

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100920\
 Data File : BF121906.D
 Acq On : 9 Oct 2020 15:09
 Operator : JU/CG
 Sample : L4329-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-1

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF091020.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.963	485	489	499	rBV	58588	88050	3.68%	0.477%
2	5.381	556	560	582	rBV	1951174	2021953	84.58%	10.950%
3	6.398	729	733	761	rBV	1930446	2069998	86.59%	11.210%
4	6.598	764	767	785	rVV	38212	71572	2.99%	0.388%
5	6.757	790	794	801	rVB	695210	596611	24.96%	3.231%
6	7.322	886	890	902	rBV	1450581	1420984	59.44%	7.695%
7	8.040	1008	1012	1032	rBV	807883	802890	33.59%	4.348%
8	9.116	1190	1195	1198	rBV	2772166	2390520	100.00%	12.946%
9	9.510	1258	1262	1270	rBV	374445	325511	13.62%	1.763%
10	9.792	1306	1310	1314	rBV	988843	862030	36.06%	4.668%
11	10.339	1399	1403	1408	rBV	26665	29744	1.24%	0.161%
12	10.581	1440	1444	1454	rBV	1433779	1446375	60.50%	7.833%
13	11.275	1558	1562	1565	rBV	859435	841683	35.21%	4.558%
14	11.298	1565	1566	1570	rVB	137572	103482	4.33%	0.560%
15	11.810	1650	1653	1657	rBV2	31514	30957	1.29%	0.168%
16	12.492	1765	1769	1772	rBV	131346	119026	4.98%	0.645%
17	12.527	1772	1775	1781	rVB5	20388	27432	1.15%	0.149%
18	12.722	1802	1808	1813	rVB	86486	106285	4.45%	0.576%
19	12.863	1827	1832	1835	rBV	2106346	1990218	83.25%	10.778%
20	13.092	1867	1871	1873	rBV3	30238	36038	1.51%	0.195%
21	13.769	1982	1986	1998	rBV	258486	499325	20.89%	2.704%
22	13.916	2007	2011	2014	rBV	733269	678294	28.37%	3.673%
23	14.380	2086	2090	2092	rBV	51050	51377	2.15%	0.278%
24	14.410	2092	2095	2106	rVB4	105009	209423	8.76%	1.134%
25	14.674	2134	2140	2146	rBV7	48040	106982	4.48%	0.579%
26	14.998	2192	2195	2200	rVB	121259	133931	5.60%	0.725%
27	15.068	2201	2207	2213	rVB6	43687	107588	4.50%	0.583%
28	15.233	2232	2235	2241	rVB	21506	33002	1.38%	0.179%
29	15.292	2243	2245	2249	rVB	27198	28134	1.18%	0.152%
30	15.357	2251	2256	2264	rVB	552297	744918	31.16%	4.034%
31	15.774	2323	2327	2333	rVB2	73082	107213	4.48%	0.581%
32	15.868	2337	2343	2345	rBV3	63080	126154	5.28%	0.683%
33	17.357	2592	2596	2604	rVB8	59588	130894	5.48%	0.709%
34	17.639	2638	2644	2661	rVB8	32167	127043	5.31%	0.688%

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ClientSampleId :
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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-BF091020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

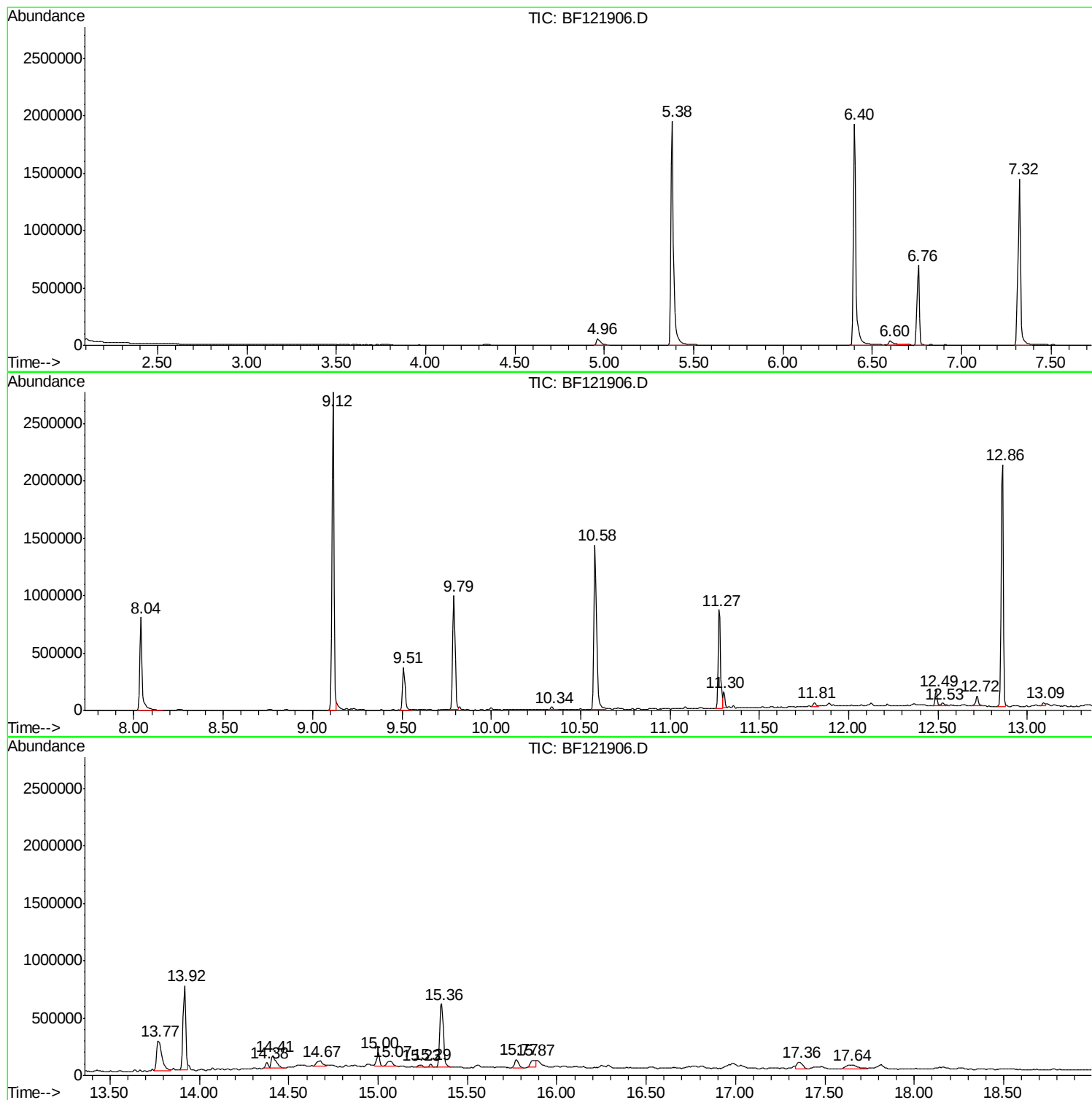
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF091020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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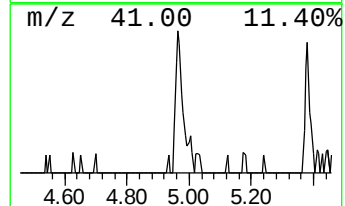
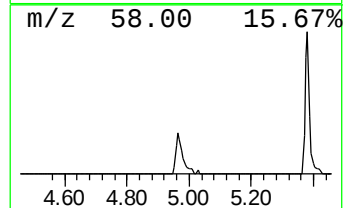
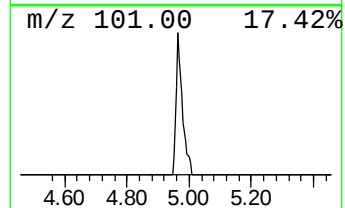
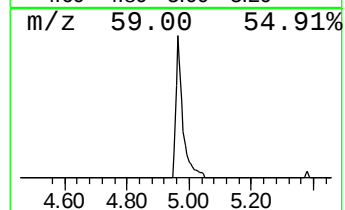
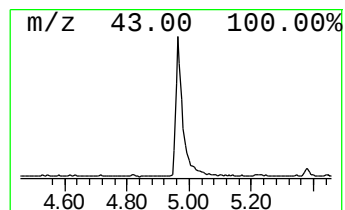
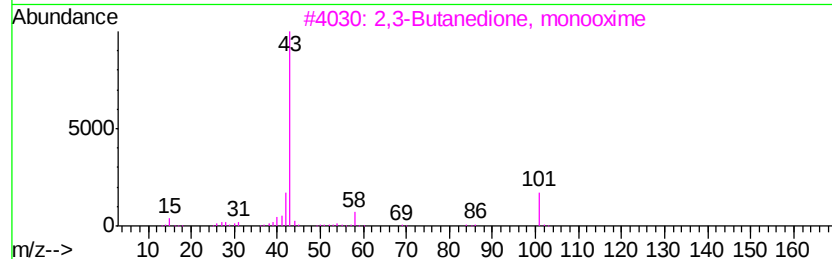
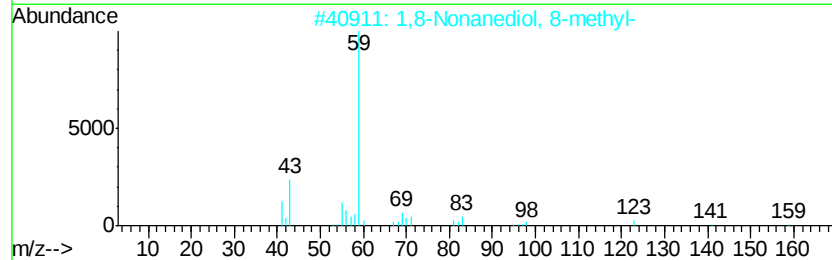
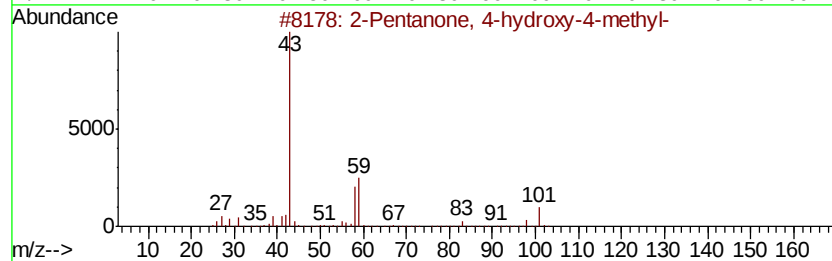
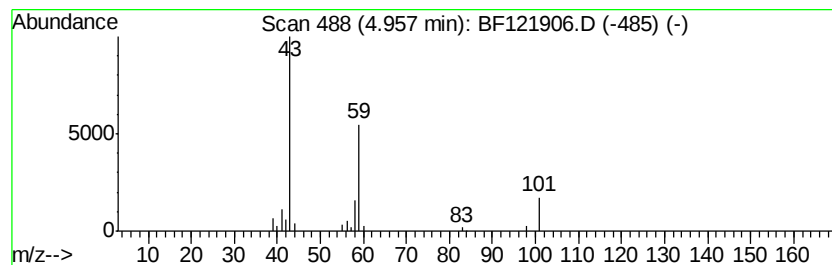
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	2.95 ng	88050	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	9
5			Acetic acid, 1,1-dimethylethyl ester	116	C6H12O2	000540-88-5	9



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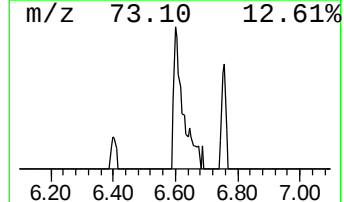
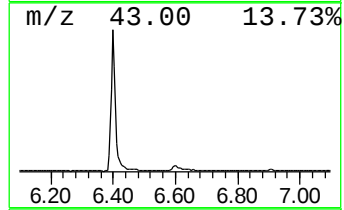
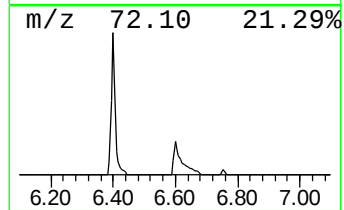
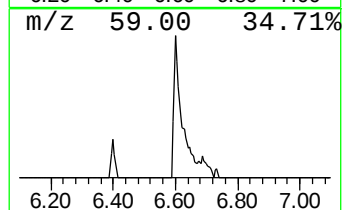
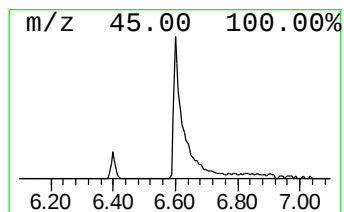
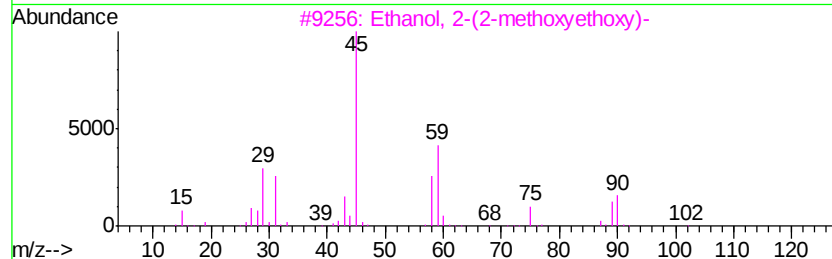
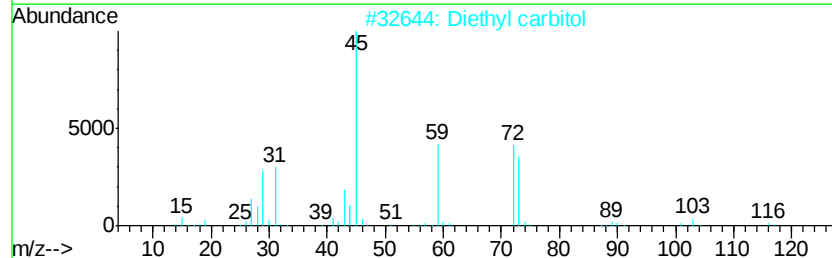
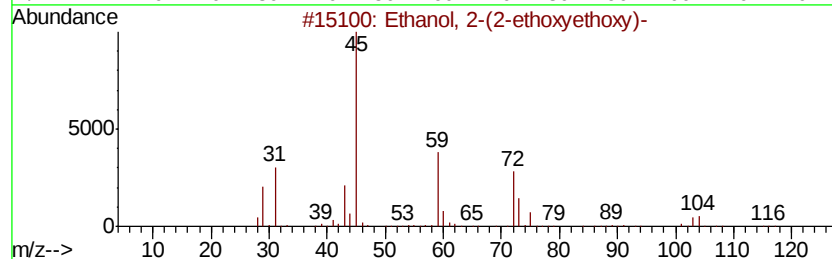
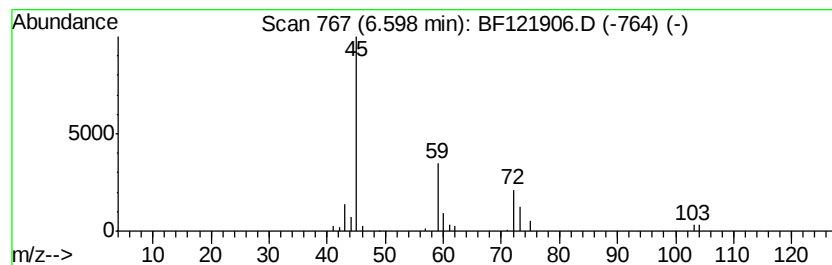
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	2.40 ng	71572	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
2		Diethyl carbitol	162	C8H18O3	000112-36-7	72
3		Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	45
4		2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	38
5		.beta.-Methoxyethoxymethyl chloride	124	C4H9ClO2	003970-21-6	36



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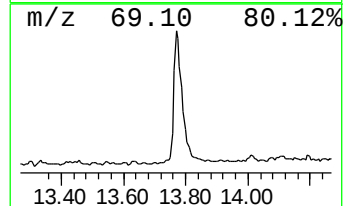
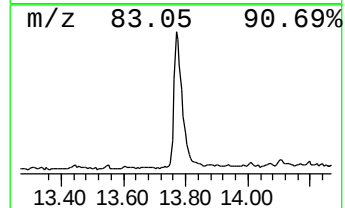
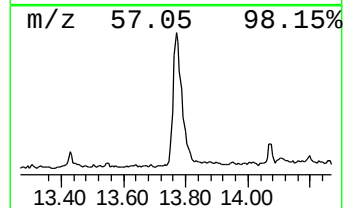
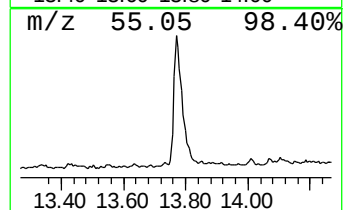
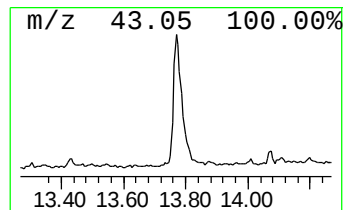
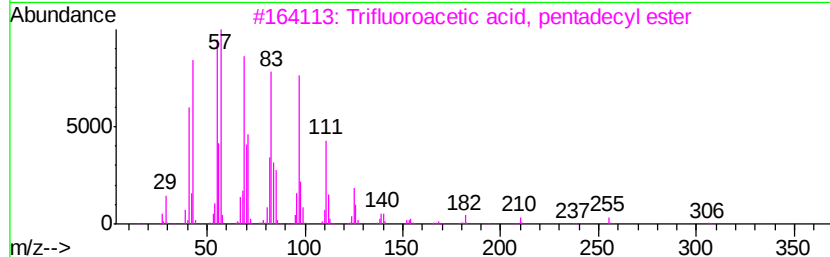
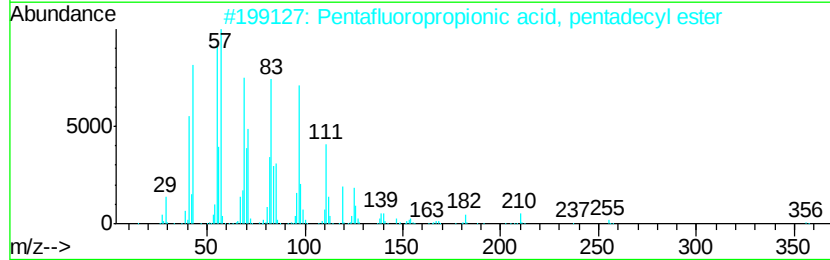
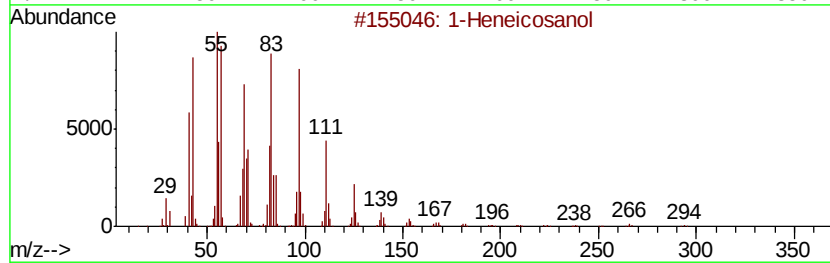
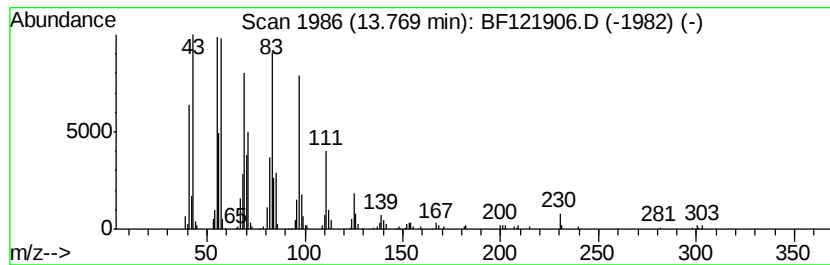
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Heneicosanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.77	14.72 ng	499325	Chrysene-d12	13.92		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	95
2		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
3		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
4		Heptafluorobutvric acid, pentade...	424	C19H31F7O2	959261-23-5	94
5		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94



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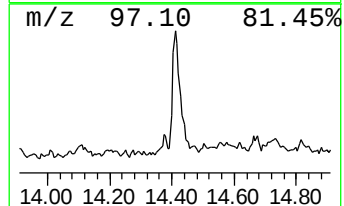
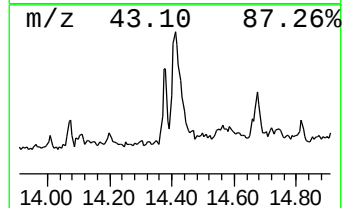
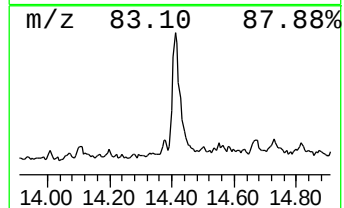
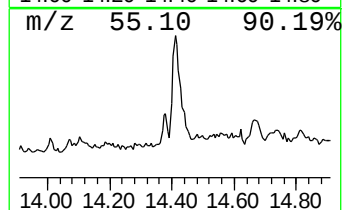
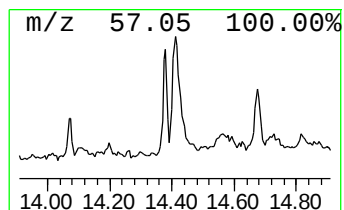
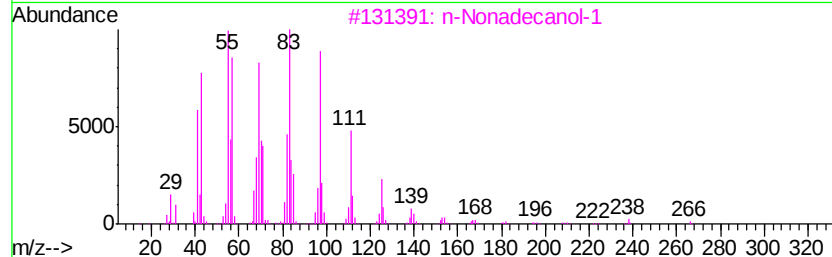
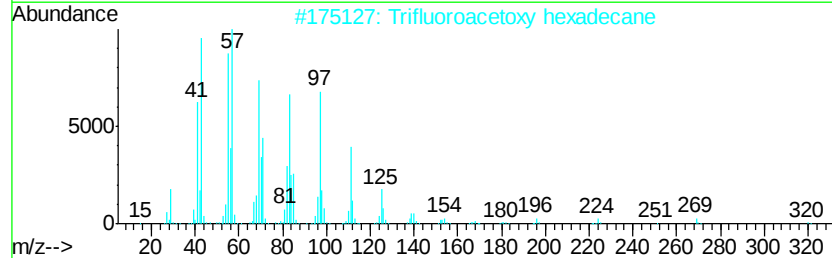
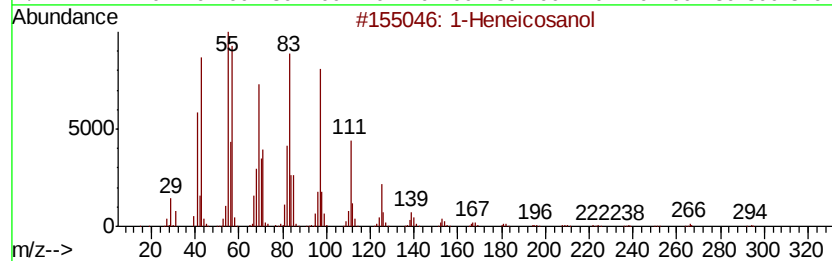
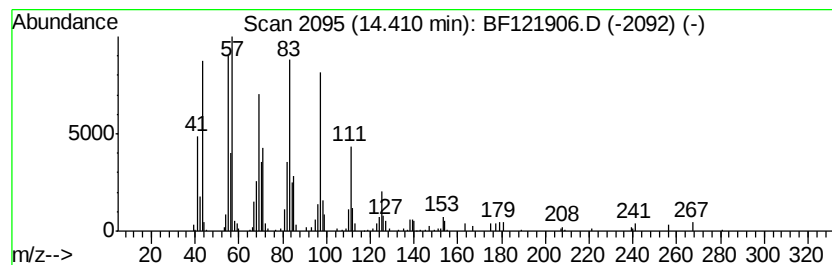
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Trifluoroacetoxy hexadecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	6.17 ng	209423	Chrysene-d12	13.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	95
2		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	93
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93



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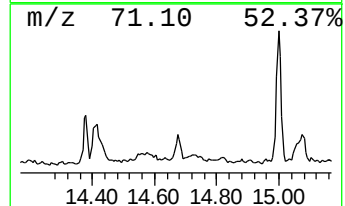
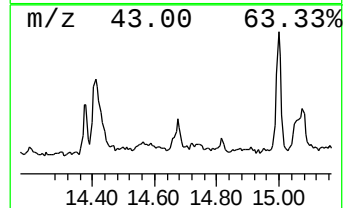
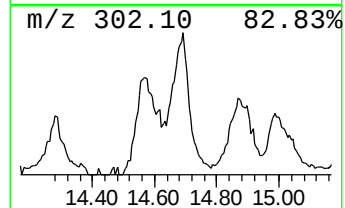
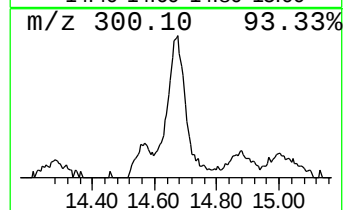
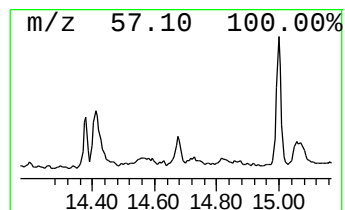
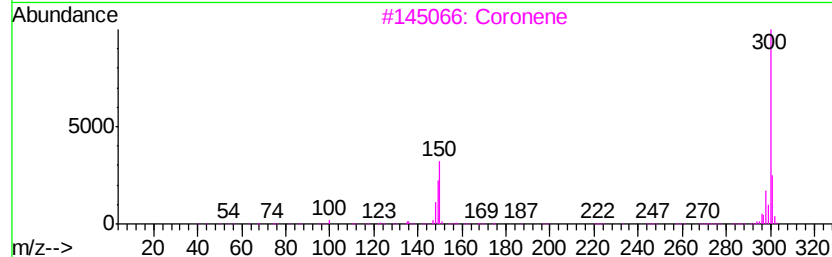
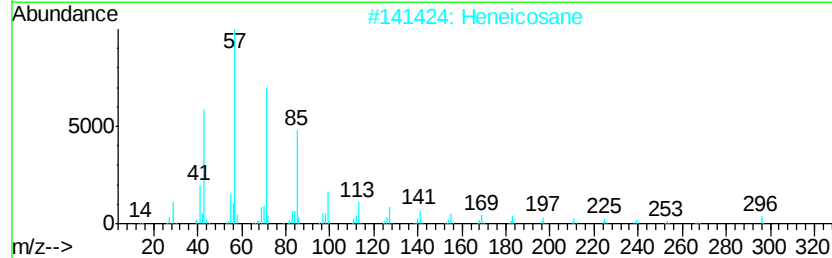
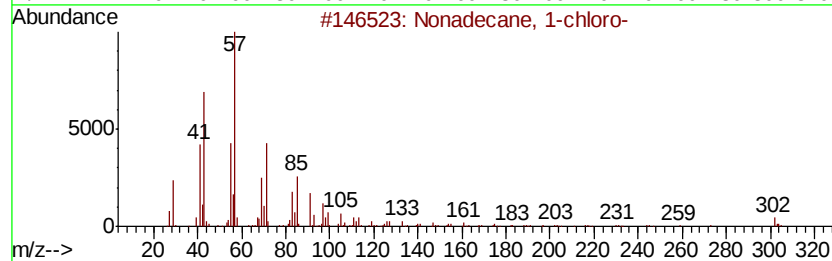
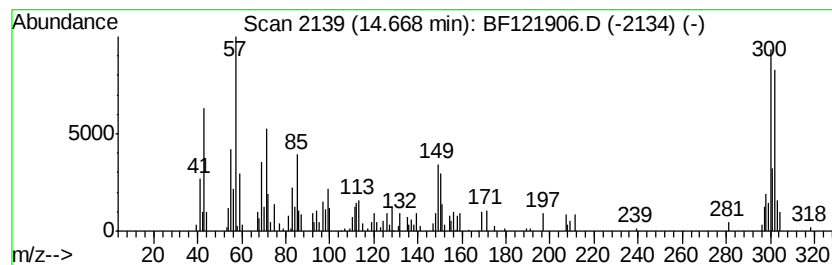
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown14.67 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	2.87 ng	106982	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	45
2		Heneicosane	296	C21H44	000629-94-7	44
3		Coronene	300	C24H12	000191-07-1	41
4		1,2,9,10-Dibenzopyrene	302	C24H14	000191-30-0	38
5		Heptacosane	380	C27H56	000593-49-7	30



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ALS Vial : 7 Sample Multiplier: 1

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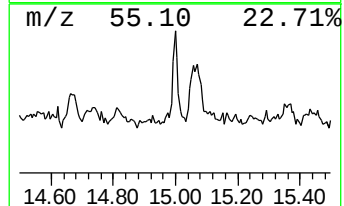
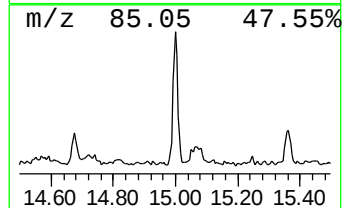
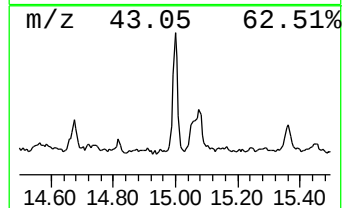
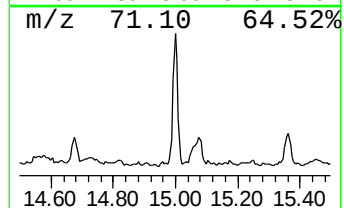
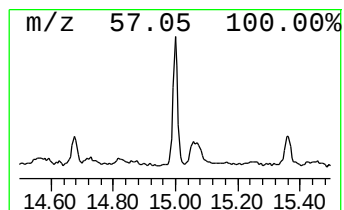
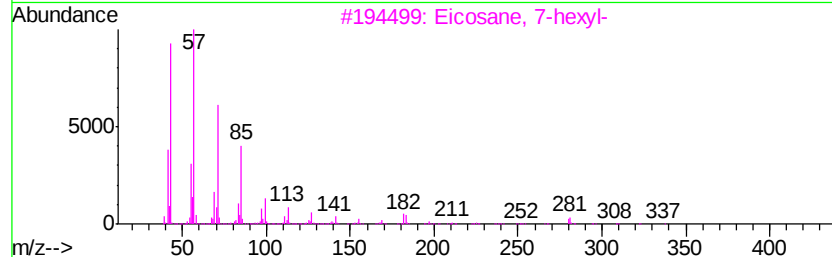
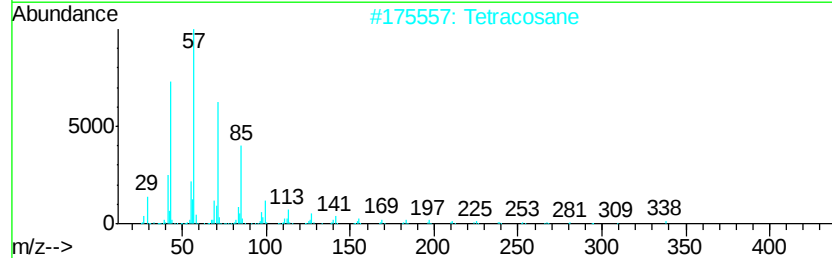
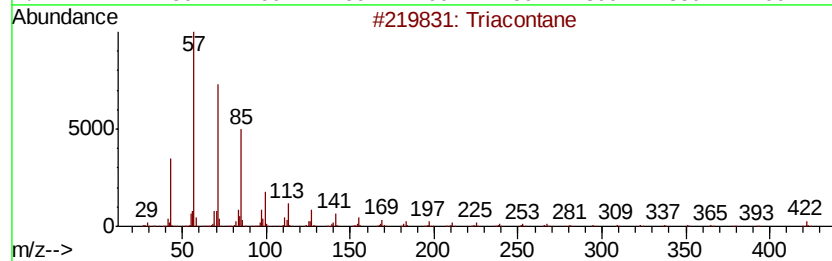
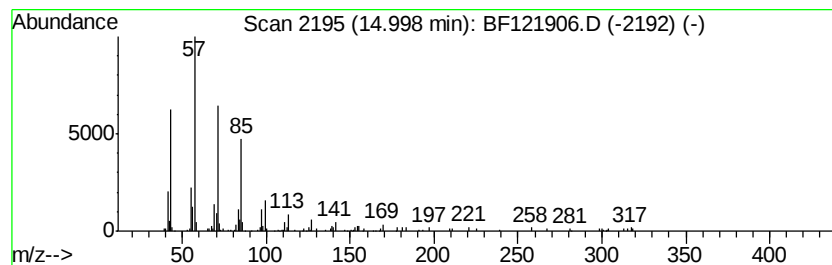
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Triacontane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	3.60 ng	133931	Perylene-d12	15.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontane	422	C30H62	000638-68-6	87
2		Tetracosane	338	C24H50	000646-31-1	87
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100920\
Data File : BF121906.D
Acq On : 9 Oct 2020 15:09
Operator : JU/CG
Sample : L4329-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-1

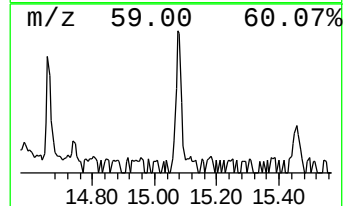
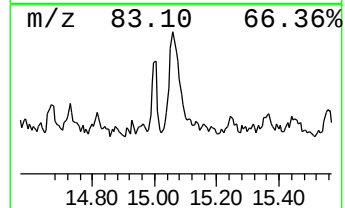
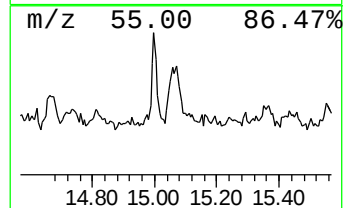
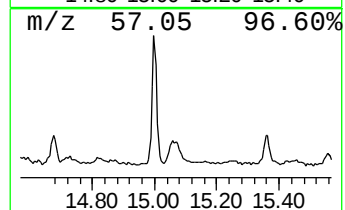
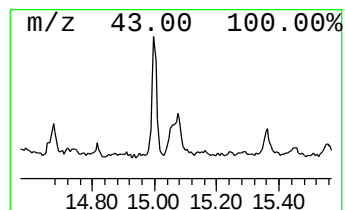
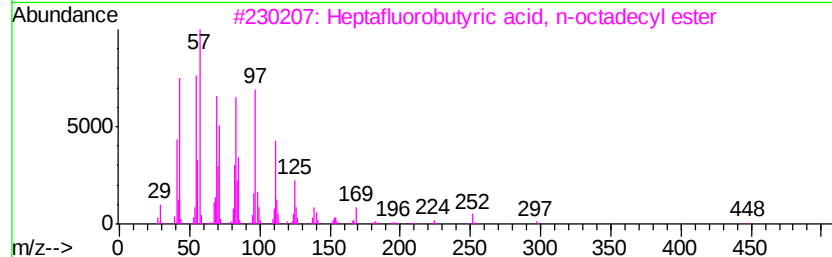
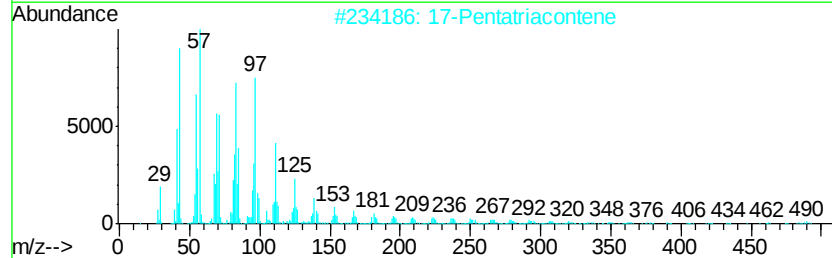
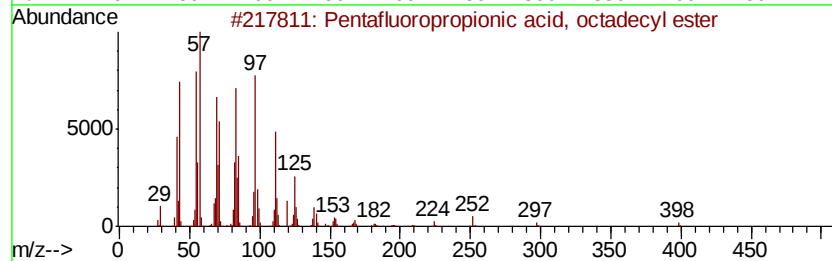
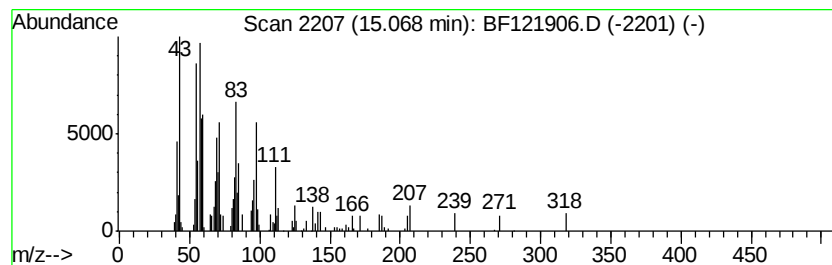
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentafluoropropionic acid, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	2.89 ng	107588	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	64
2		17-Pentatriacontene	491	C35H70	006971-40-0	64
3		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	64
4		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	58
5		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100920\
Data File : BF121906.D
Acq On : 9 Oct 2020 15:09
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Sample : L4329-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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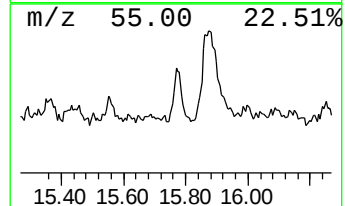
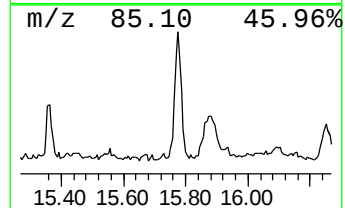
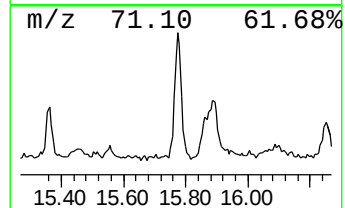
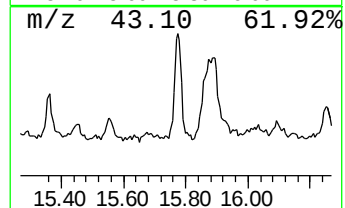
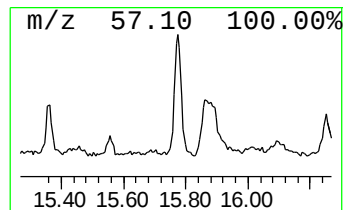
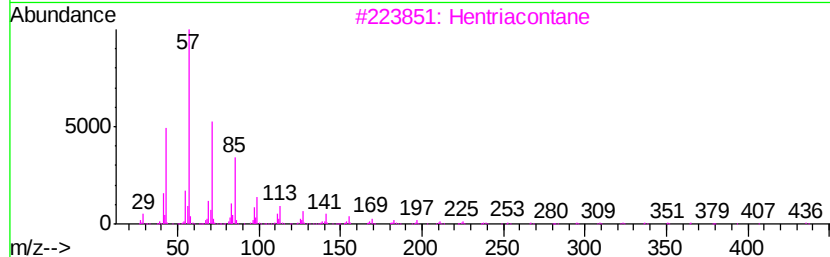
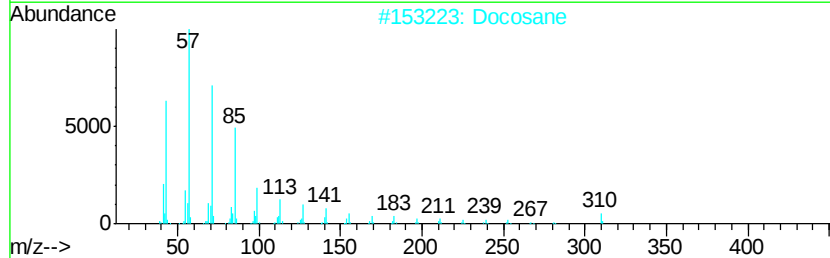
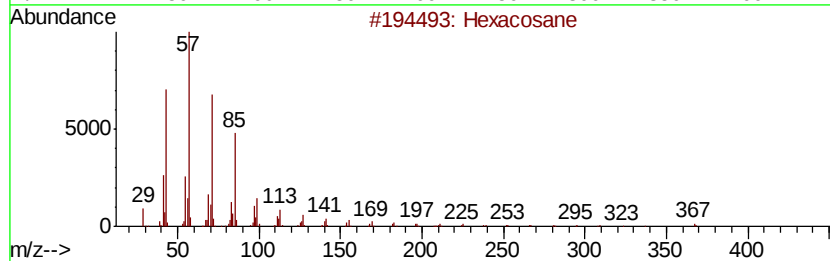
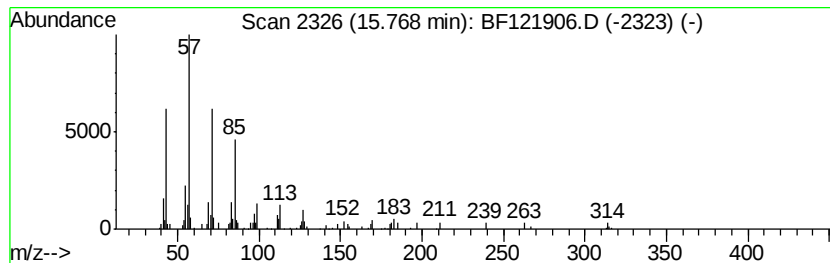
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	2.88 ng	107213	Perylene-d12	15.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	90
2	Docosane		310	C22H46	000629-97-0	87
3	Hentriacontane		437	C31H64	000630-04-6	87
4	Tetracosane		338	C24H50	000646-31-1	87
5	triacontane		422	C30H62	000638-68-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100920\
Data File : BF121906.D
Acq On : 9 Oct 2020 15:09
Operator : JU/CG
Sample : L4329-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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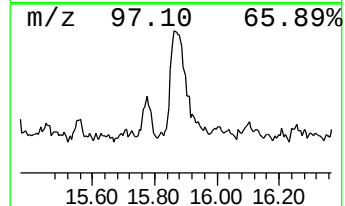
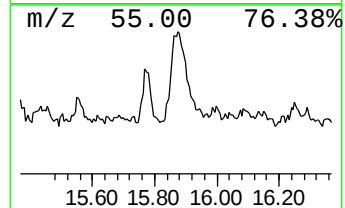
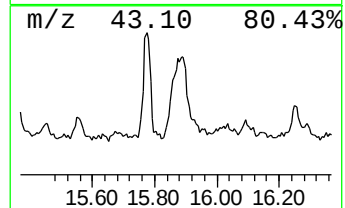
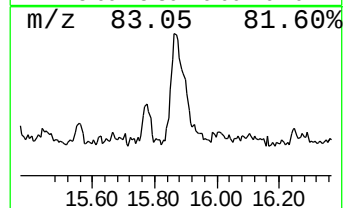
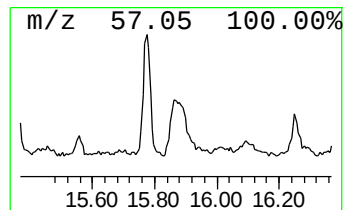
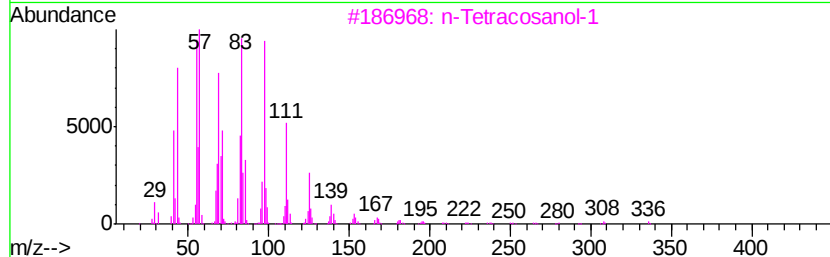
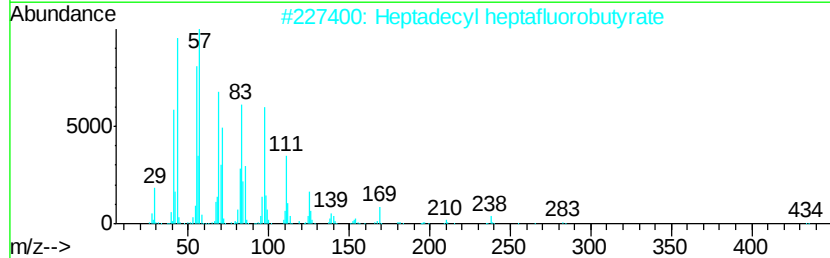
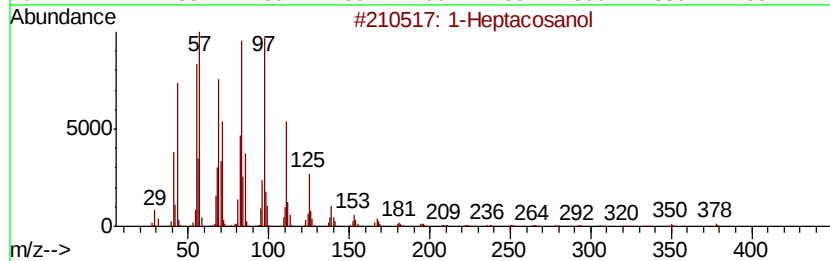
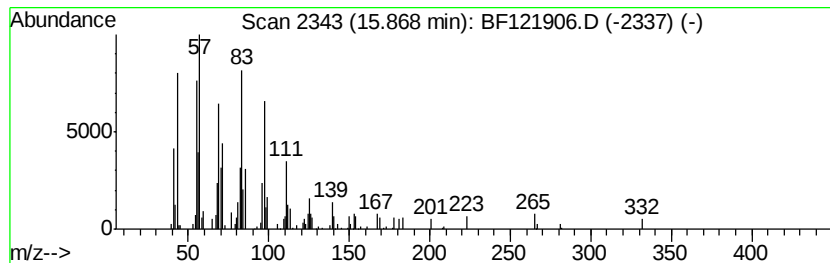
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Heptacosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	3.39 ng	126154	Perylene-d12	15.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol		396	C27H56O	002004-39-9	90
2	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	90
3	n-Tetracosanol-1		354	C24H50O	000506-51-4	90
4	Behenic alcohol		326	C22H46O	000661-19-8	87
5	Carbonic acid, hexadecyl 2,2,2-t...		416	C19H35Cl3O3	1000314-56-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100920\
Data File : BF121906.D
Acq On : 9 Oct 2020 15:09
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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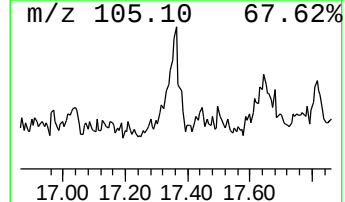
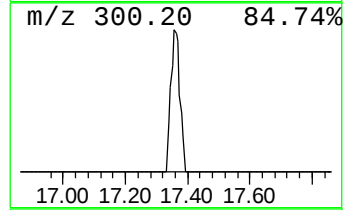
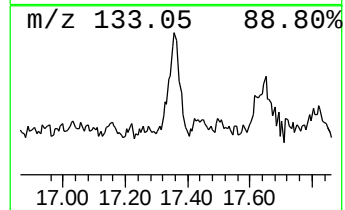
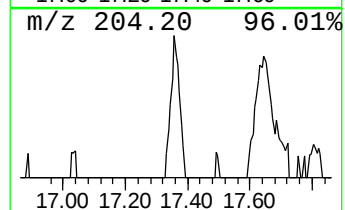
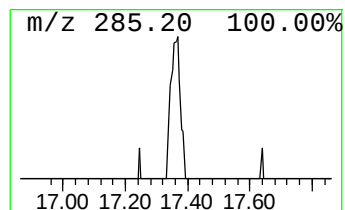
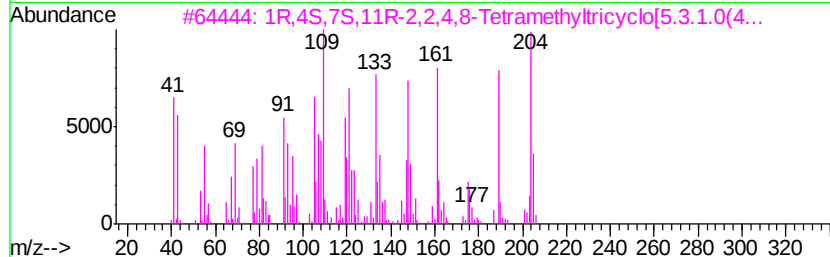
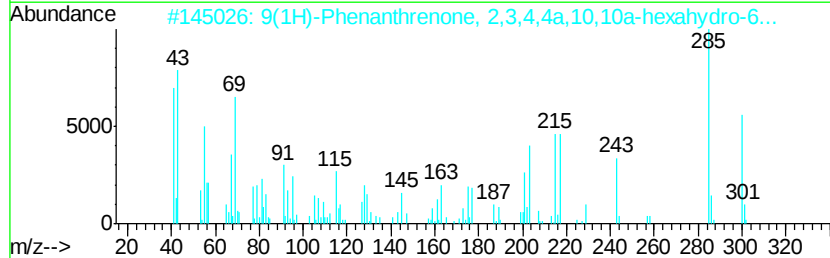
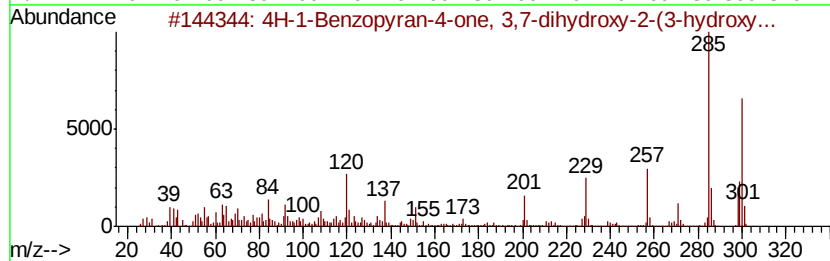
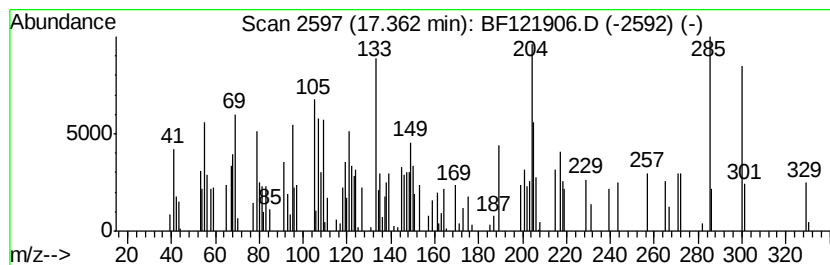
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown17.36 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.36	3.51 ng	130894	Perylene-d12	15.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-1-Benzopyran-4-one, 3,7-dihyd...	300	C16H12O6	057396-72-2	42
2			9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	42
3			1R,4S,7S,11R-2,2,4,8-Tetramethyl...	204	C15H24	1000140-07-6	27
4			4-(2,4,4-Trimethyl-bicyclo[4.1.0...	206	C14H22O	1000194-96-6	25
5			1,3-Benzodioxole, 5-(1-(4-ethoxy...	300	C18H20O4	090632-70-5	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100920\
Data File : BF121906.D
Acq On : 9 Oct 2020 15:09
Operator : JU/CG
Sample : L4329-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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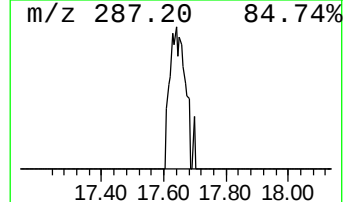
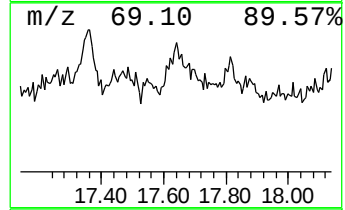
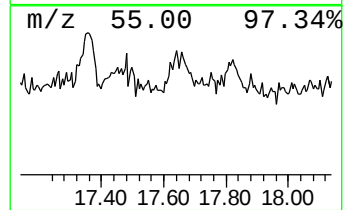
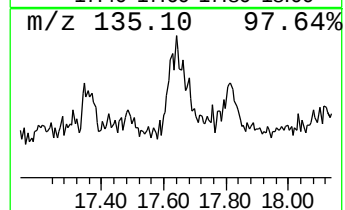
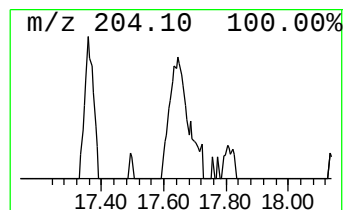
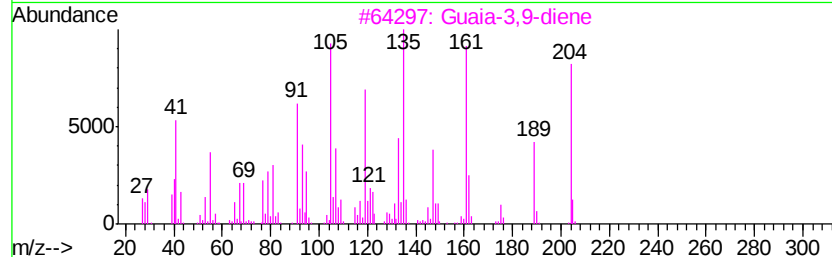
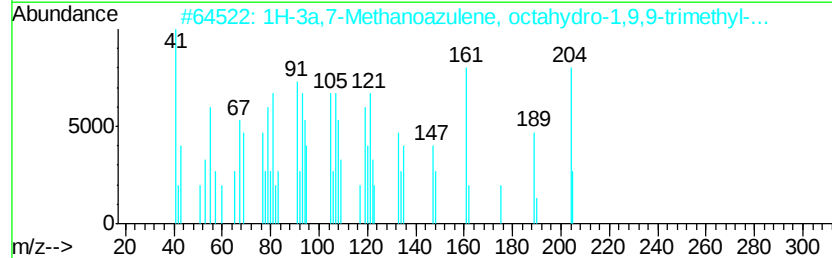
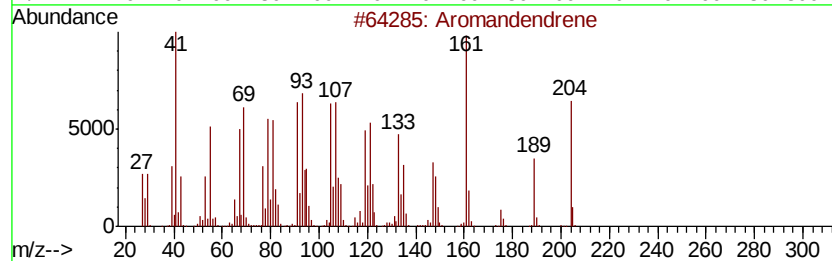
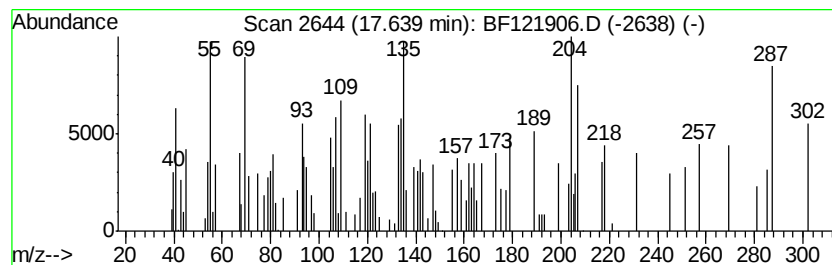
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.64 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.64	3.41 ng	127043	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Aromandendrene	204	C15H24	000489-39-4	40
2		1H-3a,7-Methanoazulene, octahydr...	204	C15H24	000508-55-4	40
3		Guaia-3,9-diene	204	C15H24	000489-83-8	35
4		2,5,5,8a-Tetramethyl-6,7,8,8a-te...	204	C14H20O	124957-09-1	20
5		1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF100920\
 Data File : BF121906.D
 Acq On : 9 Oct 2020 15:09
 Operator : JU/CG
 Sample : L4329-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF091020.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-hy...	4.96	3.0	ng	88050	1	6.76	596611	20.0
Ethanol, 2-(2-eth...	6.60	2.4	ng	71572	1	6.76	596611	20.0
1-Heneicosanol	13.77	14.7	ng	499325	5	13.92	678294	20.0
Trifluoroacetoxy ...	14.41	6.2	ng	209423	5	13.92	678294	20.0
unknown14.67	14.67	2.9	ng	106982	6	15.36	744918	20.0
Triacontane	15.00	3.6	ng	133931	6	15.36	744918	20.0
Pentafluoropropio...	15.07	2.9	ng	107588	6	15.36	744918	20.0
Hexacosane	15.77	2.9	ng	107213	6	15.36	744918	20.0
1-Heptacosanol	15.87	3.4	ng	126154	6	15.36	744918	20.0
unknown17.36	17.36	3.5	ng	130894	6	15.36	744918	20.0
unknown17.64	17.64	3.4	ng	127043	6	15.36	744918	20.0