

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
 Data File : BF125156.D
 Acq On : 23 Aug 2021 15:32
 Operator : JU/CG
 Sample : M3465-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CON-C

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-BF081821.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125156.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.216	15	22	28	rVB	1002106	1348049	34.65%	4.103%
2	3.681	265	271	277	rBV	53415	78736	2.02%	0.240%
3	4.540	408	417	420	rBV	2476481	2950433	75.83%	8.980%
4	5.145	514	520	523	rBV	2107834	2453456	63.06%	7.467%
5	5.534	582	586	590	rBV	2060920	1865449	47.94%	5.678%
6	6.204	697	700	704	rVB	894436	689814	17.73%	2.100%
7	6.539	752	757	760	rBV	2957667	2815315	72.36%	8.569%
8	6.922	818	822	825	rBV	1202643	946003	24.31%	2.879%
9	7.239	872	876	879	rBV	94525	72444	1.86%	0.220%
10	7.481	912	917	920	rBV	2440681	2258081	58.04%	6.873%
11	8.098	1019	1022	1036	rBV	309404	286843	7.37%	0.873%
12	8.204	1036	1040	1043	rBV	1565808	1229617	31.60%	3.742%
13	9.280	1218	1223	1226	rBV	4658211	3890892	100.00%	11.842%
14	9.963	1335	1339	1342	rBV	1496138	1381025	35.49%	4.203%
15	10.386	1408	1411	1413	rVB	85646	65278	1.68%	0.199%
16	10.604	1444	1448	1451	rBV	62331	63905	1.64%	0.195%
17	10.745	1466	1472	1478	rBV	532984	471376	12.11%	1.435%
18	10.857	1488	1491	1496	rBV	336574	348406	8.95%	1.060%
19	11.051	1521	1524	1526	rBV3	44345	45949	1.18%	0.140%
20	11.145	1537	1540	1543	rBV	45823	48762	1.25%	0.148%
21	11.310	1564	1568	1570	rBV	453941	385187	9.90%	1.172%
22	11.339	1570	1573	1575	rVB	115278	105287	2.71%	0.320%
23	11.445	1588	1591	1595	rBV	1391543	1310294	33.68%	3.988%
24	11.733	1637	1640	1644	rBV	437199	325717	8.37%	0.991%
25	11.833	1655	1657	1662	rVB	70827	60983	1.57%	0.186%
26	11.963	1677	1679	1683	rBV2	65851	61645	1.58%	0.188%
27	12.139	1706	1709	1713	rBV	266036	217618	5.59%	0.662%
28	12.533	1773	1776	1781	rBV	130293	138955	3.57%	0.423%
29	12.904	1836	1839	1841	rBV	66342	49160	1.26%	0.150%
30	13.033	1857	1861	1864	rBV	3661887	3379426	86.85%	10.286%
31	13.257	1897	1899	1902	rBV	51468	46099	1.18%	0.140%
32	13.462	1931	1934	1938	rBV	400554	377667	9.71%	1.149%
33	13.510	1938	1942	1950	rVB2	369289	408143	10.49%	1.242%
34	13.927	2010	2013	2024	rVB	222893	437539	11.25%	1.332%
35	14.092	2037	2041	2044	rVB	991457	934197	24.01%	2.843%
36	14.486	2106	2108	2112	rVB4	55453	51253	1.32%	0.156%

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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-BF081821.M
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37	14.551	2117	2119	2122	rBV	46046	44713	1.15%	0.136%
38	14.951	2184	2187	2190	rBV	129670	124448	3.20%	0.379%
39	15.221	2231	2233	2237	rVB3	46707	45286	1.16%	0.138%
40	15.592	2291	2296	2305	rVB2	801284	1042567	26.80%	3.173%

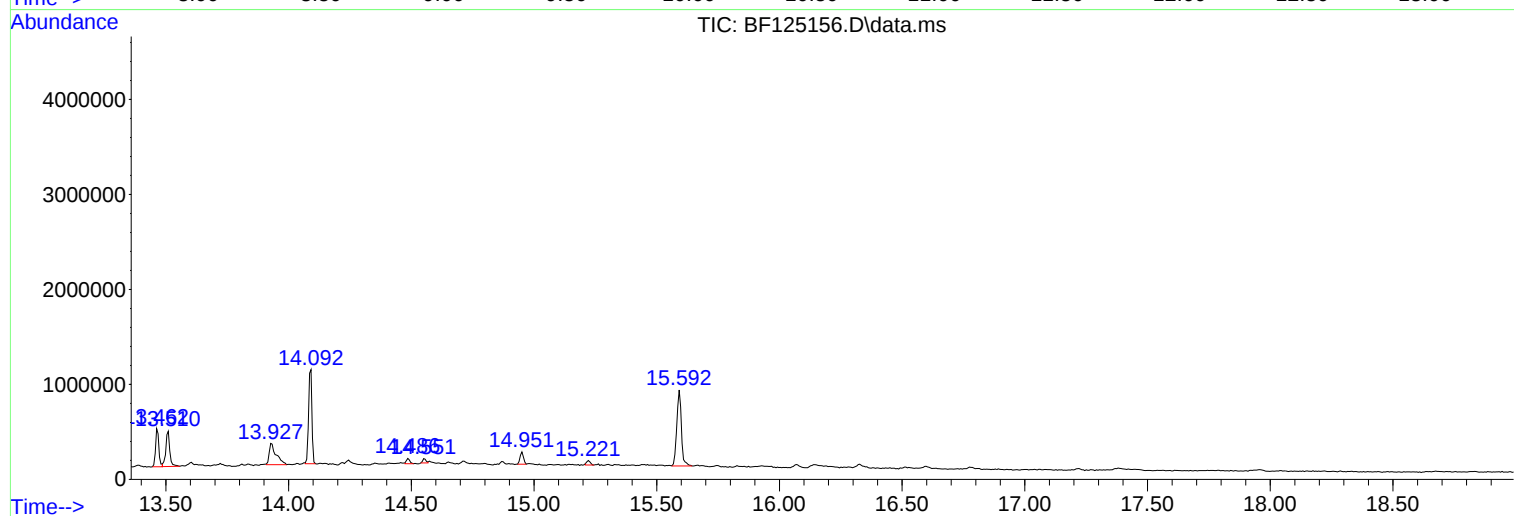
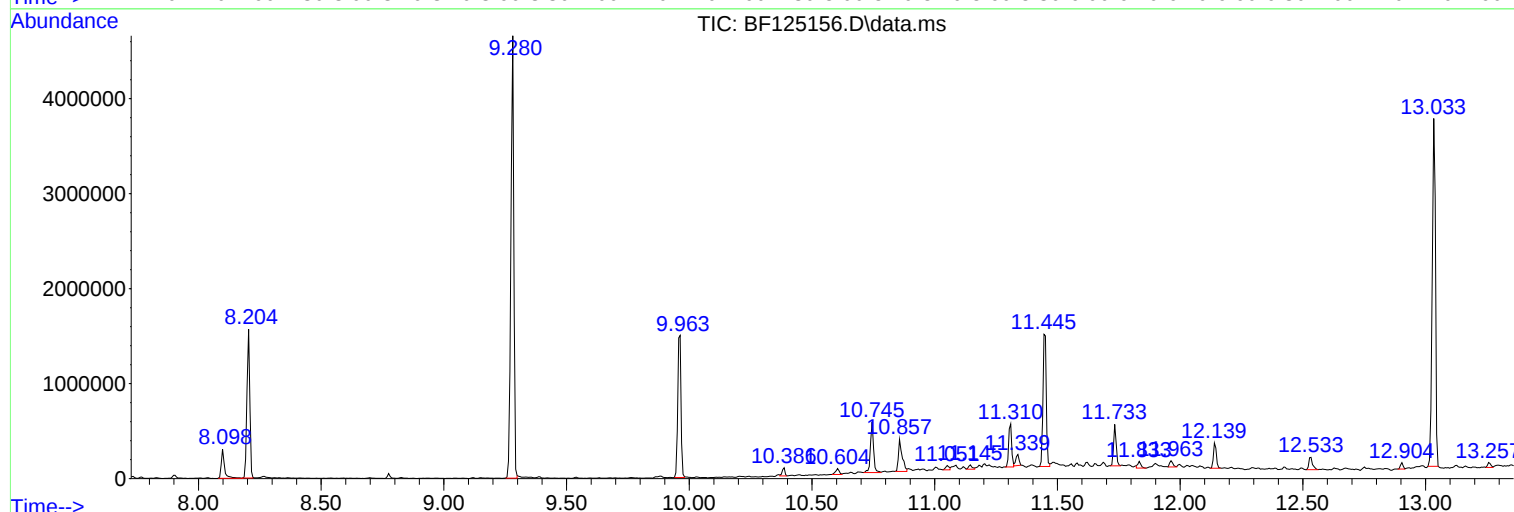
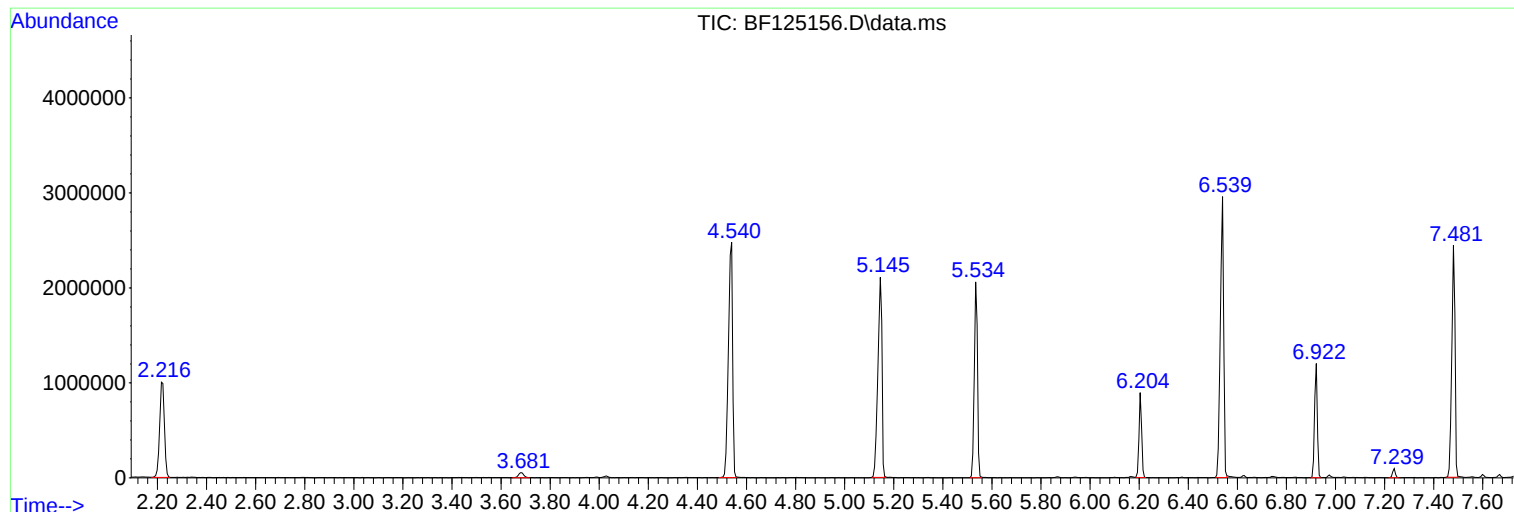
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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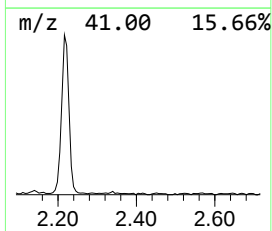
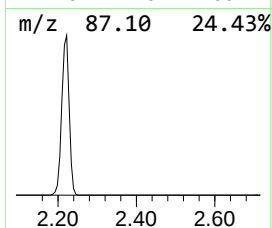
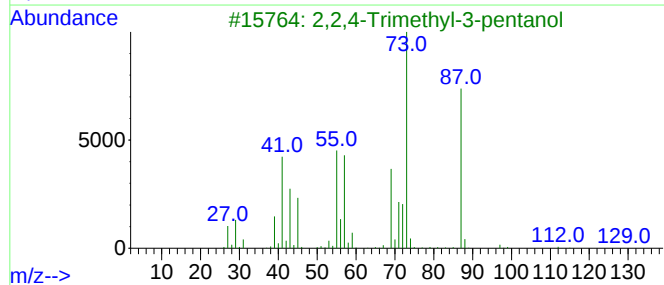
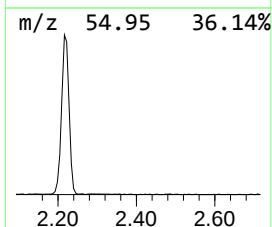
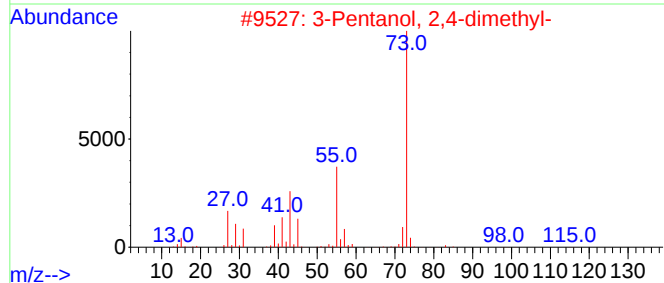
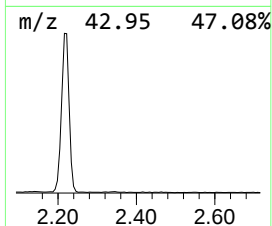
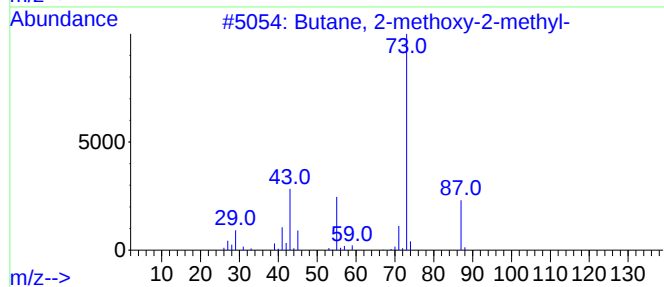
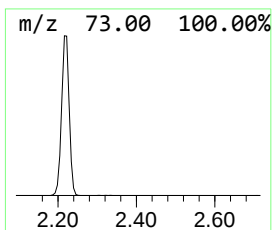
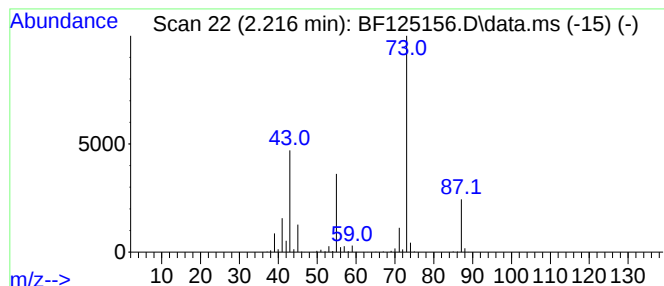
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.216	28.50 ng	1348050	1,4-Dichlorobenzene-d4	6.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25



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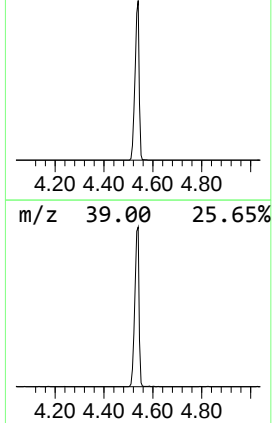
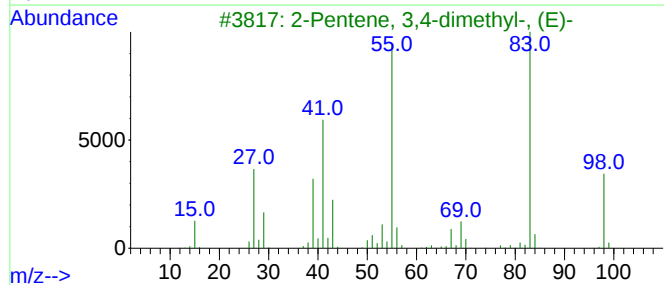
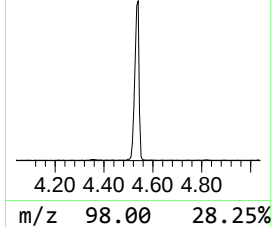
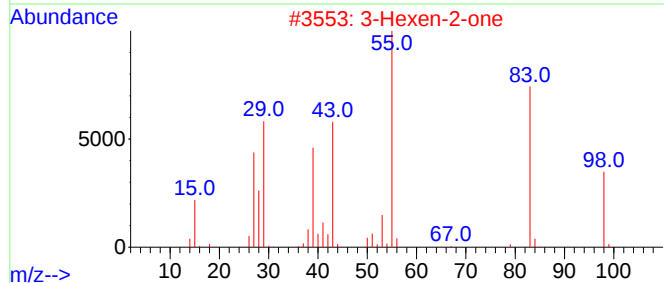
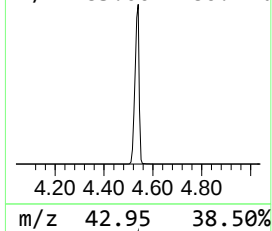
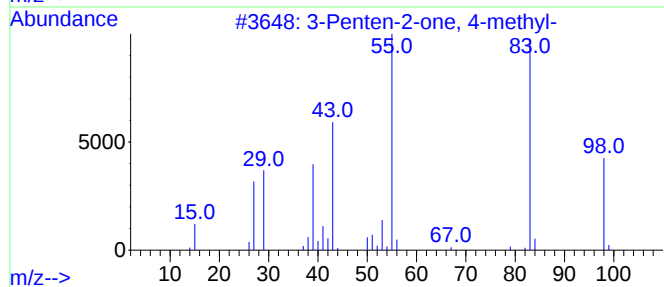
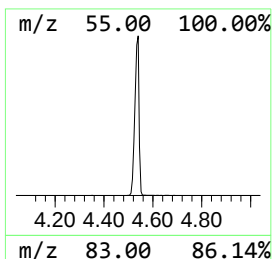
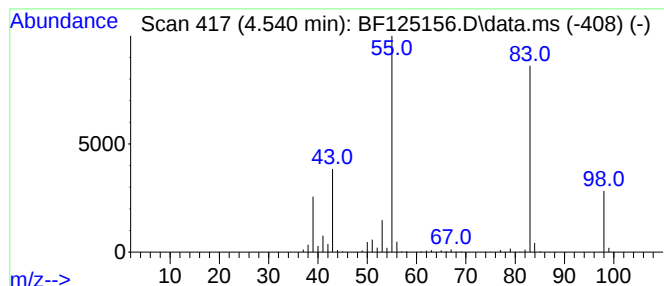
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Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.540	62.38 ng	2950430	1,4-Dichlorobenzene-d4	6.922

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	72



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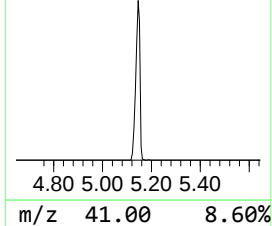
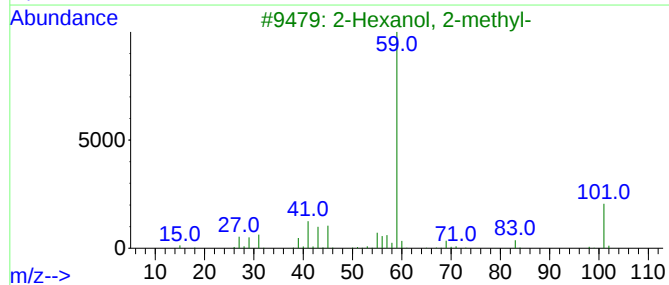
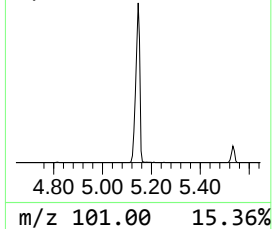
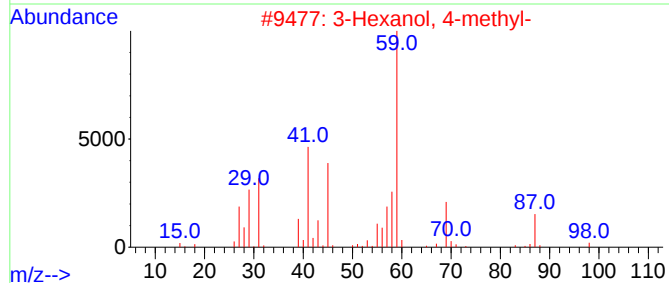
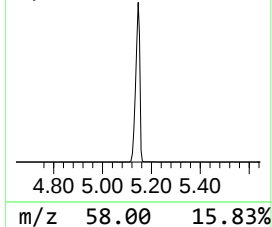
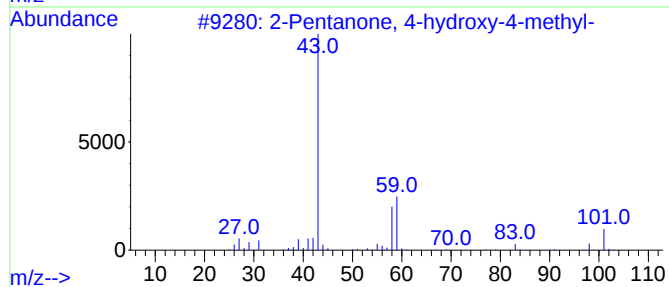
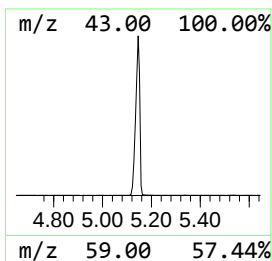
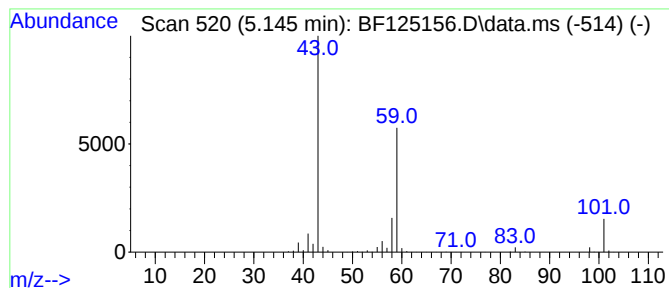
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.145	51.87 ng	2453460	1,4-Dichlorobenzene-d4	6.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28		
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	17		



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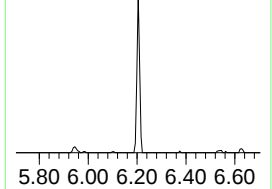
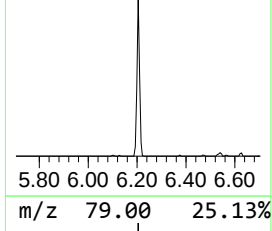
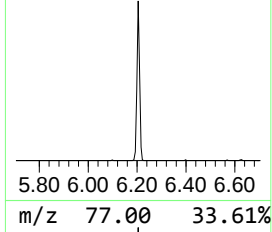
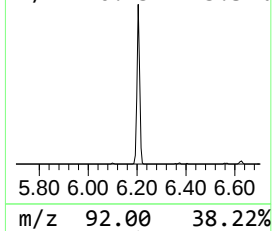
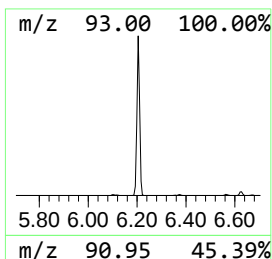
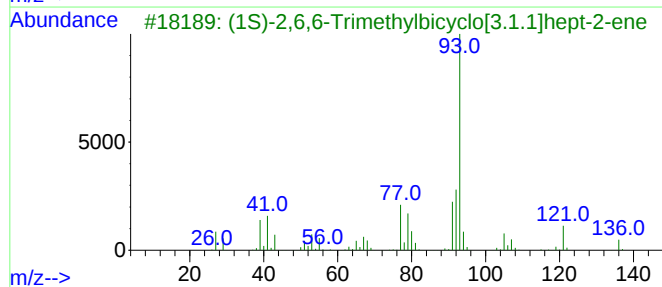
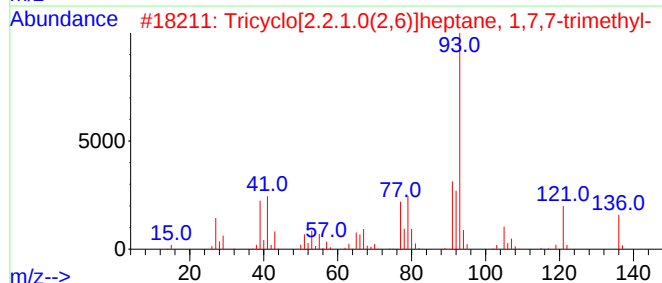
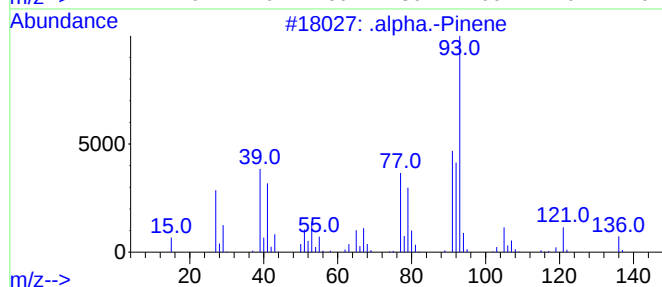
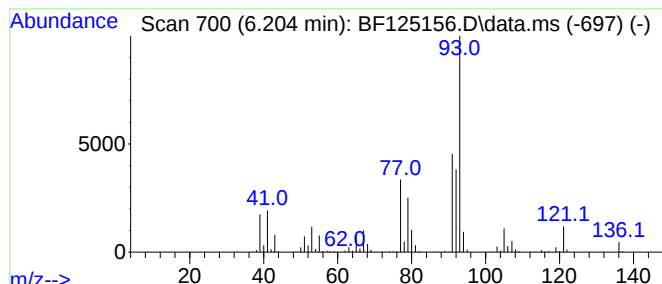
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 .alpha.-Pinene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.204	14.58 ng	689814	1,4-Dichlorobenzene-d4	6.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.		.alpha.-Pinene	136	C10H16	000080-56-8	95
2	Tricyclo[2.2.1.0(2,6)]heptane, 1...			136	C10H16	000508-32-7	91
3	(1S)-2,6,6-Trimethylbicyclo[3.1....			136	C10H16	007785-26-4	91
4	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...			136	C10H16	004889-83-2	91
5	(1R)-2,6,6-Trimethylbicyclo[3.1....			136	C10H16	007785-70-8	90



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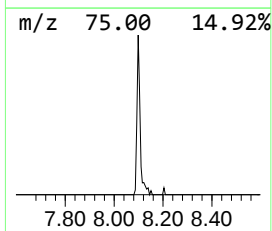
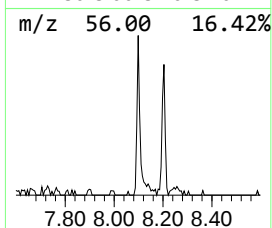
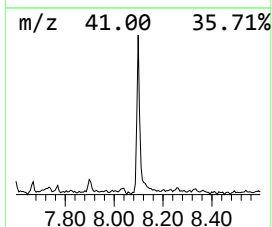
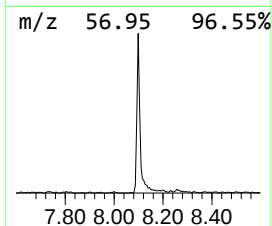
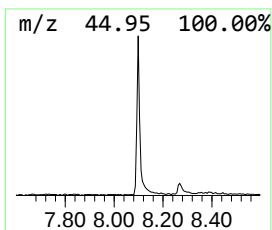
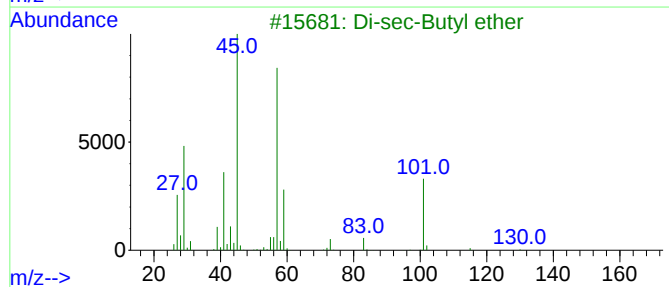
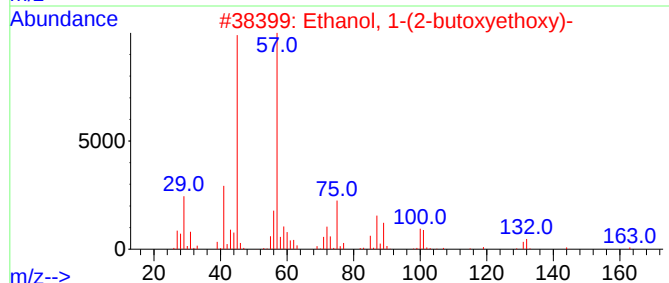
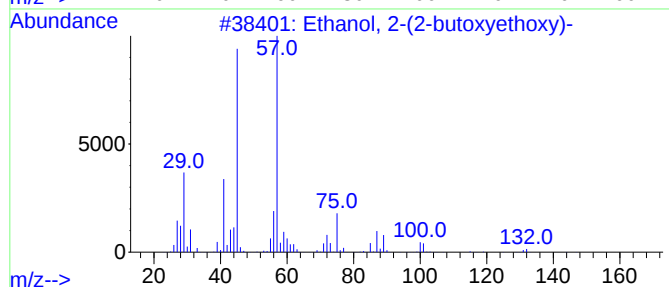
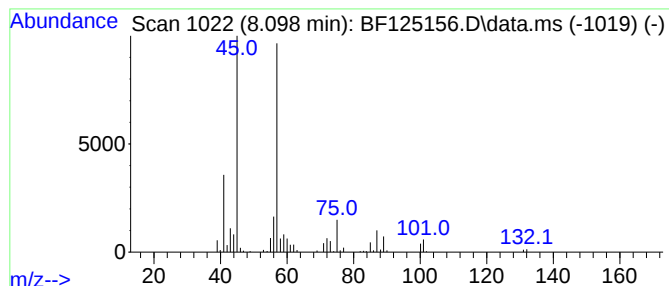
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.098	4.67 ng	286843	Naphthalene-d8	8.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Di-sec-Butyl ether	130	C8H18O	006863-58-7	59
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	50
5			1-Methoxy-3-methyl-3-butene	100	C6H12O	034752-58-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
Operator : JU/CG
Sample : M3465-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
CON-C

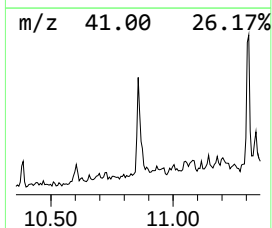
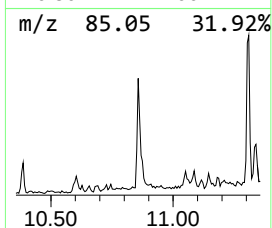
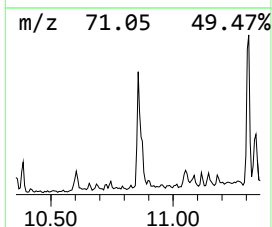
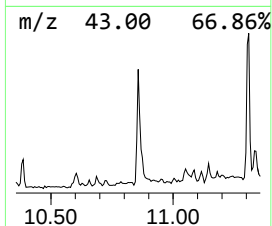
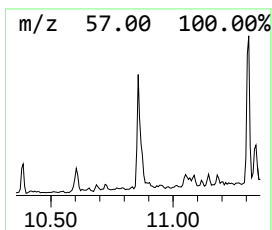
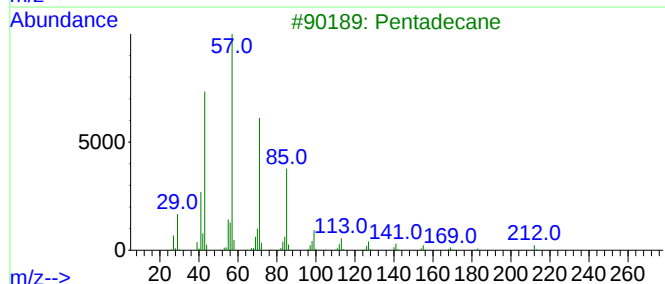
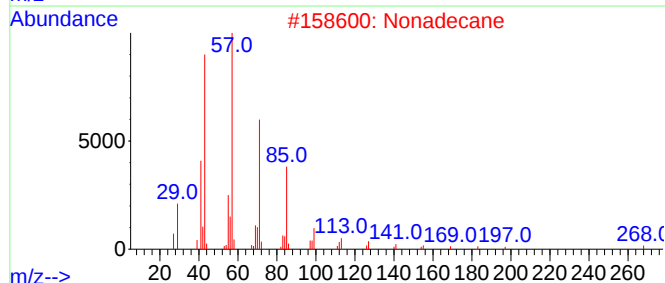
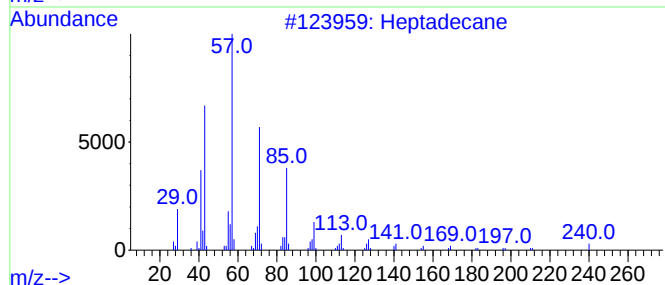
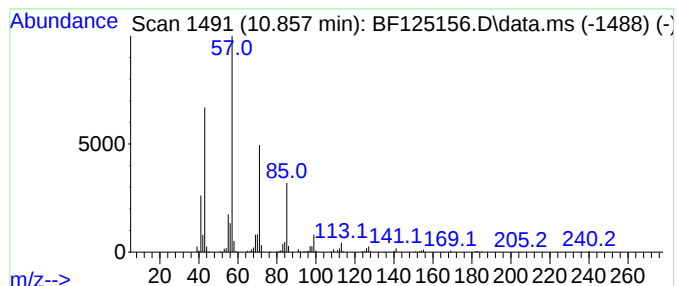
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.857	5.32 ng	348406	Phenanthrene-d10	11.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	91
2			Nonadecane	268	C19H40	000629-92-5	91
3			Pentadecane	212	C15H32	000629-62-9	90
4			Tridecane	184	C13H28	000629-50-5	90
5			Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
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Sample : M3465-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CON-C

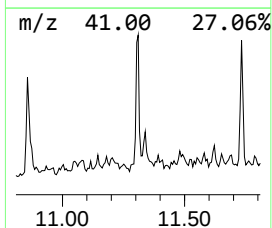
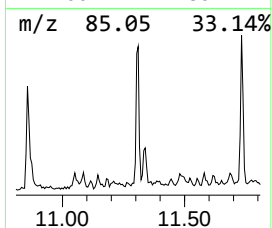
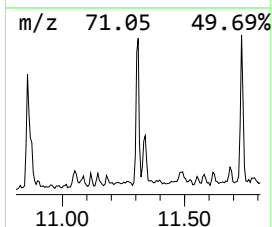
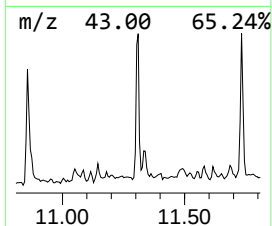
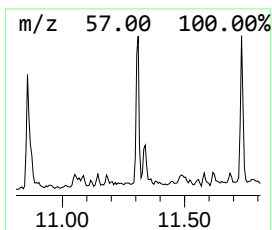
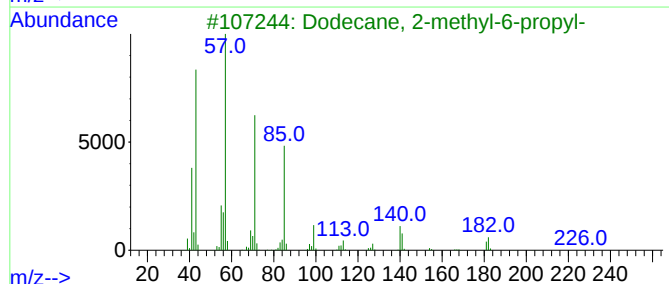
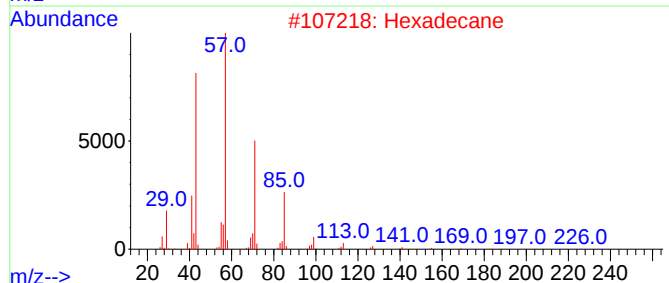
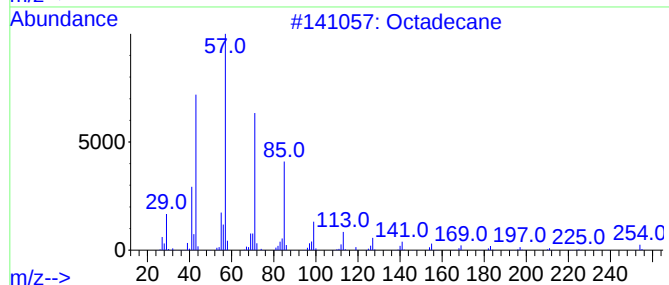
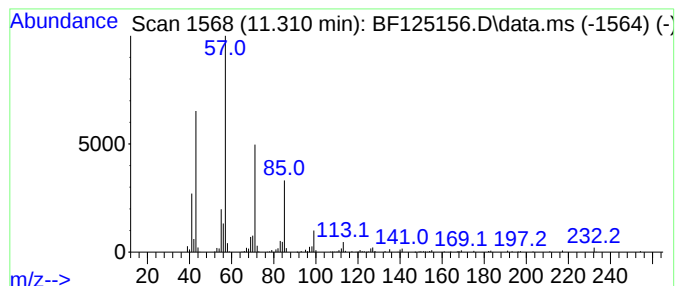
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.310	5.88 ng	385187	Phenanthrene-d10	11.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	94
2			Hexadecane	226	C16H34	000544-76-3	86
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
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Sample : M3465-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

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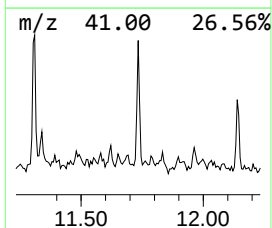
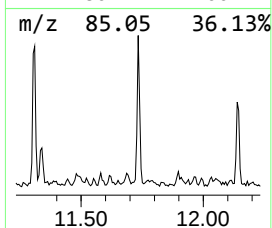
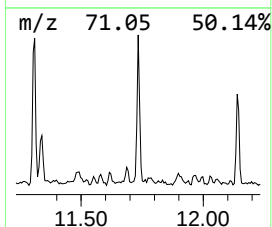
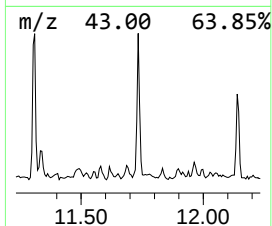
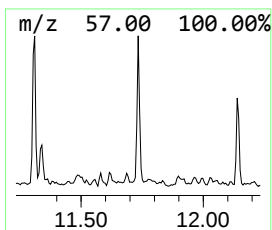
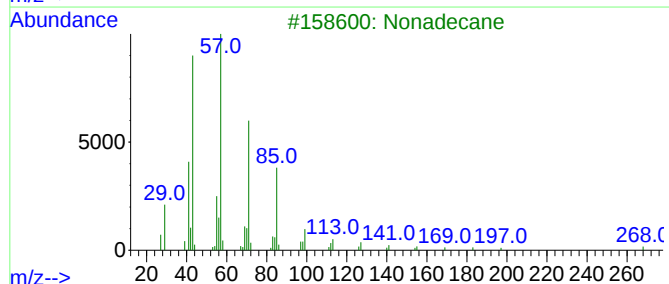
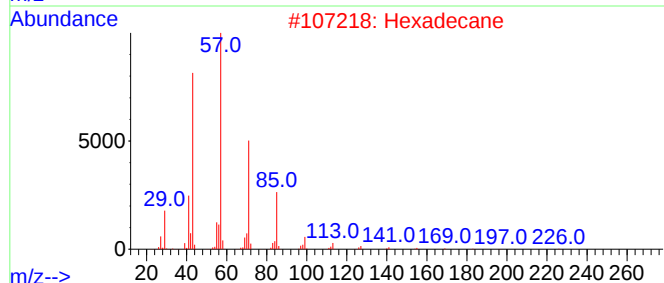
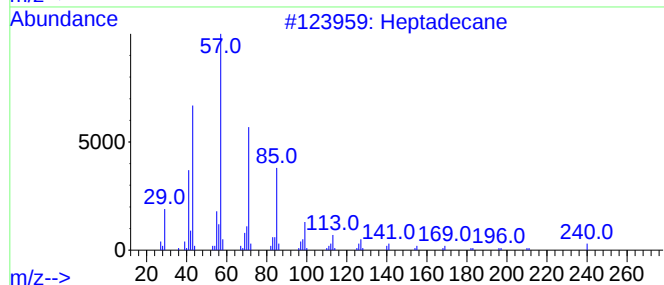
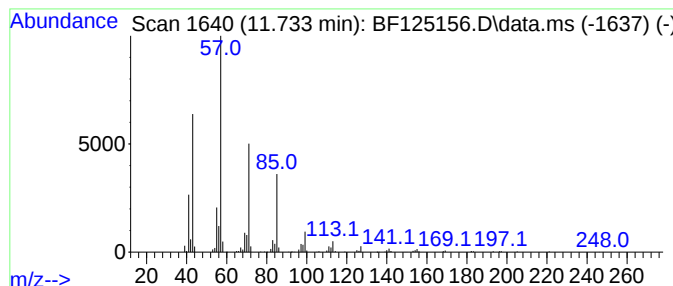
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.733	4.97 ng	325717	Phenanthrene-d10	11.451

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	90
2		Hexadecane	226	C16H34	000544-76-3	90
3		Nonadecane	268	C19H40	000629-92-5	86
4		Hexacosane	366	C26H54	000630-01-3	83
5		Heneicosane	296	C21H44	000629-94-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
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Misc :
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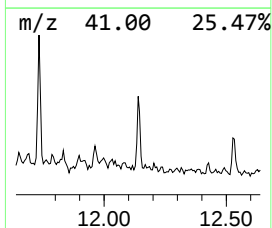
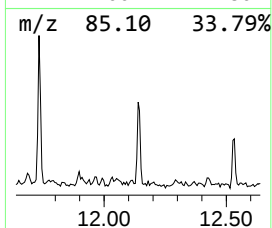
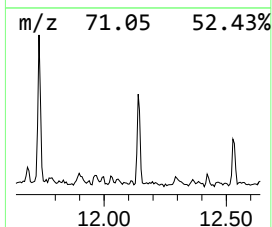
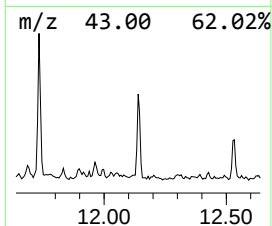
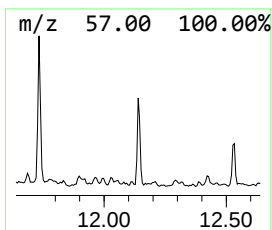
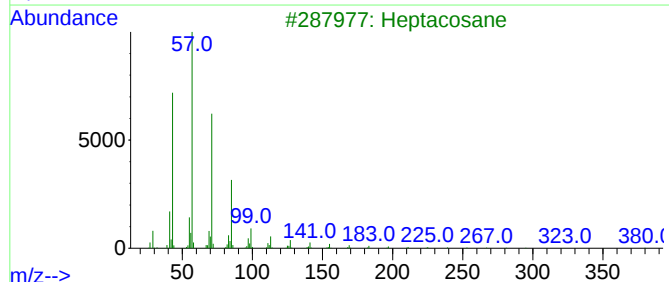
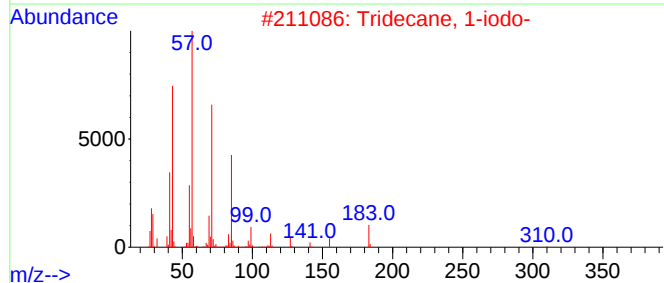
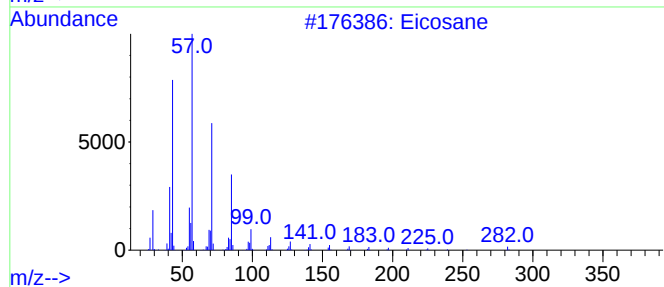
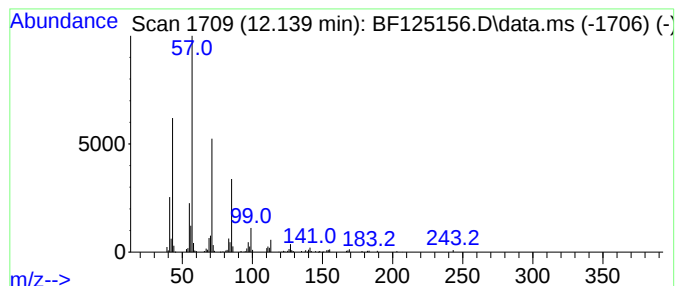
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.139	3.32 ng	217618	Phenanthrene-d10	11.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3			Heptacosane	380	C27H56	000593-49-7	87
4			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	86
5			Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-C

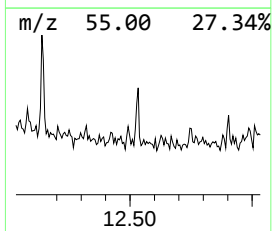
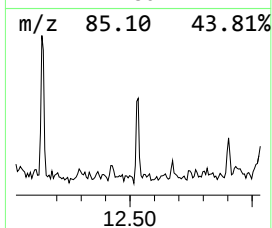
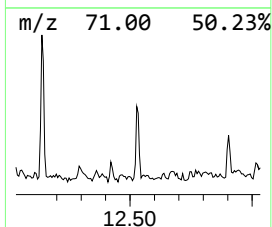
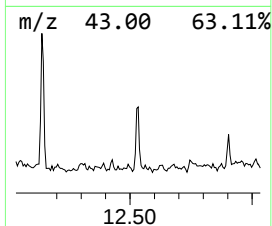
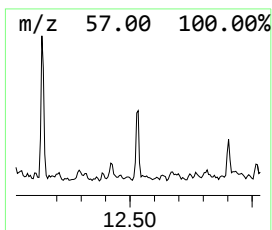
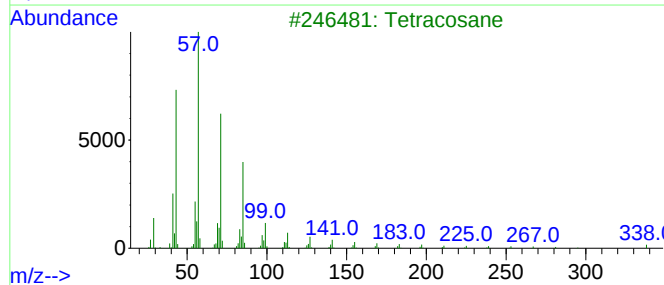
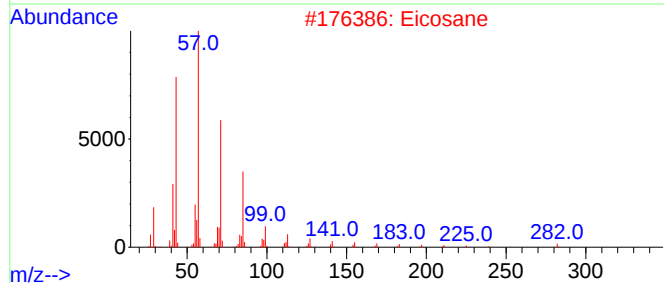
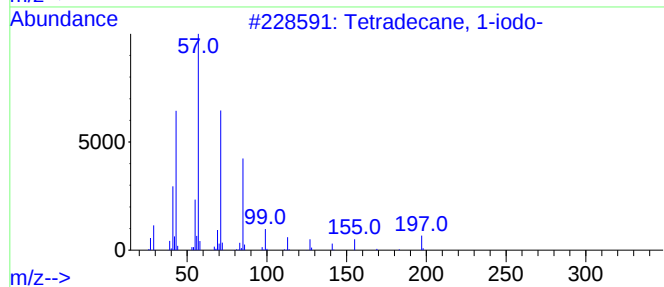
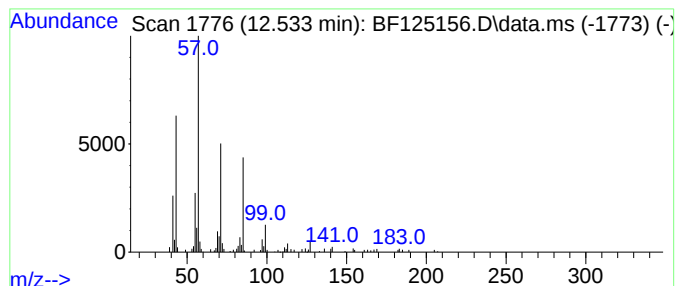
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecane, 1-iodo- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.533	2.12 ng	138955	Phenanthrene-d10	11.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	83
2			Eicosane	282	C20H42	000112-95-8	72
3			Tetracosane	338	C24H50	000646-31-1	72
4			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	72
5			Heptacosane	380	C27H56	000593-49-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
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Acq On : 23 Aug 2021 15:32
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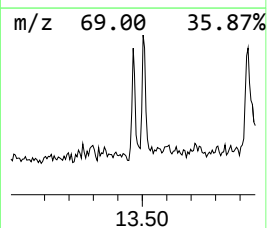
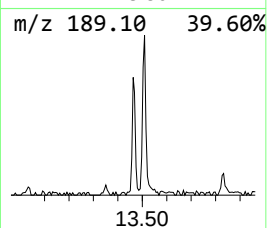
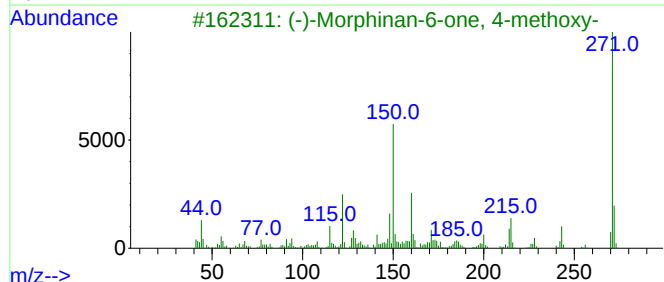
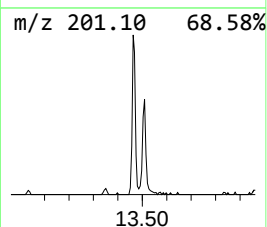
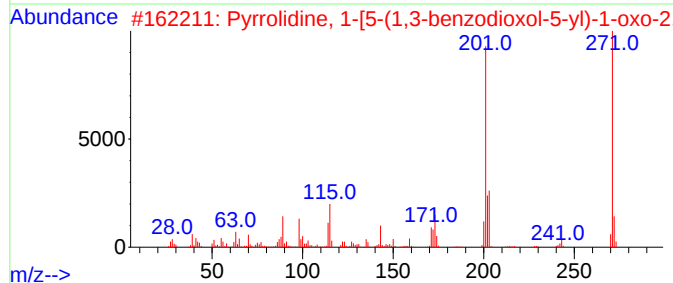
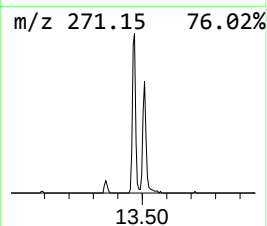
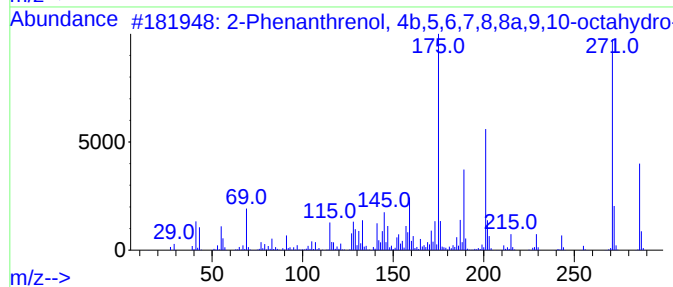
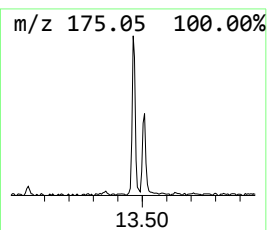
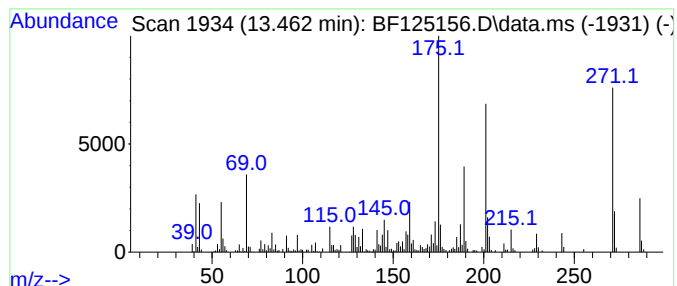
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.463	8.09 ng	377667	Chrysene-d12	14.092

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		Pyrrolidine, 1-[5-(1,3-benzodiox...	271	C16H17NO3	025924-78-1	38
3		(-)-Morphinan-6-one, 4-methoxy-	271	C17H21NO2	1000129-09-1	35
4		Naphthalene, 1,2,3,4-tetrahydro...	244	C18H28	029577-17-1	25
5		Crinan-1-one	271	C16H17NO3	098301-98-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
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Sample : M3465-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CON-C

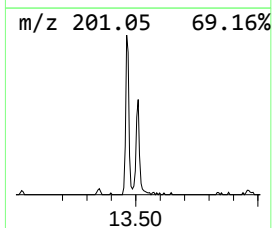
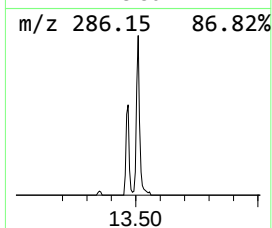
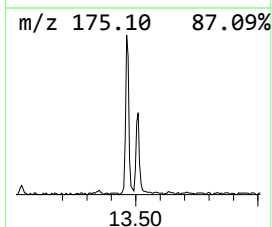
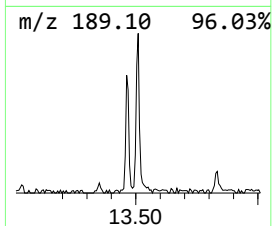
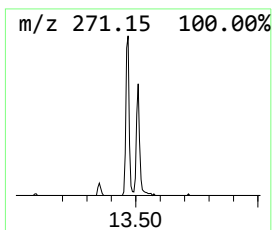
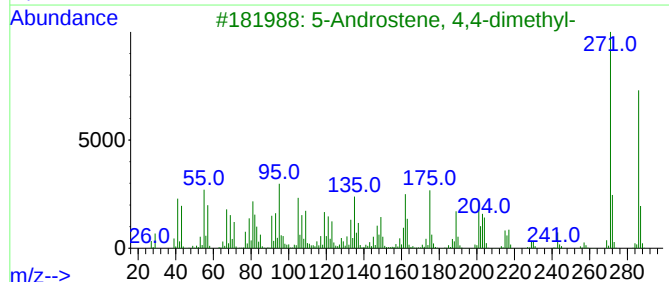
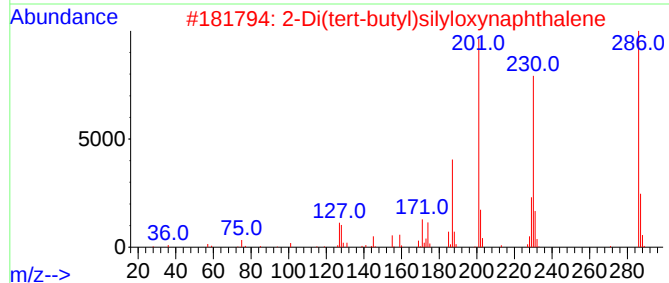
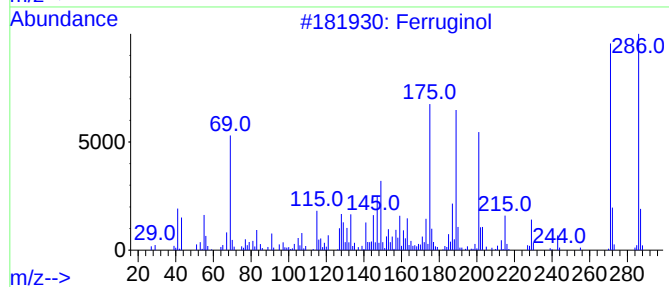
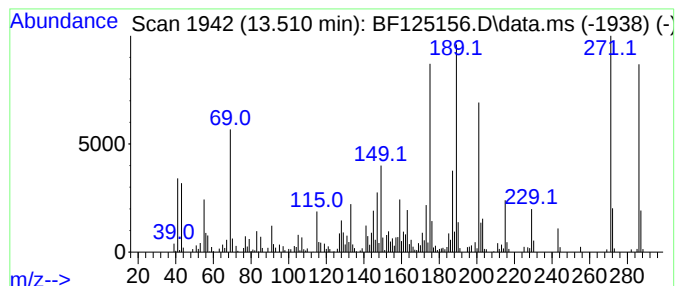
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Ferruginol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.510	8.74 ng	408143	Chrysene-d12	14.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ferruginol	286	C20H30O	000514-62-5	99
2			2-Di(tert-butyl)silyloxynaphthalene	286	C18H26OSi	1000307-94-6	50
3			5-Androstene, 4,4-dimethyl-	286	C21H34	1000194-15-4	43
4			5.beta.-Pregn-11-ene	286	C21H34	006673-73-0	38
5			2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
Operator : JU/CG
Sample : M3465-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-C

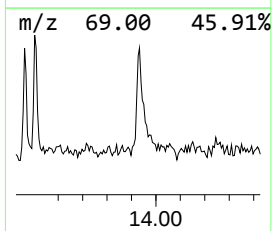
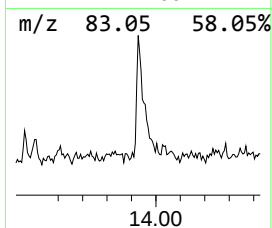
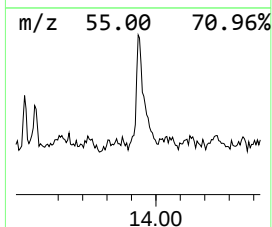
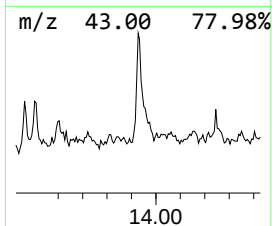
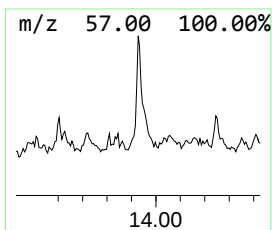
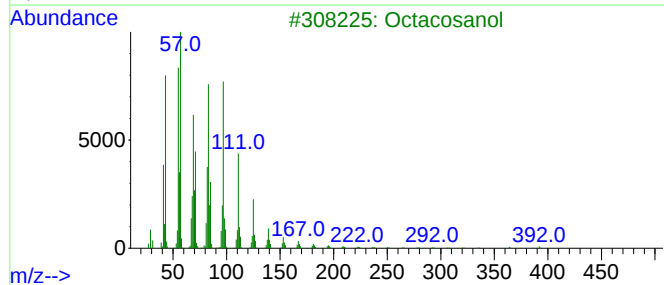
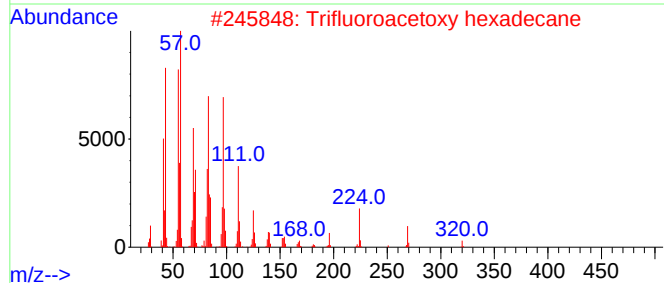
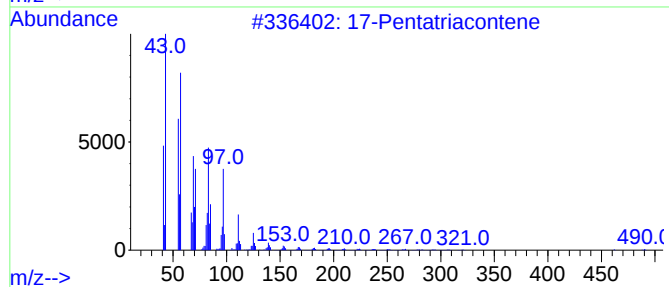
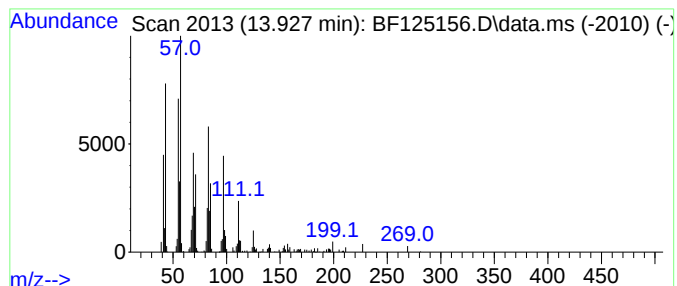
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 17-Pentatriacontene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.927	9.37 ng	437539	Chrysene-d12	14.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			17-Pentatriacontene	491	C35H70	006971-40-0	91
2			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
3			Octacosanol	410	C28H58O	000557-61-9	86
4			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	80
5			1-Octadecanol	270	C18H38O	000112-92-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
Data File : BF125156.D
Acq On : 23 Aug 2021 15:32
Operator : JU/CG
Sample : M3465-05
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-C

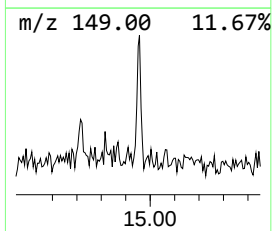
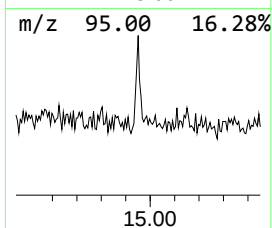
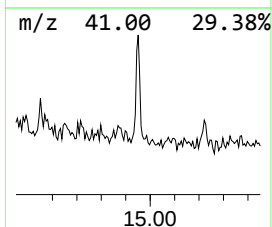
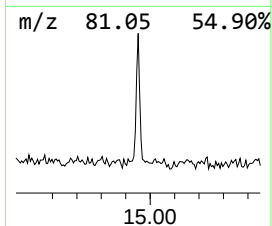
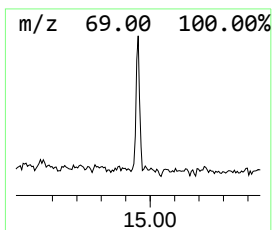
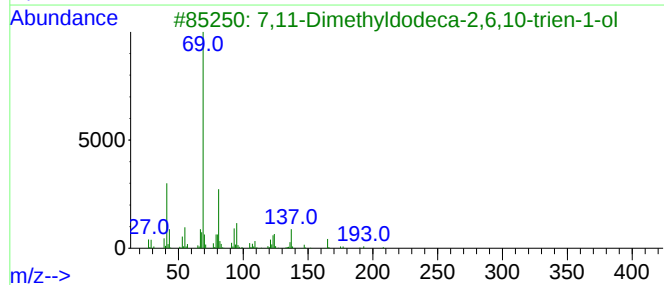
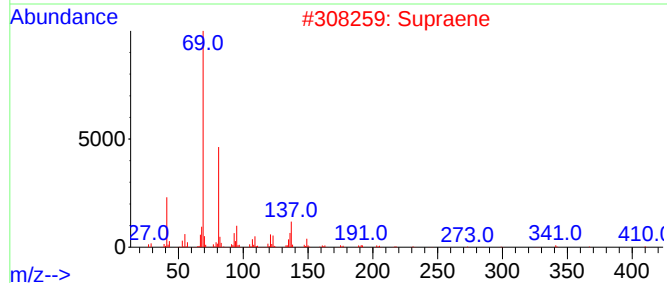
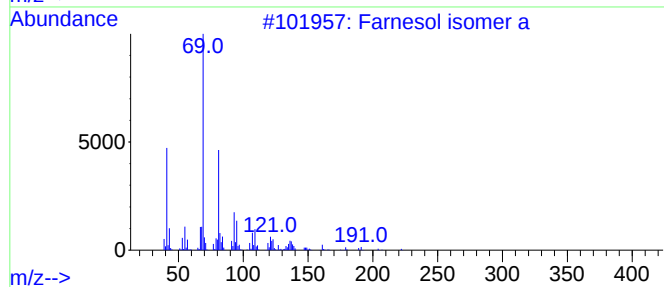
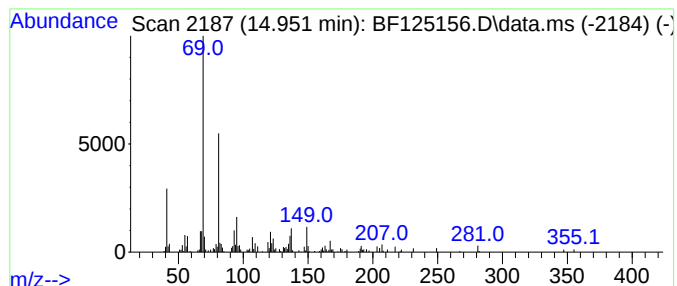
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Farnesol isomer a Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.951	2.39 ng	124448	Perylene-d12	15.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Farnesol isomer a	222	C15H26O	1000108-92-4	81
2			Supraene	410	C30H50	007683-64-9	80
3			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	74
4			2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	72
5			3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
 Data File : BF125156.D
 Acq On : 23 Aug 2021 15:32
 Operator : JU/CG
 Sample : M3465-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CON-C

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.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...	2.216	28.5	ng	1348050	1	6.922	946003	20.0
3-Penten-2-one,...	4.540	62.4	ng	2950430	1	6.922	946003	20.0
2-Pentanone, 4-...	5.145	51.9	ng	2453460	1	6.922	946003	20.0
.alpha.-Pinene	6.204	14.6	ng	689814	1	6.922	946003	20.0
Ethanol, 2-(2-b...	8.098	4.7	ng	286843	2	8.204	1229620	20.0
Heptadecane	10.857	5.3	ng	348406	4	11.451	1310290	20.0
Octadecane	11.310	5.9	ng	385187	4	11.451	1310290	20.0
Hexadecane	11.733	5.0	ng	325717	4	11.451	1310290	20.0
Eicosane	12.139	3.3	ng	217618	4	11.451	1310290	20.0
Tetradecane, 1-...	12.533	2.1	ng	138955	4	11.451	1310290	20.0
2-Phenanthrenol...	13.463	8.1	ng	377667	5	14.092	934197	20.0
Ferruginol	13.510	8.7	ng	408143	5	14.092	934197	20.0
17-Pentatriacon...	13.927	9.4	ng	437539	5	14.092	934197	20.0
Farnesol isomer a	14.951	2.4	ng	124448	6	15.592	1042570	20.0