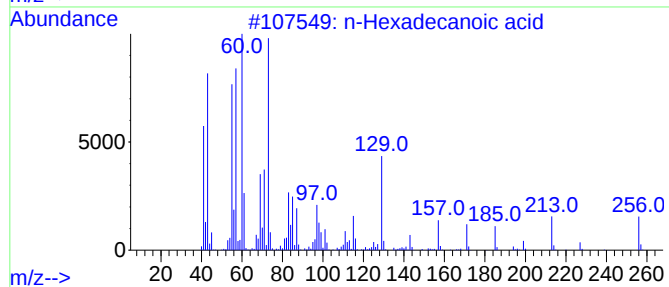
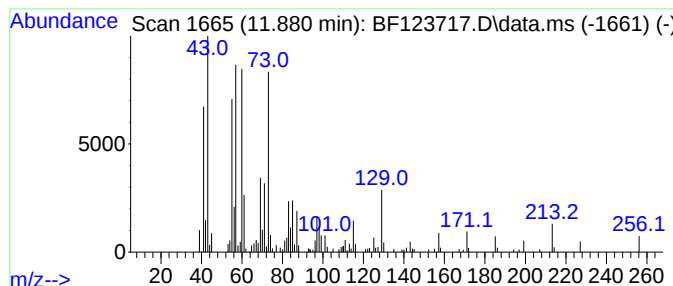
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	3.11 ng	305920	Phenanthrene-d10	11.345

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	97
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4			Undecanoic acid	186	C11H22O2	000112-37-8	87
5			n-Decanoic acid	172	C10H20O2	000334-48-5	86



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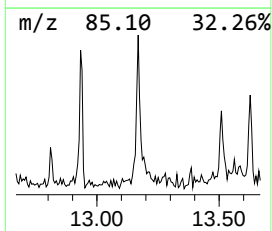
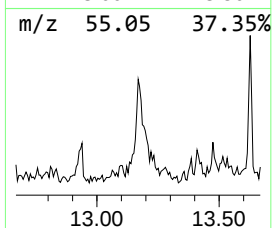
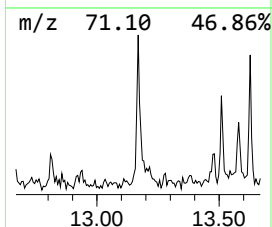
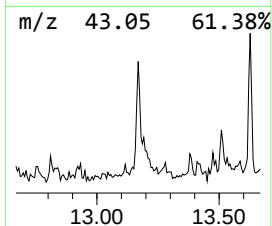
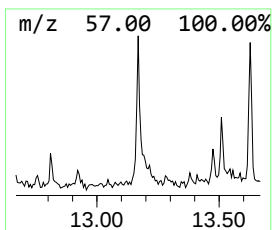
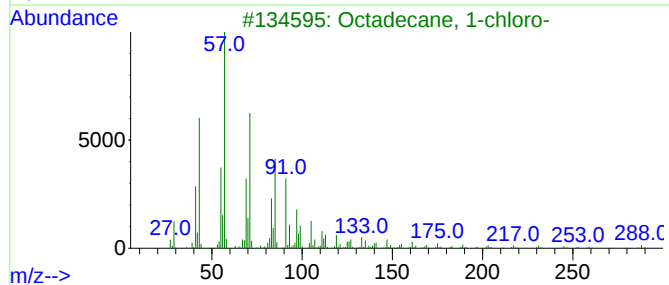
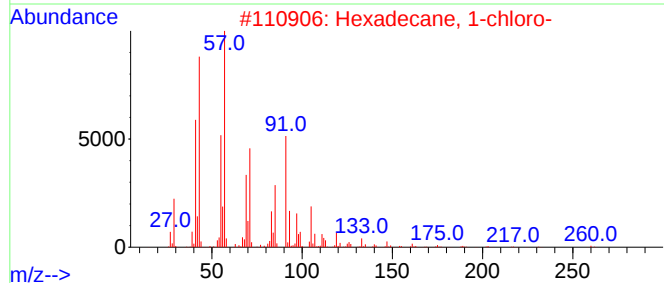
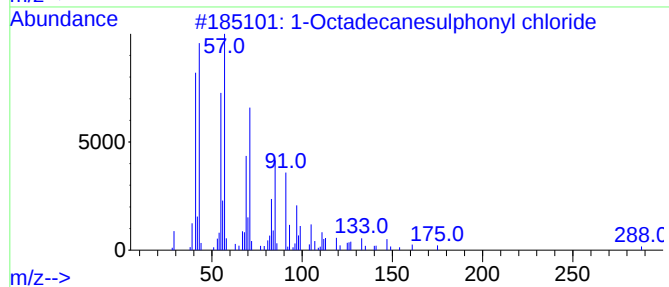
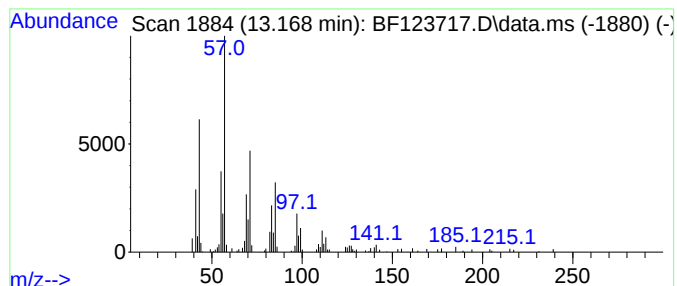
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Octadecanesulphonyl chloride Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.168	4.62 ng	404520	Chrysene-d12	13.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	90
2			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	86
3			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	86
4			Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	86
5			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	86



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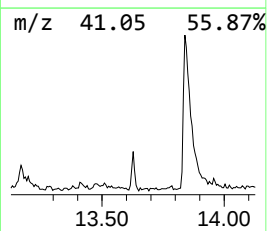
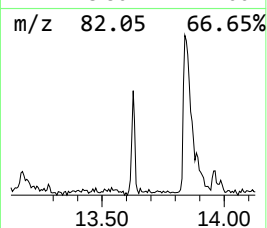
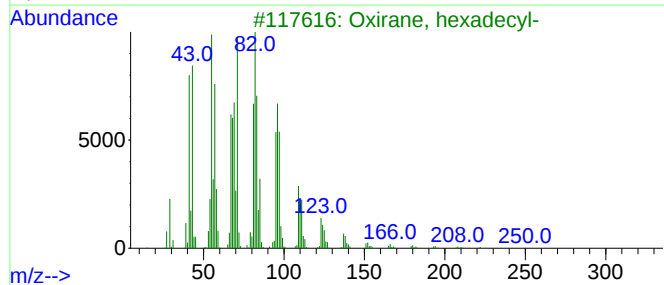
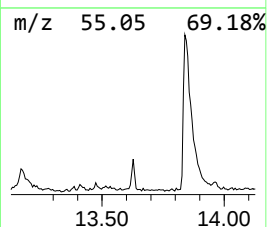
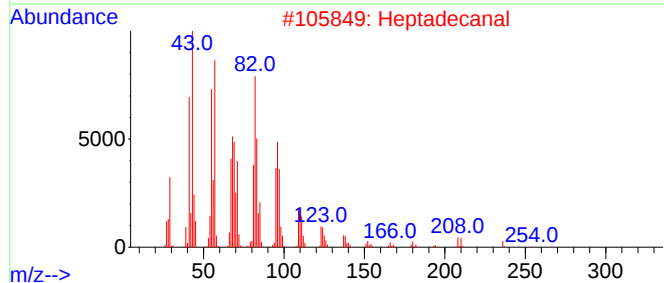
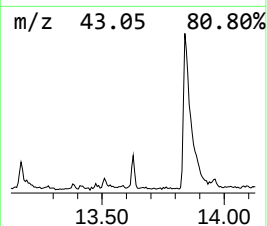
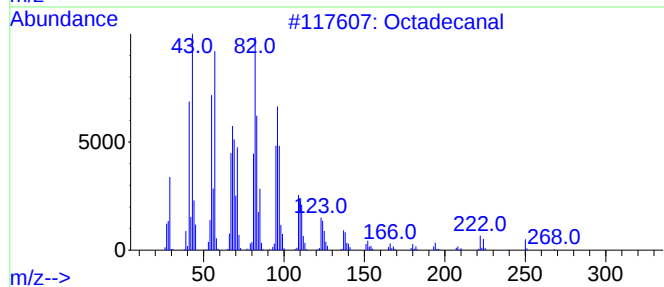
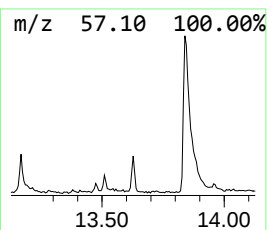
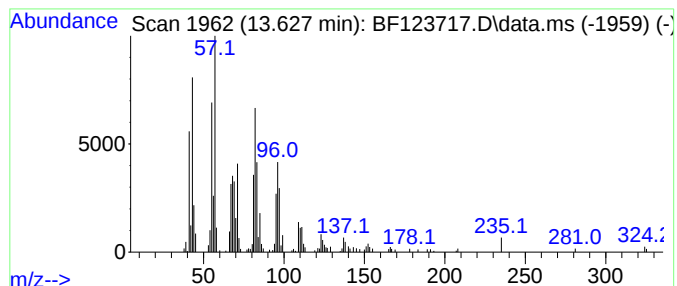
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Octadecanal Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.627	2.37 ng	207300	Chrysene-d12	13.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanal	268	C18H36O	000638-66-4	91
2			Heptadecanal	254	C17H34O	1000376-70-0	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
4			Hexadecanal	240	C16H32O	000629-80-1	87
5			1,16-Hexadecanediol	258	C16H34O2	007735-42-4	81



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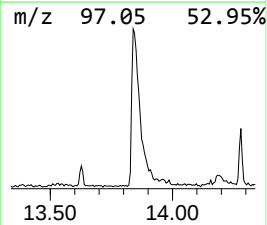
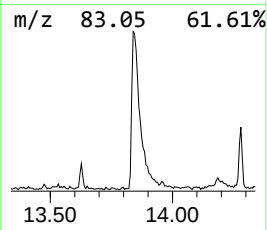
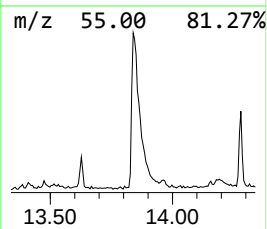
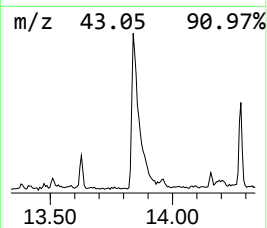
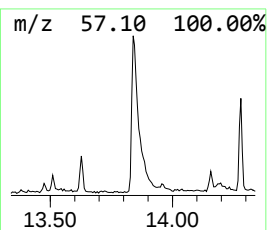
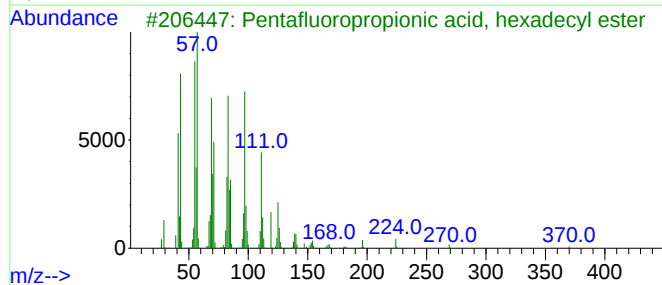
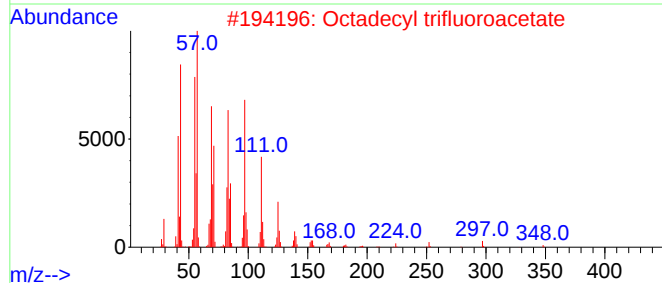
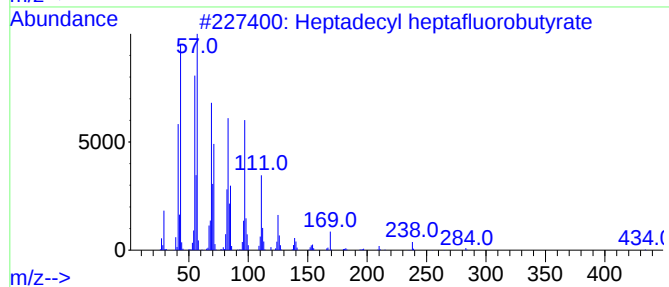
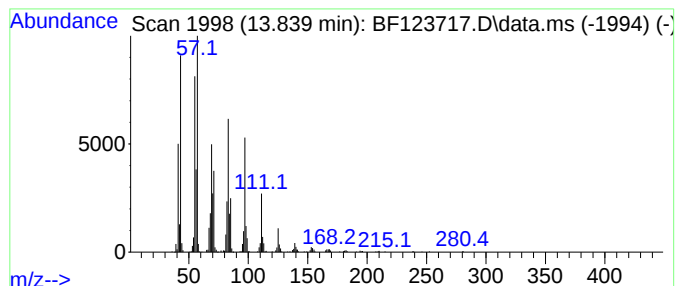
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Heptadecyl heptafluorobutyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.839	26.02 ng	2276380	Chrysene-d12	13.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
3			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
4			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
Acq On : 24 Apr 2021 15:14
Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-40-9.5-10.0-DUP

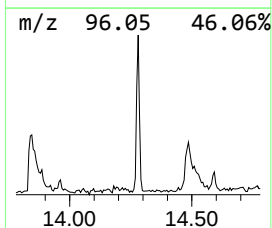
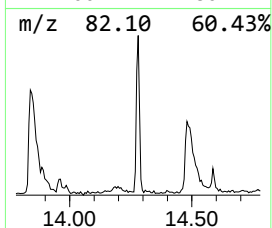
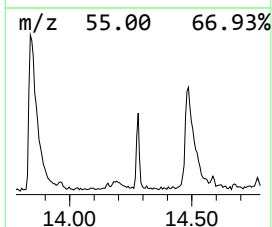
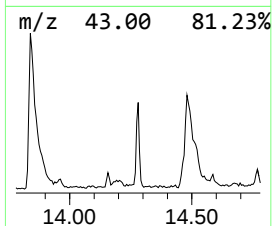
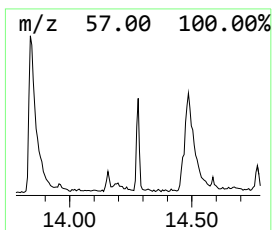
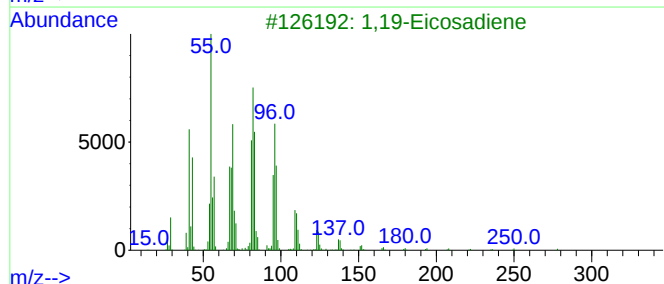
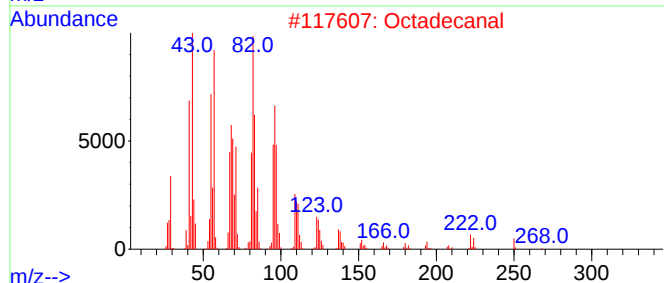
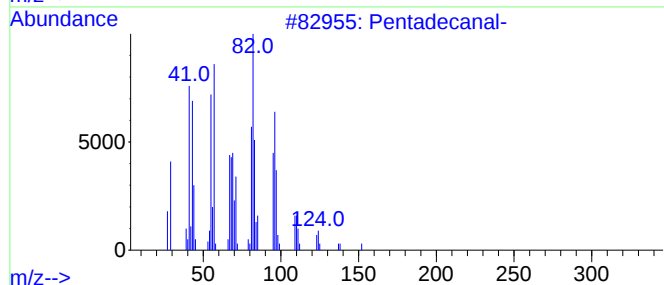
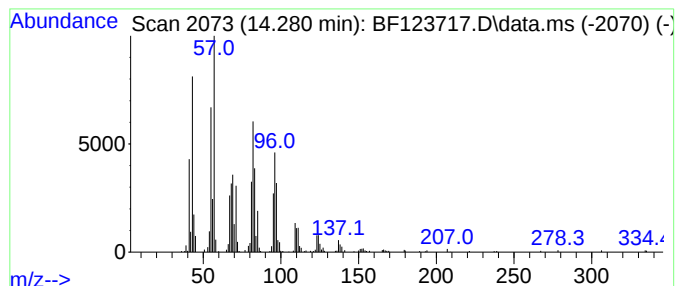
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Pentadecanal- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.280	6.00 ng	524636	Chrysene-d12	13.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanal-	226	C15H30O	002765-11-9	91
2			Octadecanal	268	C18H36O	000638-66-4	91
3			1,19-Eicosadiene	278	C20H38	014811-95-1	90
4			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
5			16-Heptadecenal	252	C17H32O	1000144-57-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
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Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

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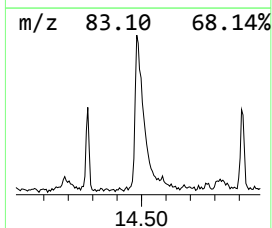
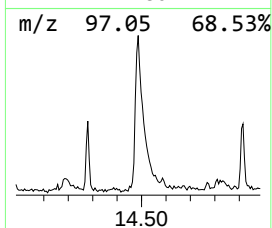
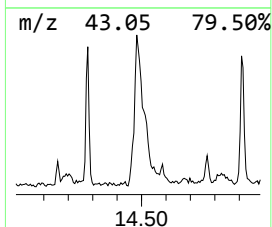
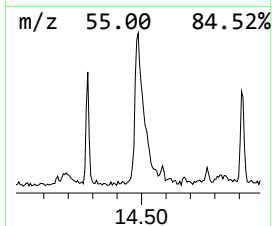
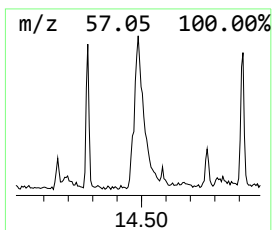
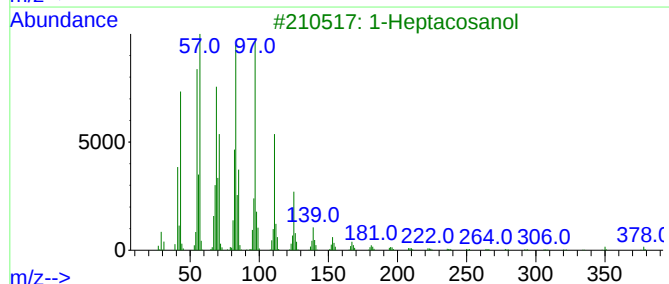
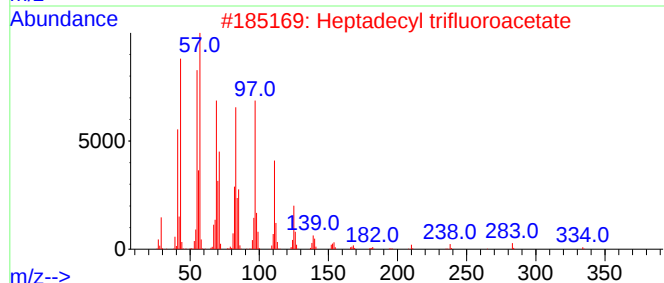
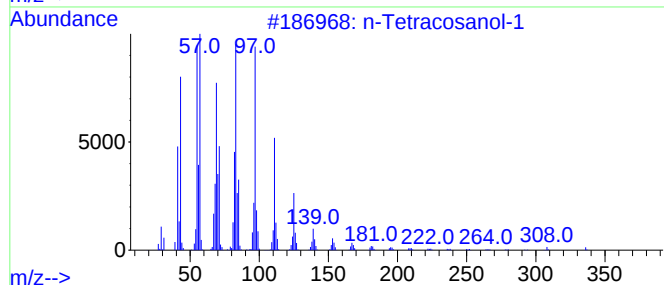
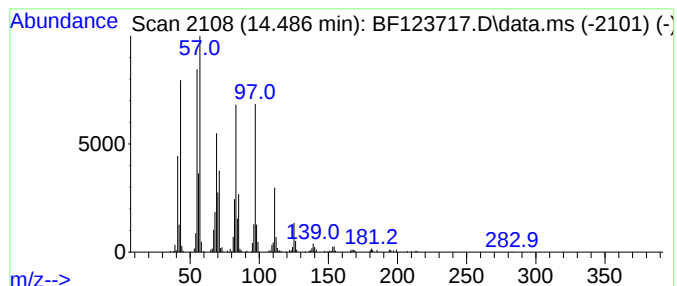
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 n-Tetracosanol-1 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.486	15.82 ng	1383970	Chrysene-d12	13.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
2			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
3			1-Heptacosanol	396	C27H56O	002004-39-9	91
4			Octacosanol	410	C28H58O	000557-61-9	91
5			1-Heneicosanol	312	C21H44O	015594-90-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
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Operator : JU/CG
Sample : M2149-16
Misc :
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Instrument :
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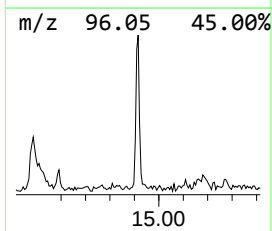
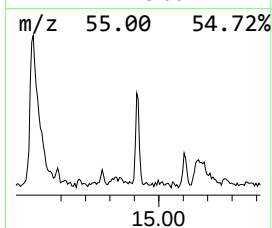
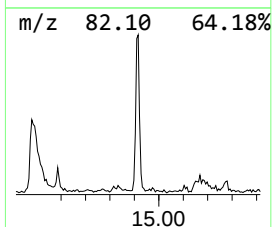
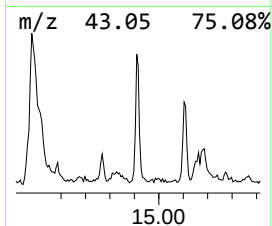
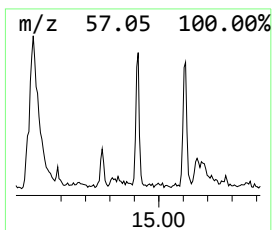
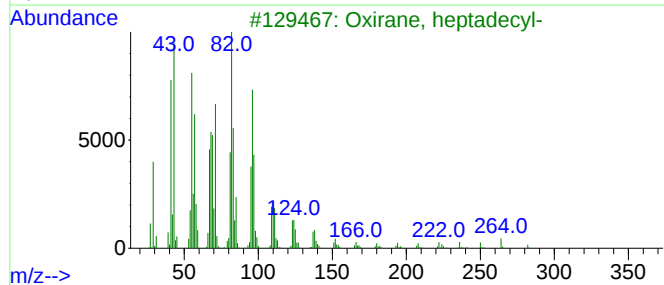
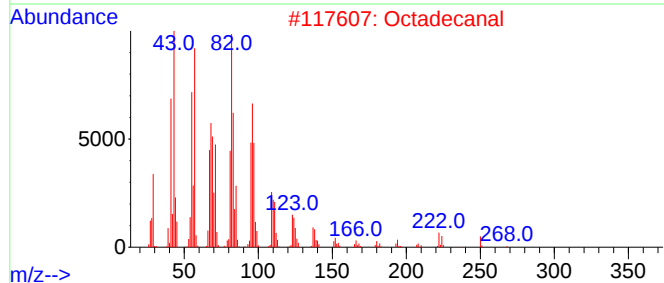
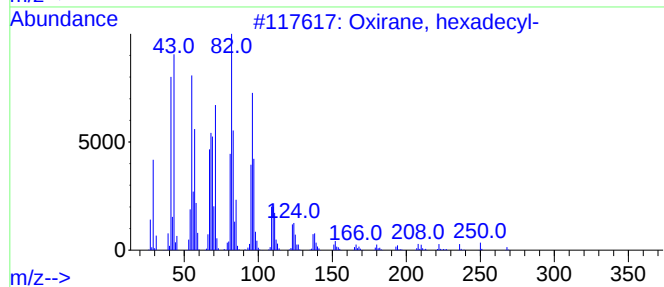
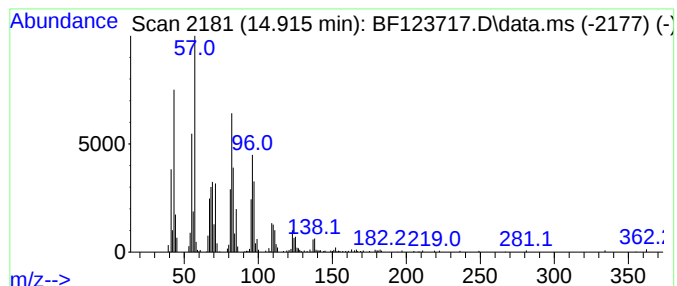
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Oxirane, hexadecyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.915	8.49 ng	577955	Perylene-d12	15.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2			Octadecanal	268	C18H36O	000638-66-4	91
3			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4			1,19-Eicosadiene	278	C20H38	014811-95-1	90
5			(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
Acq On : 24 Apr 2021 15:14
Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

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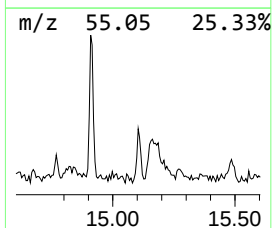
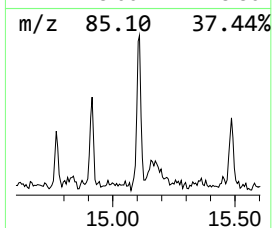
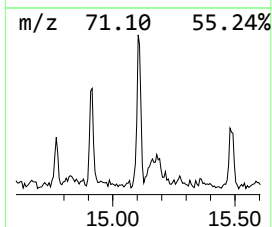
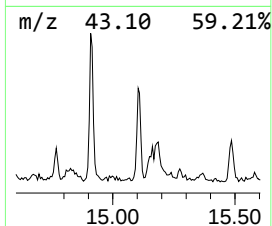
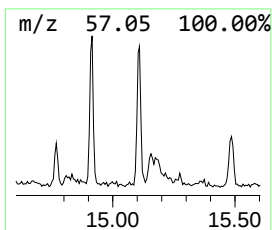
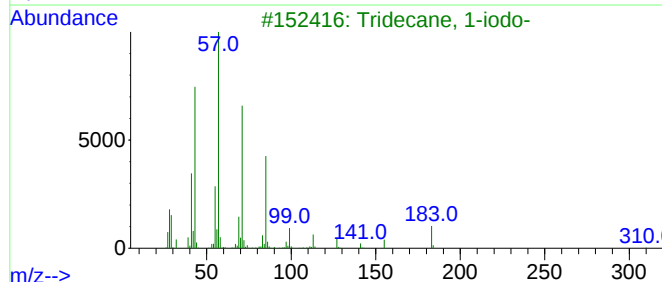
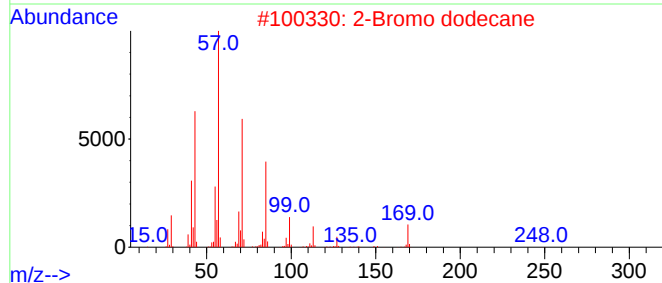
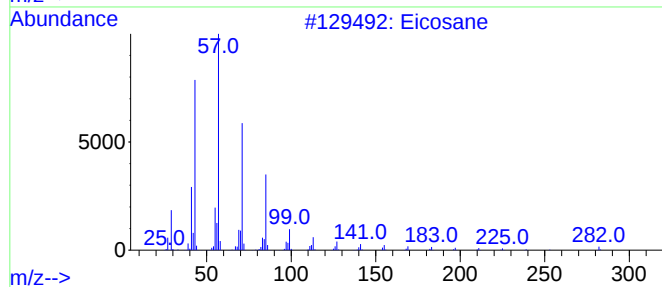
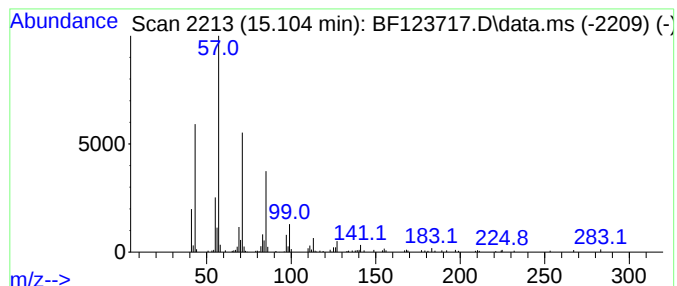
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.104	4.02 ng	273346	Perylene-d12	15.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	90
3			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5			Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
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Acq On : 24 Apr 2021 15:14
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Sample : M2149-16
Misc :
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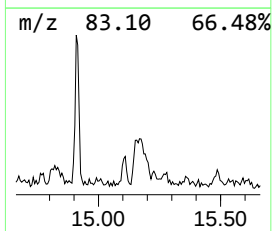
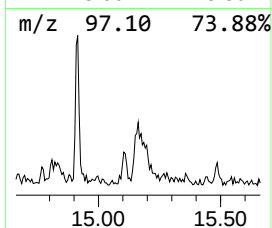
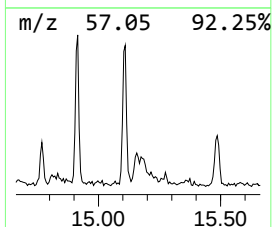
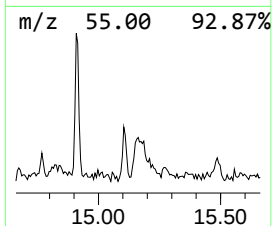
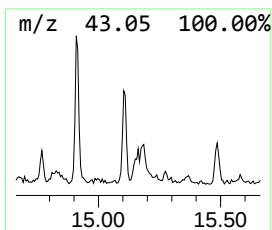
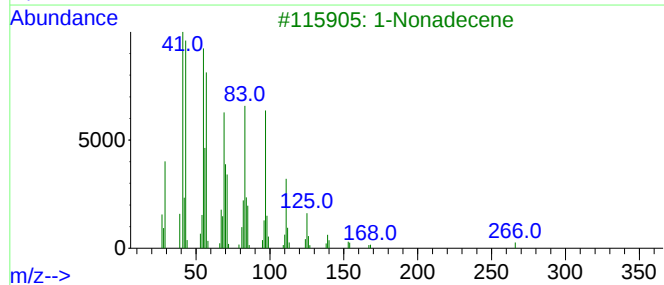
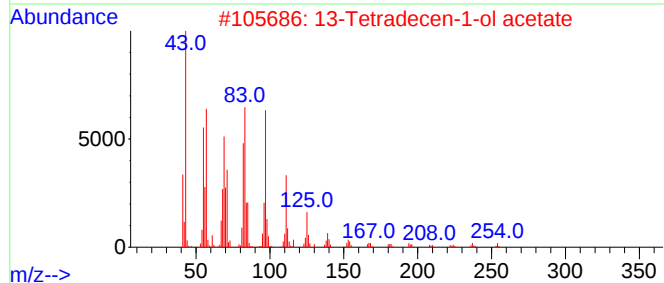
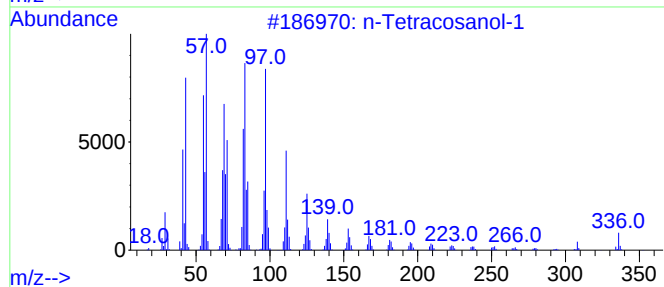
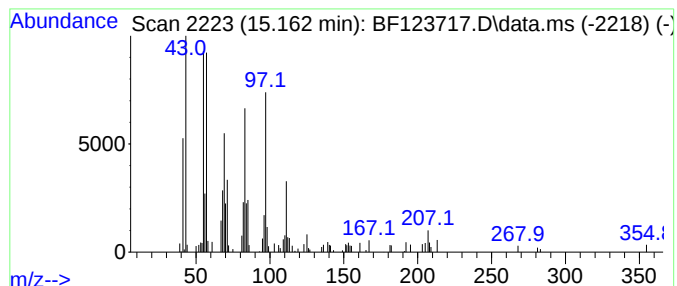
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 13-Tetradecen-1-ol acetate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.162	2.17 ng	147589	Perylene-d12	15.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	80
2	13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	80
3	1-Nonadecene	266	C19H38	018435-45-5	72
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	53
5	Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
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Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

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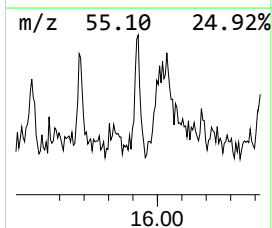
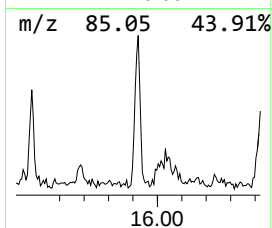
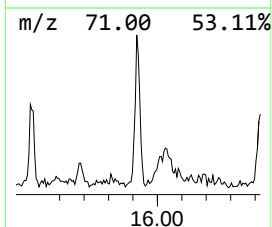
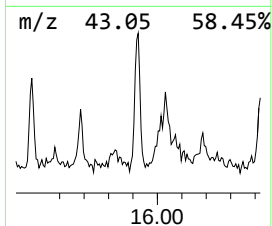
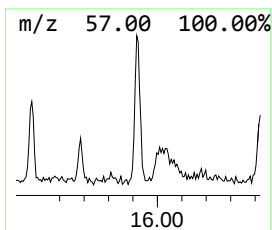
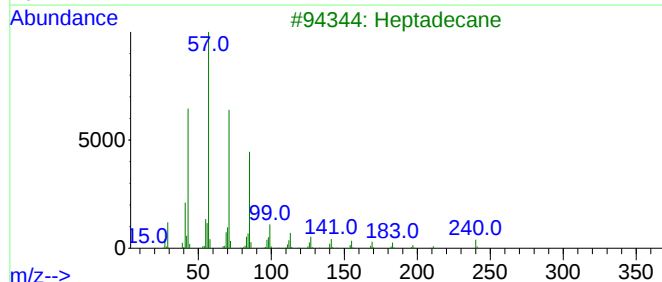
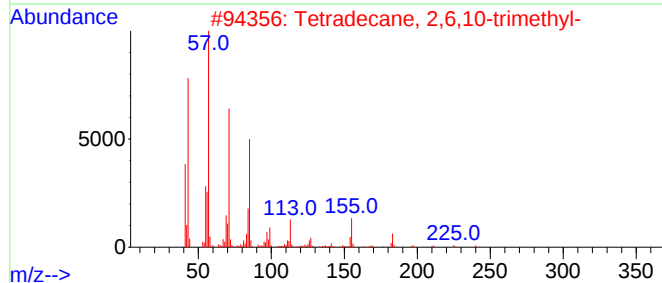
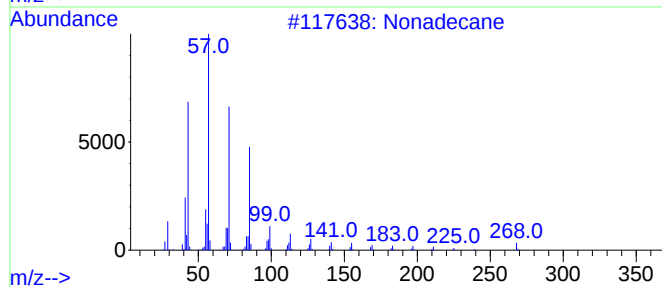
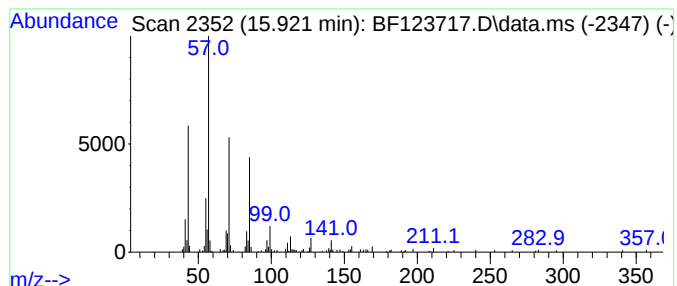
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.921	3.70 ng	251770	Perylene-d12	15.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	90
3			Heptadecane	240	C17H36	000629-78-7	90
4			Hexadecane	226	C16H34	000544-76-3	90
5			Tetradecane	198	C14H30	000629-59-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
Acq On : 24 Apr 2021 15:14
Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

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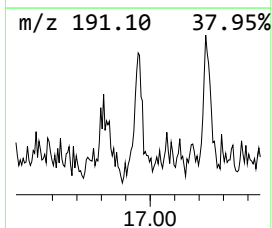
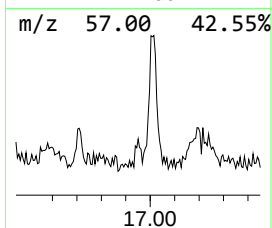
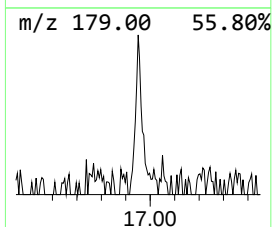
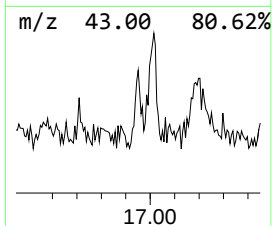
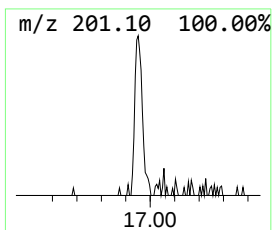
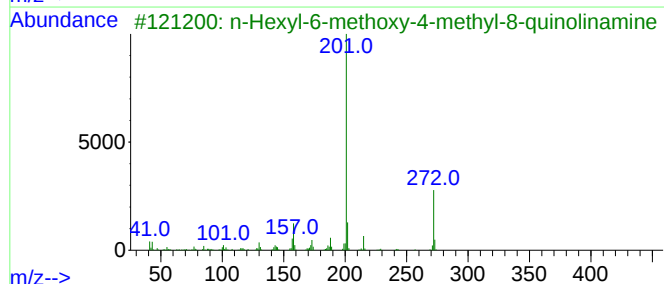
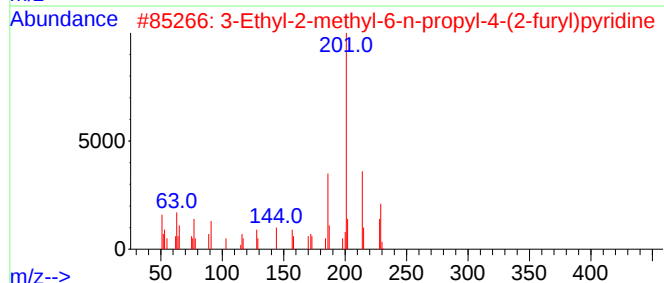
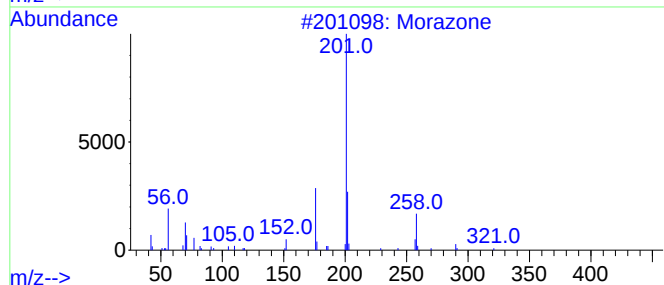
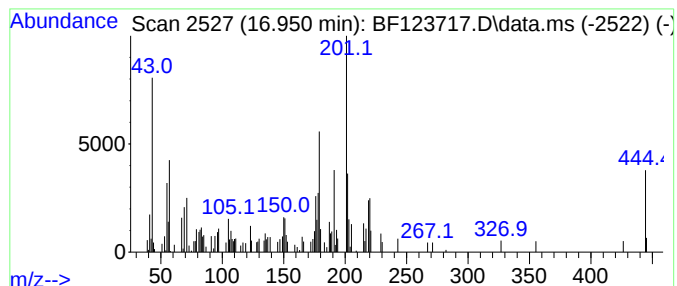
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown16.951 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.951	2.10 ng	142695	Perylene-d12	15.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morazone	377	C23H27N3O2	006536-18-1	27
2			3-Ethyl-2-methyl-6-n-propyl-4-(2-furyl)pyridine	229	C15H19NO	078563-70-9	25
3			n-Hexyl-6-methoxy-4-methyl-8-quinolinamine	272	C17H24N2O	083547-16-4	22
4			1H-Pyrrole-2,5-dione, 2,5-dihydro-1,3-dimethyl-4-(2-furyl)-	201	C12H11N2O2	065833-09-2	22
5			Phosphinous acid, diphenyl-	202	C12H11OP	1000296-41-9	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
Acq On : 24 Apr 2021 15:14
Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-40-9.5-10.0-DUP

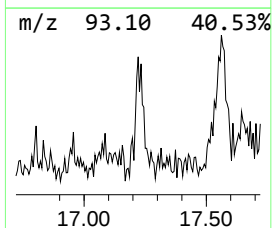
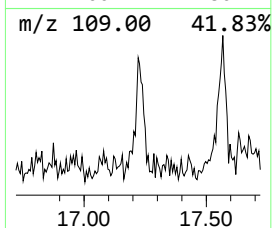
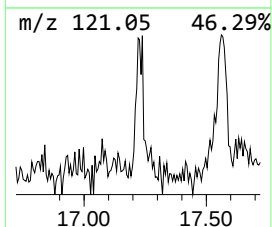
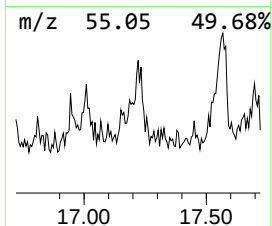
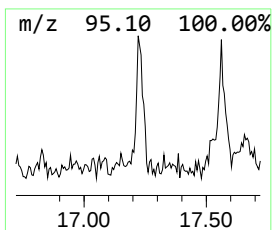
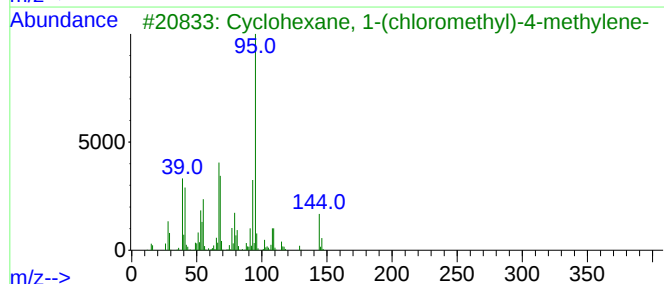
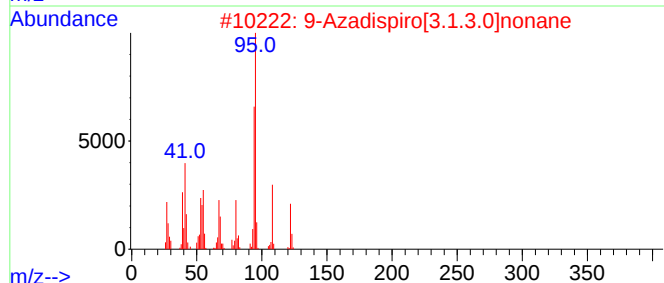
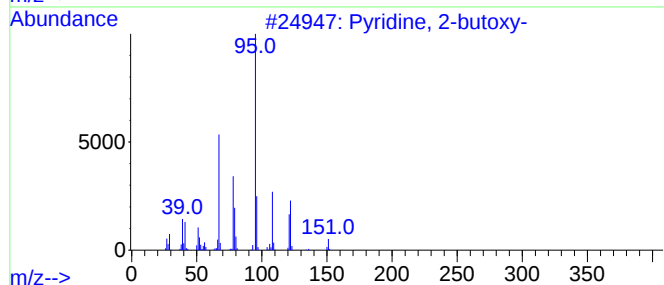
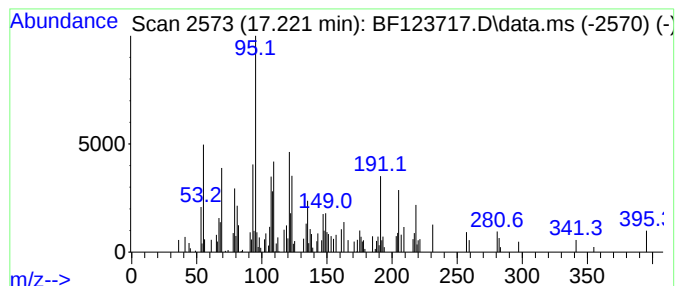
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown17.221 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.221	3.78 ng	257368	Perylene-d12	15.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine, 2-butoxy-	151	C9H13NO	027361-16-6	27
2			9-Azadispiro[3.1.3.0]nonane	123	C8H13N	135763-48-3	27
3			Cyclohexane, 1-(chloromethyl)-4-...	144	C8H13Cl	000823-83-6	22
4			Cyclohexanemethanol, 4-(1-methyl...	156	C10H20O	013828-37-0	22
5			Cyclopentene, 1,4-dimethyl-5-(1-...	138	C10H18	061142-33-4	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
Acq On : 24 Apr 2021 15:14
Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-40-9.5-10.0-DUP

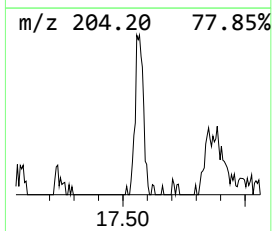
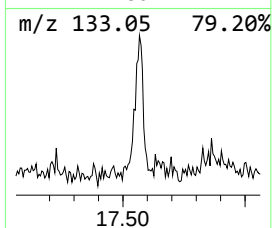
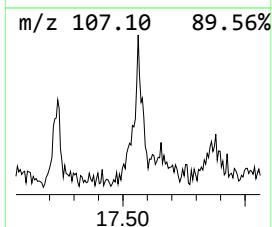
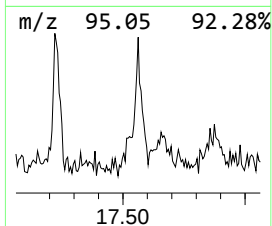
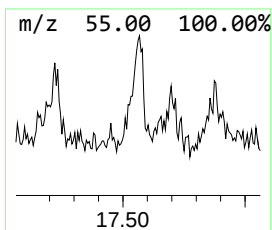
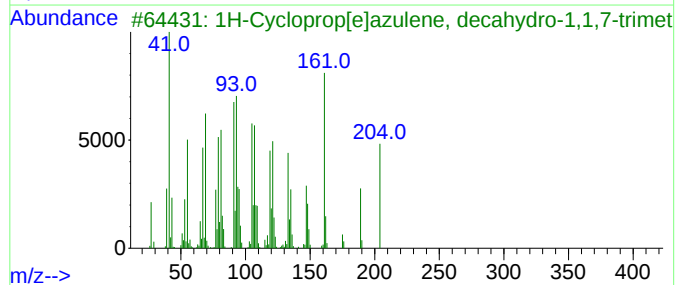
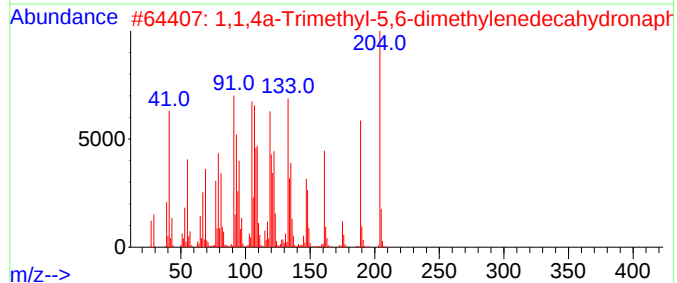
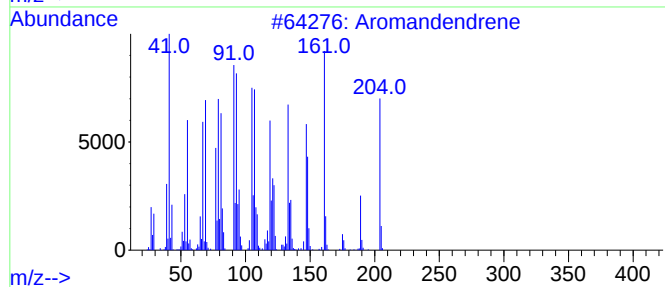
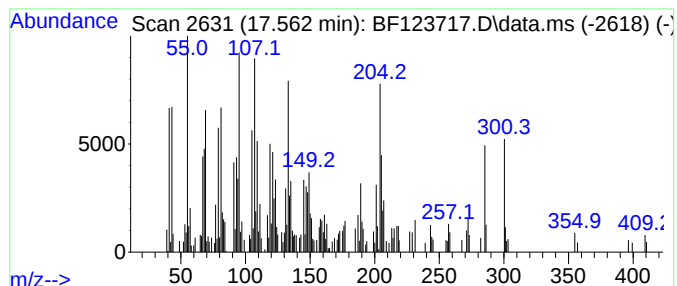
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Aromandendrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.562	9.80 ng	666910	Perylene-d12	15.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Aromandendrene	204	C15H24	000489-39-4	55
2			1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	49
3			1H-Cycloprop[e]azulene, decahydr...	204	C15H24	072747-25-2	43
4			Alloaromadendrene	204	C15H24	025246-27-9	38
5			.alpha.-Guaiene	204	C15H24	003691-12-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042421\
Data File : BF123717.D
Acq On : 24 Apr 2021 15:14
Operator : JU/CG
Sample : M2149-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-40-9.5-10.0-DUP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.034	2.4	ng	162376	1	6.822	1345920	20.0
n-Hexadecanoic ...	11.880	3.1	ng	305920	4	11.345	1970260	20.0
1-Octadecanesul...	13.168	4.6	ng	404520	5	13.986	1749900	20.0
Octadecanal	13.627	2.4	ng	207300	5	13.986	1749900	20.0
Heptadecyl hept...	13.839	26.0	ng	2276380	5	13.986	1749900	20.0
Pentadecanal-	14.280	6.0	ng	524636	5	13.986	1749900	20.0
n-Tetracosanol-1	14.486	15.8	ng	1383970	5	13.986	1749900	20.0
Oxirane, hexade...	14.915	8.5	ng	577955	6	15.451	1361340	20.0
Eicosane	15.104	4.0	ng	273346	6	15.451	1361340	20.0
13-Tetradecen-1...	15.162	2.2	ng	147589	6	15.451	1361340	20.0
Nonadecane	15.921	3.7	ng	251770	6	15.451	1361340	20.0
unknown16.951	16.951	2.1	ng	142695	6	15.451	1361340	20.0
unknown17.221	17.221	3.8	ng	257368	6	15.451	1361340	20.0
Aromandendrene	17.562	9.8	ng	666910	6	15.451	1361340	20.0