

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062320\
 Data File : BP002493.D
 Acq On : 23 Jun 2020 09:24
 Operator : CG/JU
 Sample : L3082-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-2

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_P\METHODS\8270-BP060520.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	3.040	22	26	39	rVB	77537	118240	1.86%	0.321%
2	5.205	385	394	406	rBV	1497076	2172918	34.20%	5.900%
3	6.775	653	661	676	rBV	1481447	2331189	36.69%	6.330%
4	7.587	791	799	809	rBV	434081	682854	10.75%	1.854%
5	7.905	846	853	863	rBV	1301843	2043730	32.17%	5.549%
6	8.728	985	993	1006	rBV	997143	1637275	25.77%	4.446%
7	10.357	1261	1270	1281	rBV	582037	955081	15.03%	2.593%
8	12.851	1687	1694	1712	rBV	2413028	3830100	60.28%	10.400%
9	13.151	1739	1745	1752	rBV	178632	280509	4.41%	0.762%
10	13.692	1832	1837	1846	rBV	145732	224380	3.53%	0.609%
11	13.922	1864	1876	1877	rBV	38655	99535	1.57%	0.270%
12	14.234	1924	1929	1935	rVB	901862	1389333	21.87%	3.772%
13	14.687	2003	2006	2014	rVB	42557	64801	1.02%	0.176%
14	15.734	2174	2184	2195	rBV	2025349	3122075	49.14%	8.477%
15	16.398	2292	2297	2300	rBV3	54048	95941	1.51%	0.261%
16	16.992	2392	2398	2406	rBV	1177380	1686646	26.55%	4.580%
17	17.922	2551	2556	2565	rBV	505824	770353	12.12%	2.092%
18	18.175	2594	2599	2603	rBV	126895	198454	3.12%	0.539%
19	18.216	2603	2606	2626	rVB	199127	387454	6.10%	1.052%
20	19.163	2764	2767	2774	rBV2	64075	104572	1.65%	0.284%
21	19.580	2833	2838	2849	rVB	5258435	6353614	100.00%	17.252%
22	19.798	2870	2875	2882	rBV2	66622	137398	2.16%	0.373%
23	20.645	3013	3019	3025	rBV	100672	221354	3.48%	0.601%
24	20.845	3048	3053	3057	rBV	498841	735138	11.57%	1.996%
25	20.892	3057	3061	3068	rVB	589993	755272	11.89%	2.051%
26	21.098	3091	3096	3108	rBV	1717001	2169875	34.15%	5.892%
27	21.345	3133	3138	3144	rVV	126575	245362	3.86%	0.666%
28	21.710	3197	3200	3208	rVB3	99805	142873	2.25%	0.388%
29	22.133	3268	3272	3276	rBV2	142524	266920	4.20%	0.725%
30	22.292	3295	3299	3314	rVB	211093	475016	7.48%	1.290%
31	22.674	3360	3364	3367	rBV	172076	286193	4.50%	0.777%
32	23.239	3456	3460	3464	rBV	122709	194674	3.06%	0.529%
33	23.298	3464	3470	3482	rVB	1296775	2450438	38.57%	6.654%
34	23.886	3567	3570	3578	rBV5	82181	199111	3.13%	0.541%

LSC Area Percent Report

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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_P\METHODS\8270-BP060520.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

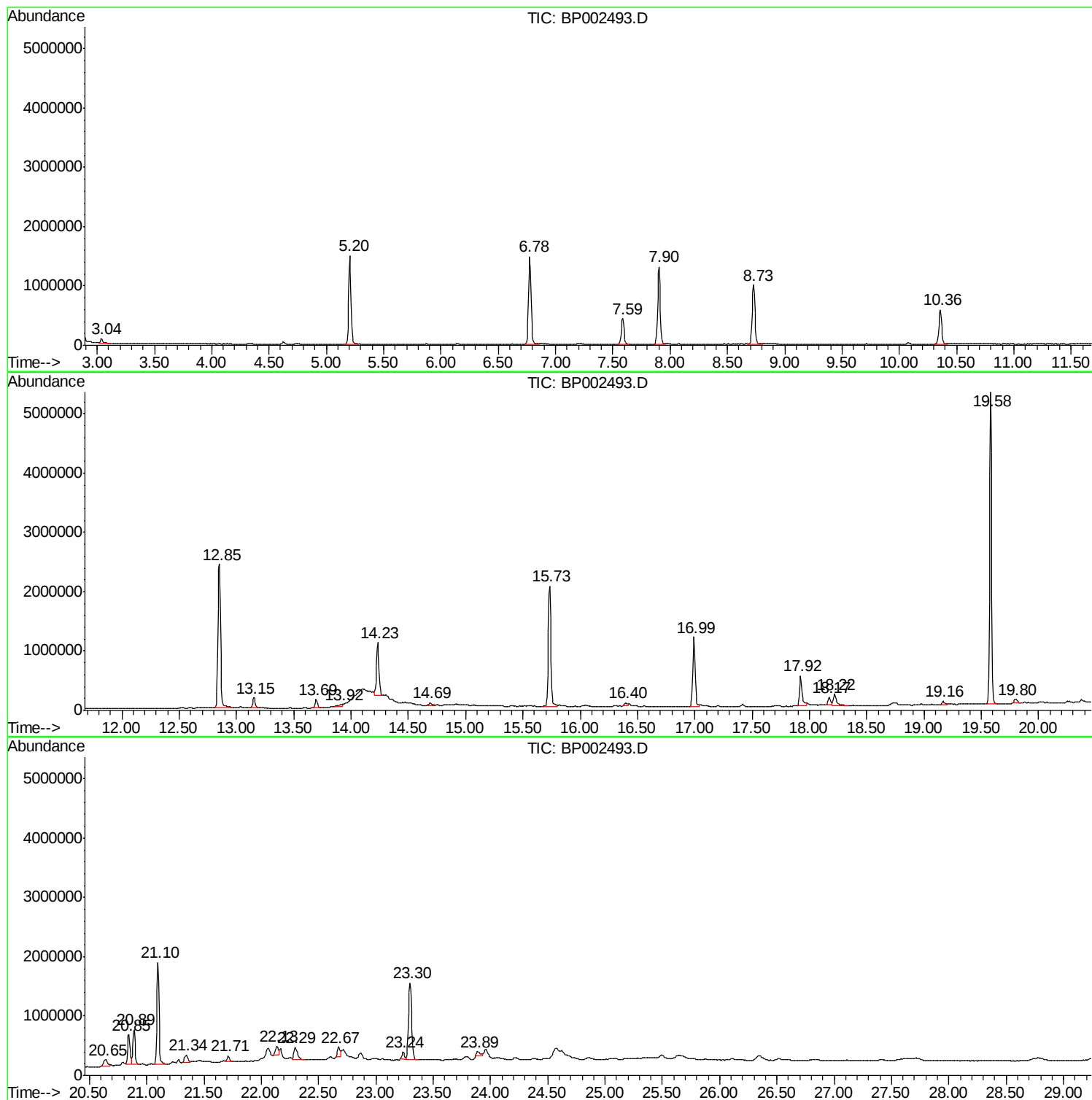
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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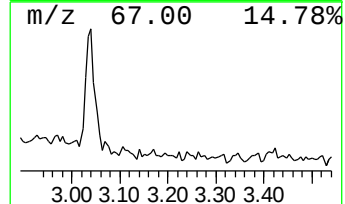
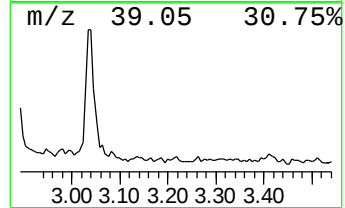
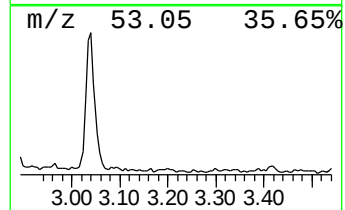
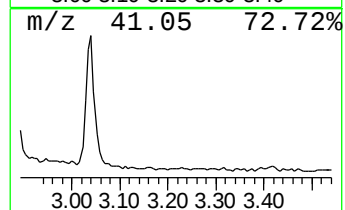
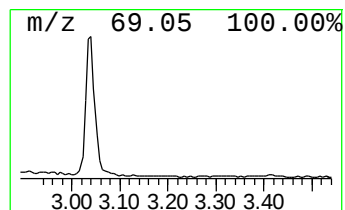
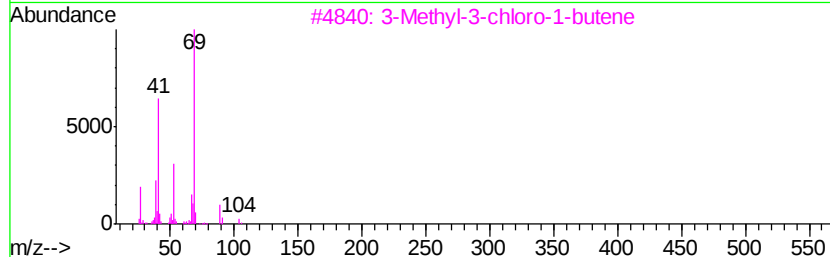
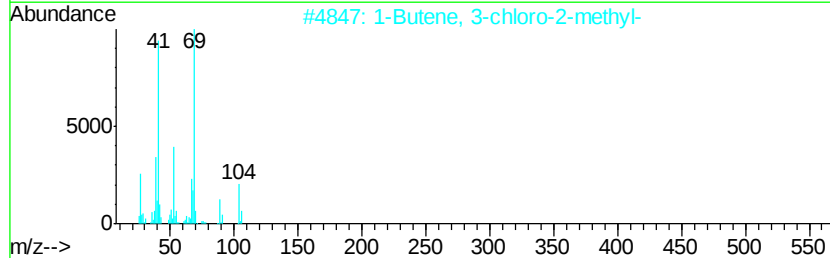
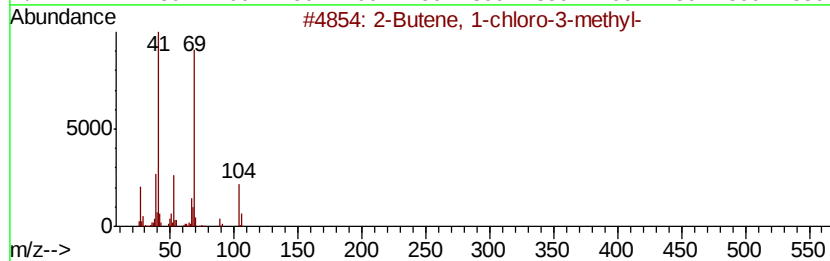
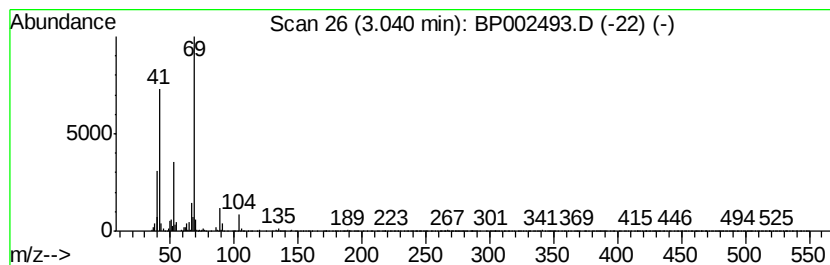
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Butene, 1-chloro-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.04	3.46 ng	118240	1,4-Dichlorobenzene-d4	7.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butene, 1-chloro-3-methyl-	104	C5H9Cl	000503-60-6	72
2			1-Butene, 3-chloro-2-methyl-	104	C5H9Cl	005166-35-8	64
3			3-Methyl-3-chloro-1-butene	104	C5H9Cl	002190-48-9	64
4			2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	64
5			Cyclopentane, bromo-	148	C5H9Br	000137-43-9	53



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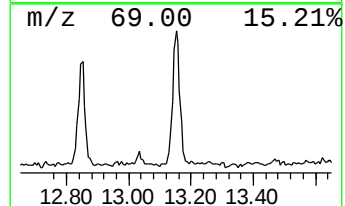
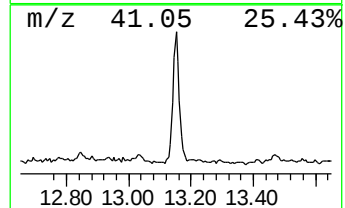
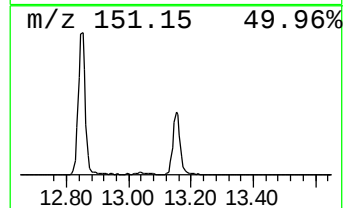
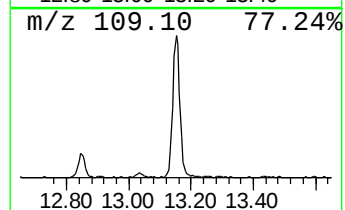
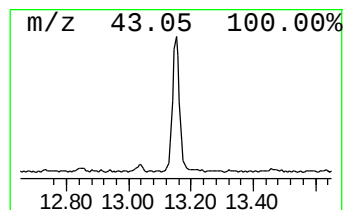
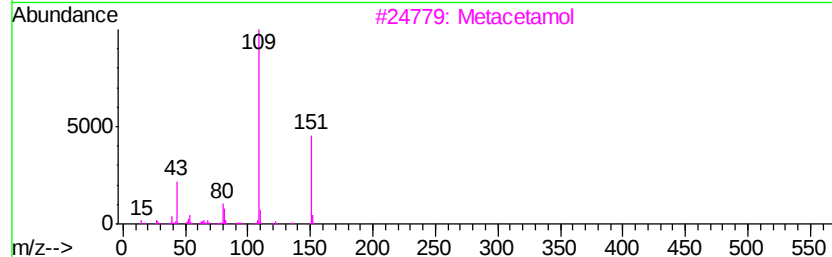
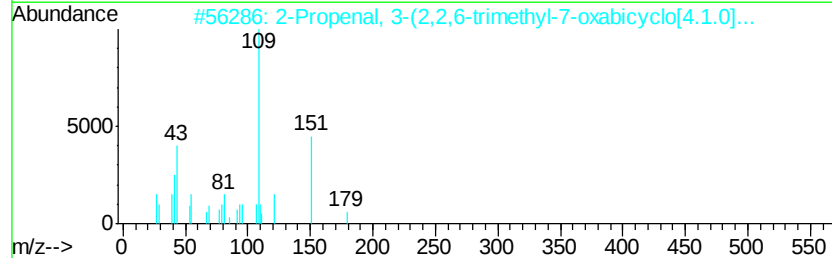
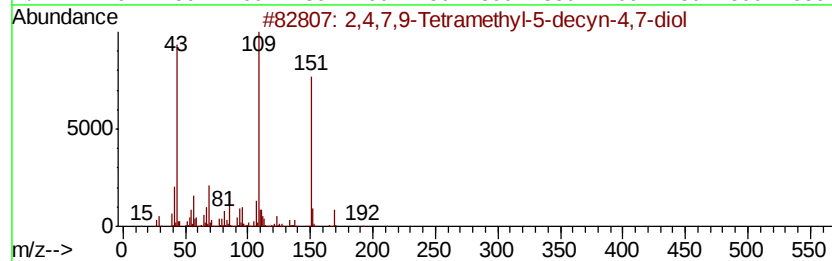
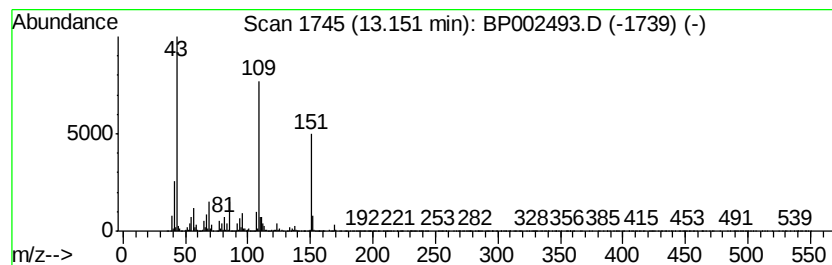
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	4.04 ng	280509	Acenaphthene-d10	14.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	53		
3	Metacetamol	151	C8H9NO2	000621-42-1	53		
4	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	43		
5	1,2-Diazaspiro[4.4]nonen-3-carbo...	238	C13H22N2O2	1000164-25-8	43		



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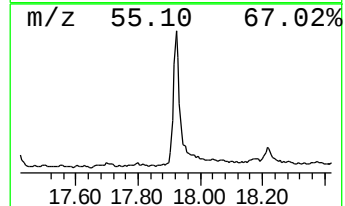
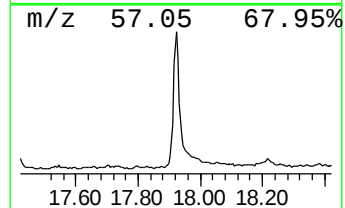
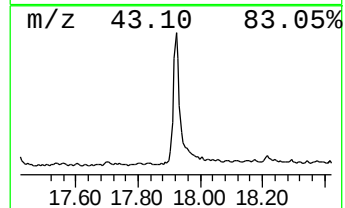
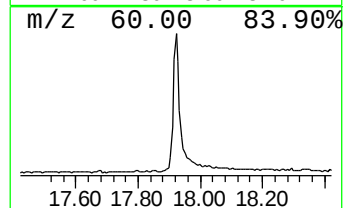
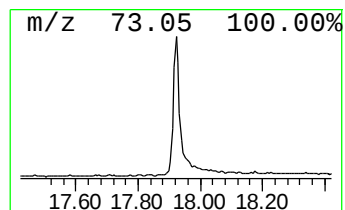
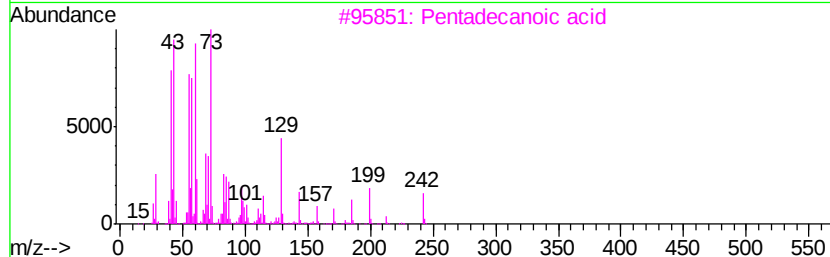
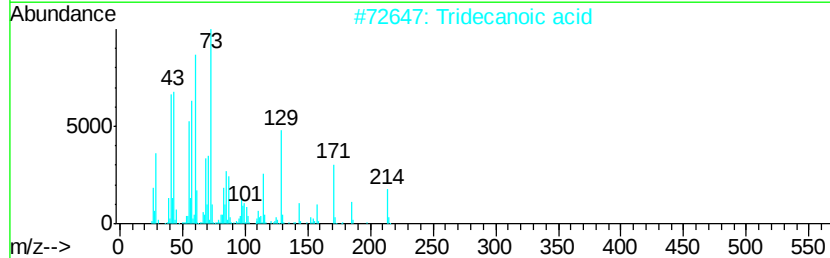
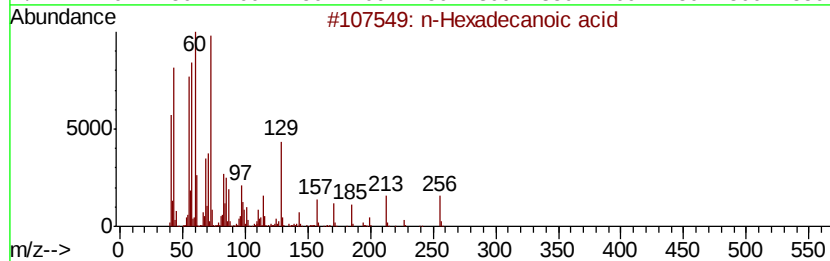
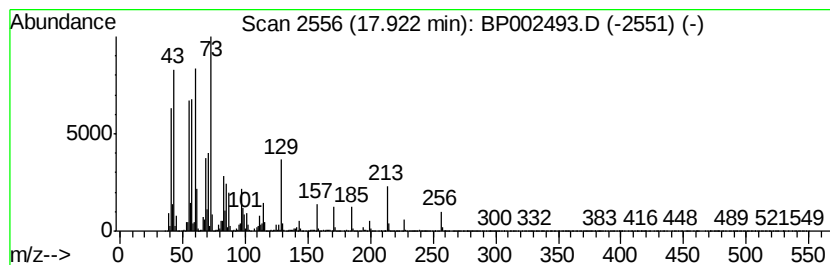
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	9.13 ng	770353	Phenanthrene-d10	16.99

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	95
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		Dodecanoic acid	200	C12H24O2	000143-07-7	59



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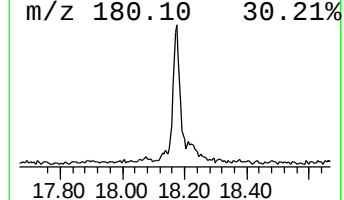
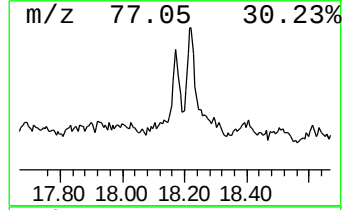
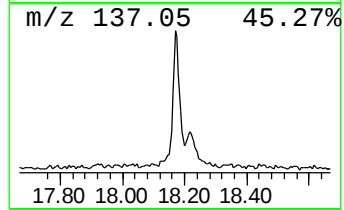
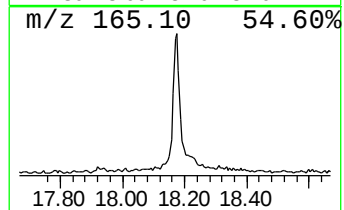
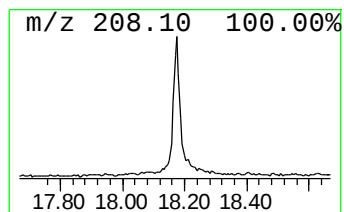
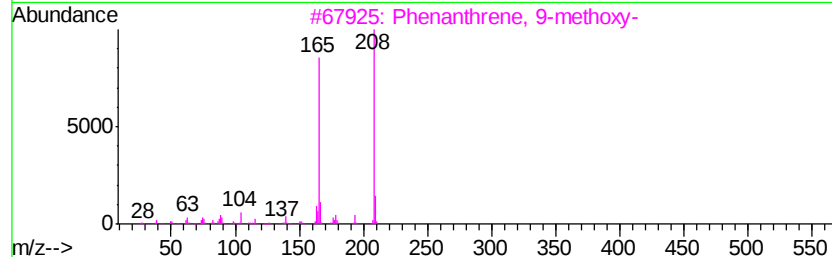
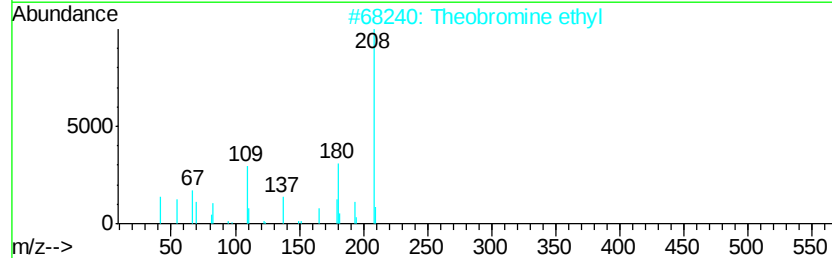
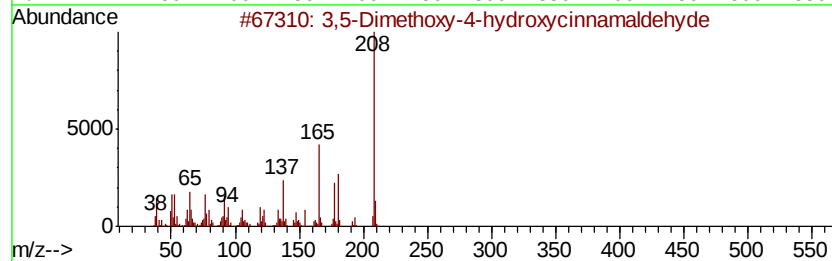
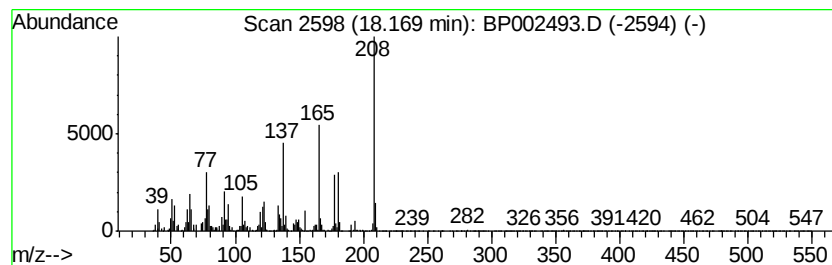
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	2.35 ng	198454	Phenanthrene-d10	16.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethoxy-4-hydroxycinnamaldehyde	208	C11H12O4	087345-53-7	95	
2	Theobromine ethyl	208	C9H12N4O2	007464-74-6	43	
3	Phenanthrene, 9-methoxy-	208	C15H12O	005085-74-5	43	
4	2-Propenoic acid, 3-(2,3-dimethoxy-5-hydroxyphenyl)-	208	C11H12O4	007345-82-6	43	
5	trans-2,5-Dimethoxycinnamic acid	208	C11H12O4	038489-74-6	43	



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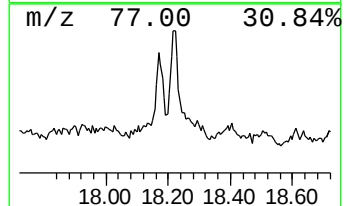
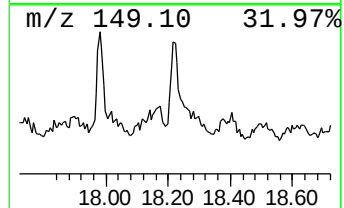
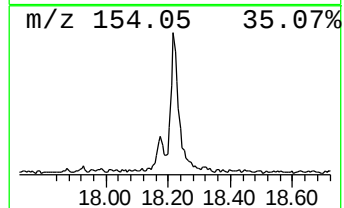
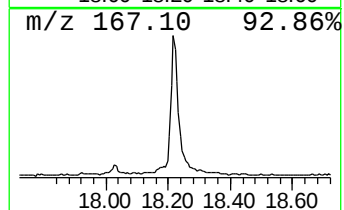
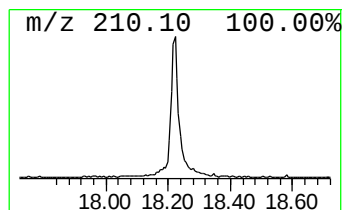
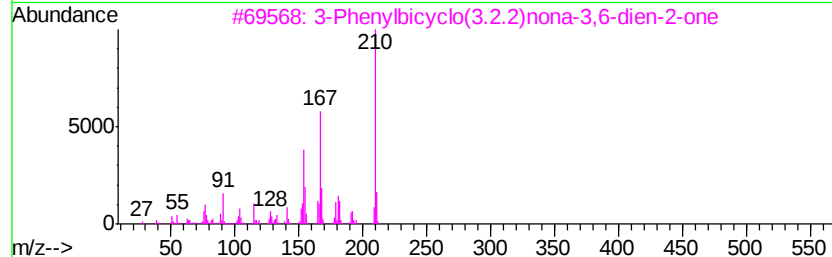
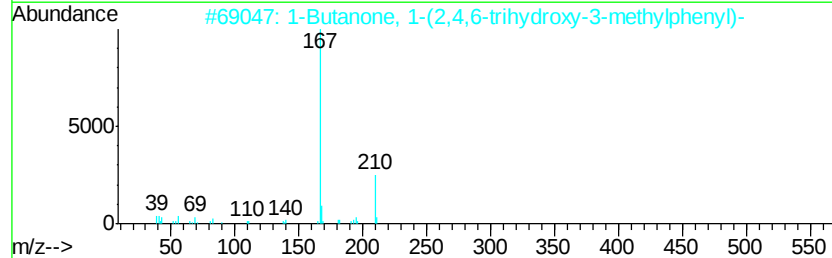
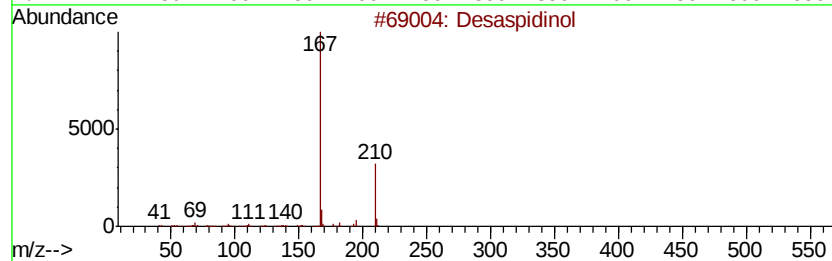
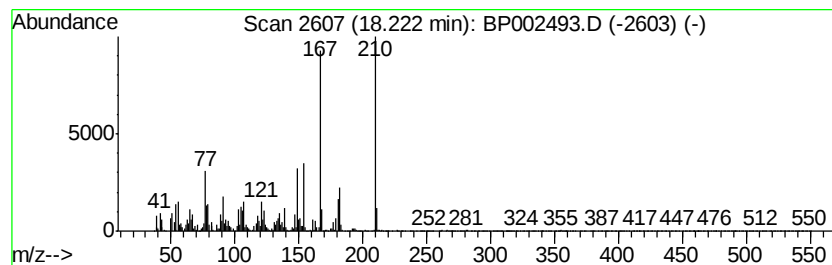
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown18.22 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.22	4.59 ng	387454	Phenanthrene-d10	16.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Desaspidinol	210	C11H14O4	000437-72-9	47
2		1-Butanone, 1-(2,4,6-trihydroxyv-...	210	C11H14O4	001509-06-4	47
3		3-Phenylbicyclo(3.2.2)nona-3,6-d...	210	C15H14O	073830-83-8	47
4		2-Pentanone, 1-(2,4,6-trihydroxyv...	210	C11H14O4	1000116-22-3	47
5		9-Methyl-1-phenazinyll acetate	252	C15H12N2O2	013860-44-1	42



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Data File : BP002493.D
Acq On : 23 Jun 2020 09:24
Operator : CG/JU
Sample : L3082-02
Misc :
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ClientSampleId :
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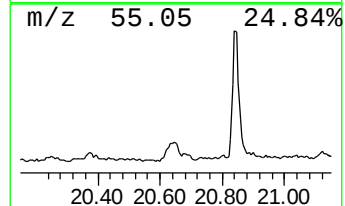
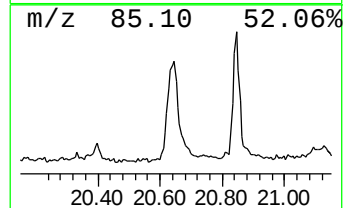
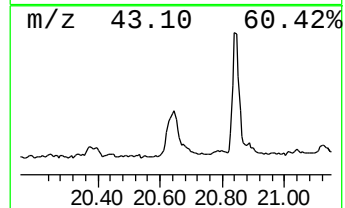
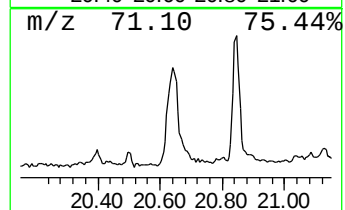
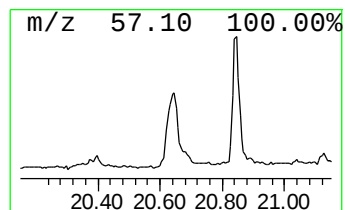
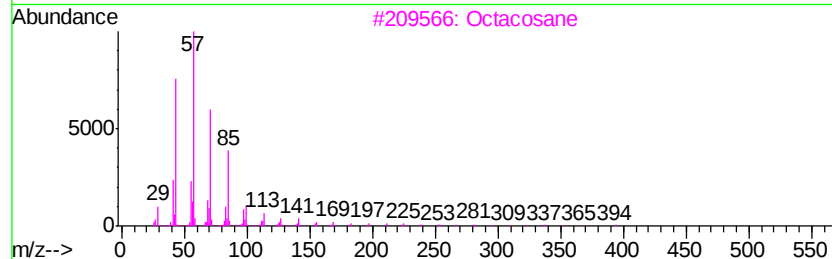
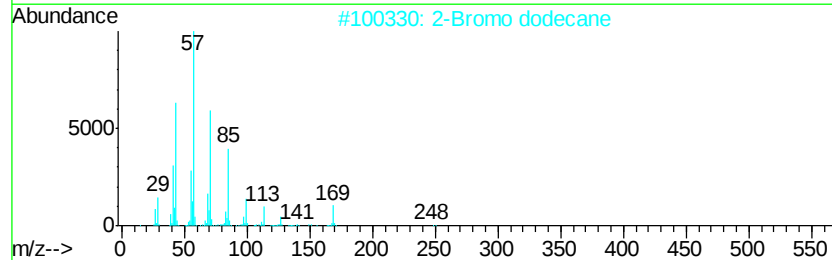
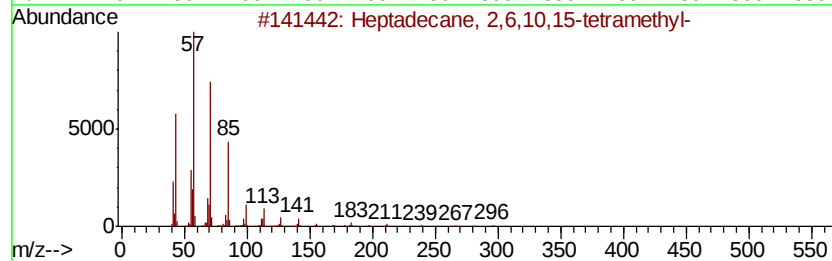
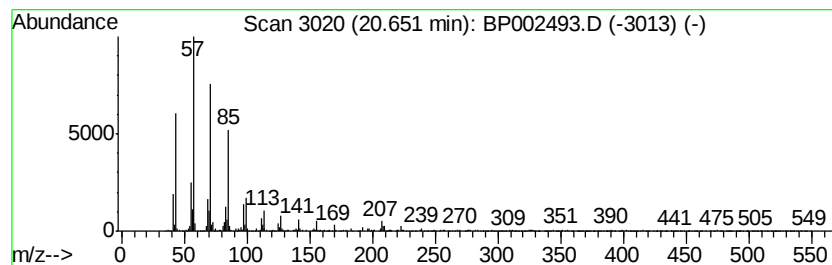
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane, 2,6,10,15-tetr... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.65	2.04 ng	221354	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Hexacosane	366	C26H54	000630-01-3	90
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062320\
Data File : BP002493.D
Acq On : 23 Jun 2020 09:24
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ClientSampleId :
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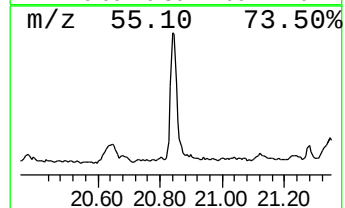
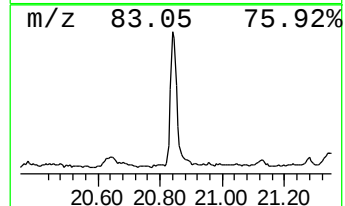
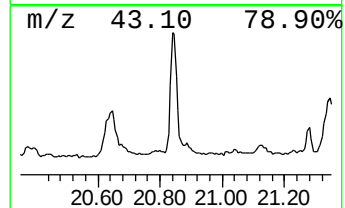
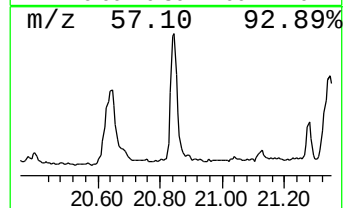
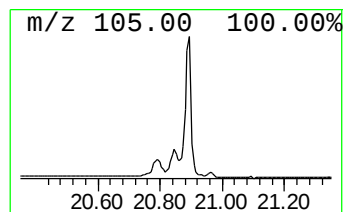
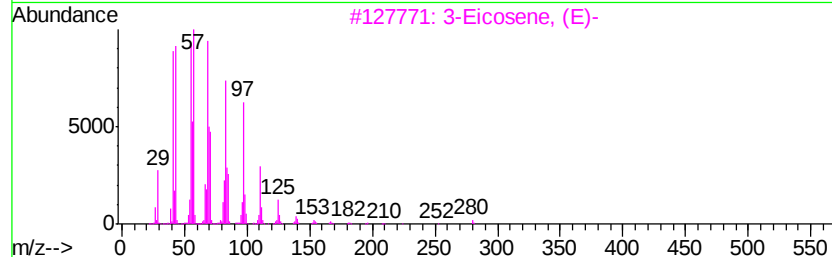
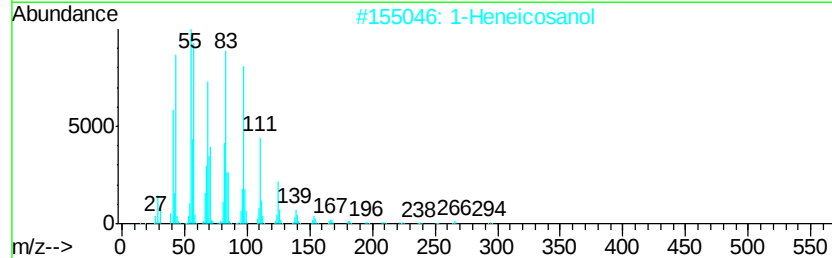
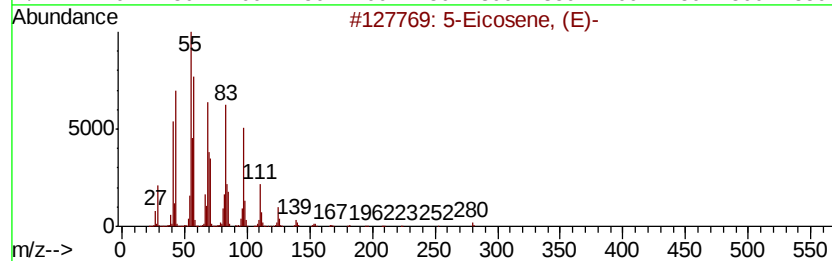
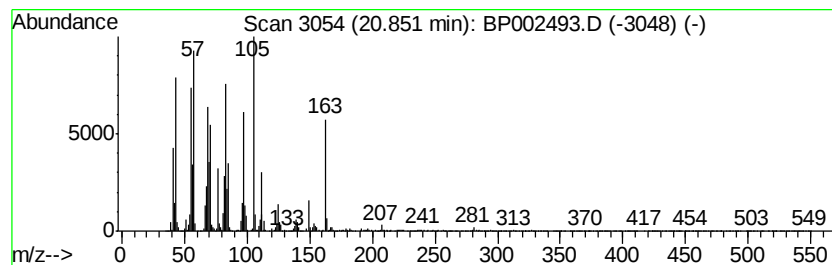
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	6.78 ng	735138	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2		1-Heneicosanol	312	C21H44O	015594-90-8	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	90
4		Trifluoroacetoxv hexadecane	338	C18H33F3O2	006222-03-3	90
5		Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062320\
Data File : BP002493.D
Acq On : 23 Jun 2020 09:24
Operator : CG/JU
Sample : L3082-02
Misc :
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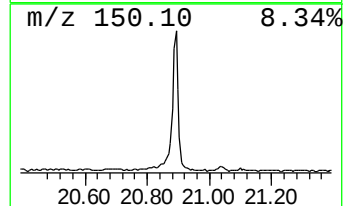
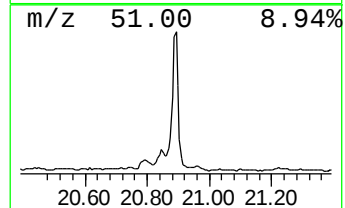
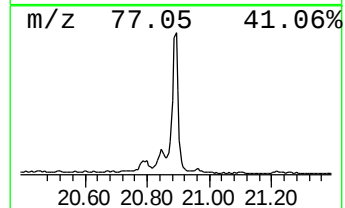
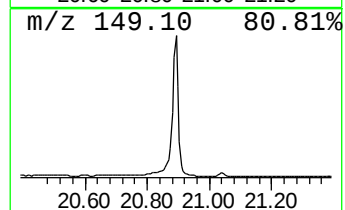
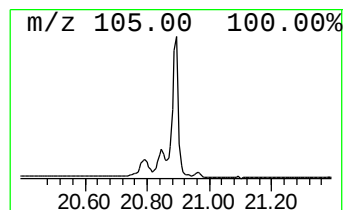
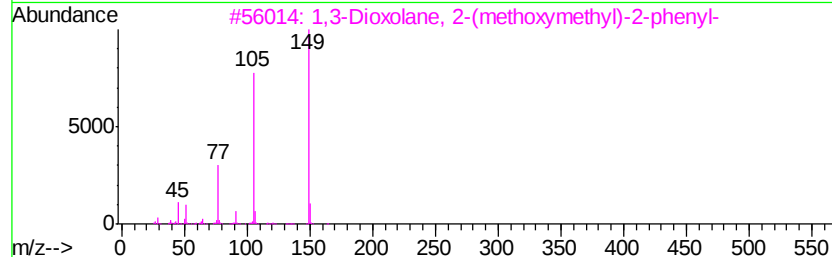
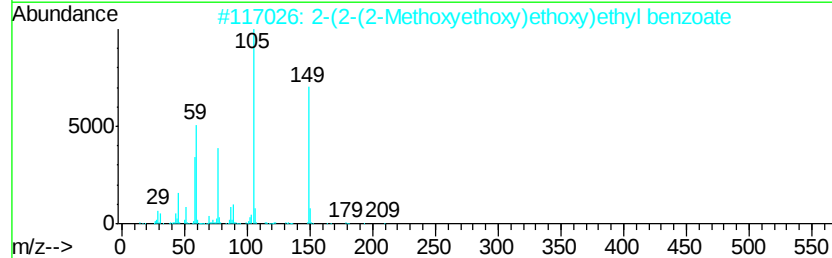
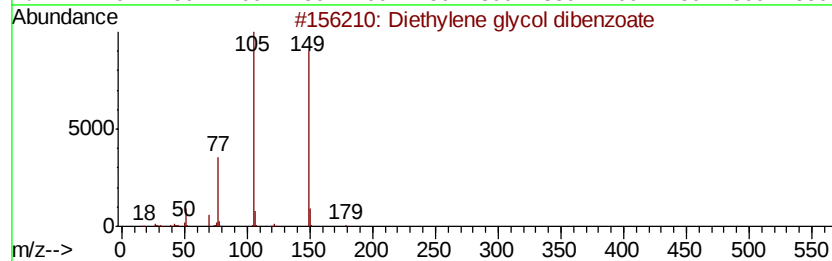
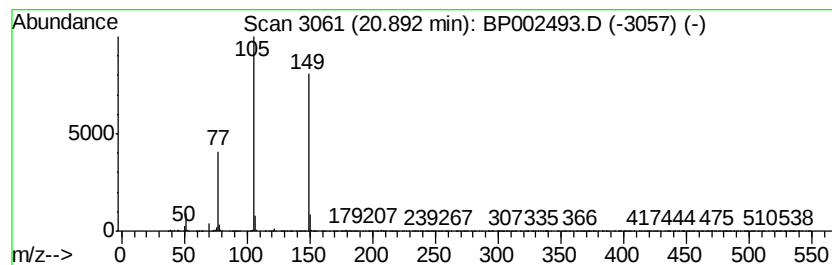
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Diethylene glycol dibenzoate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	6.96 ng	755272	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2		2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74
3		1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	74
4		2,2'-(Ethane-1,2-diylbis(oxv))bi...	358	C20H22O6	1000366-99-4	74
5		Benzoic acid, 2-(3-nitrophenyl)e...	271	C15H13NO4	1000368-22-4	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062320\
Data File : BP002493.D
Acq On : 23 Jun 2020 09:24
Operator : CG/JU
Sample : L3082-02
Misc :
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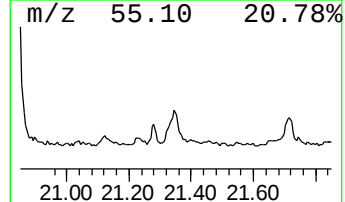
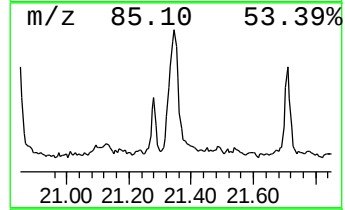
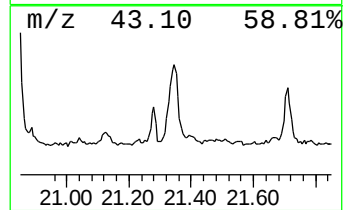
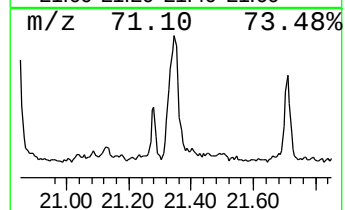
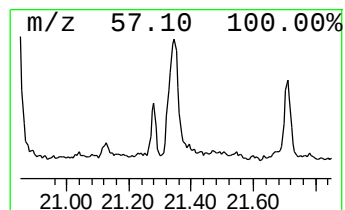
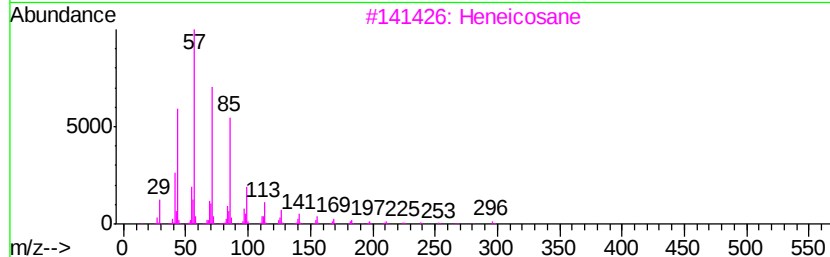
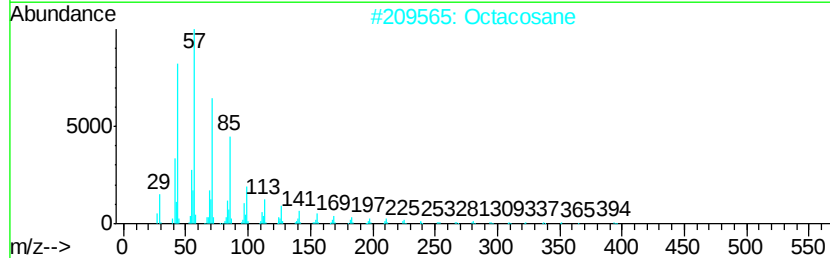
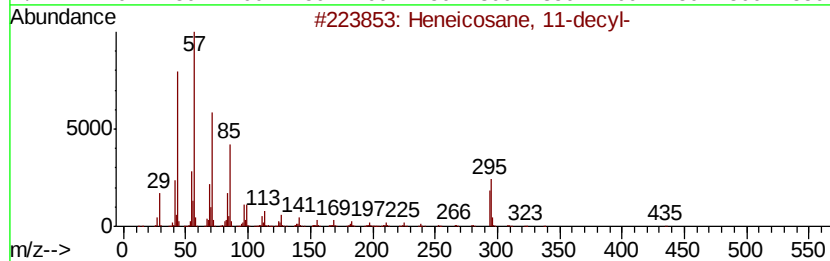
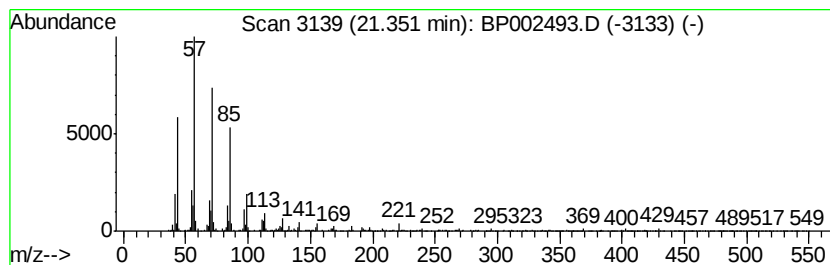
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane, 11-decyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	2.26 ng	245362	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
2		Octacosane	394	C28H58	000630-02-4	89
3		Heneicosane	296	C21H44	000629-94-7	87
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062320\
Data File : BP002493.D
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Sample : L3082-02
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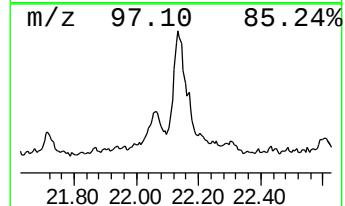
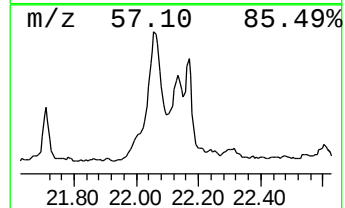
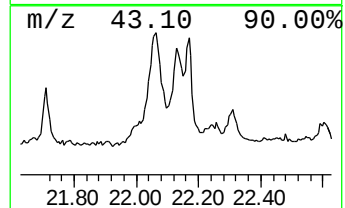
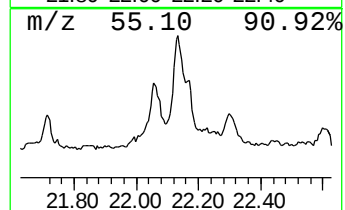
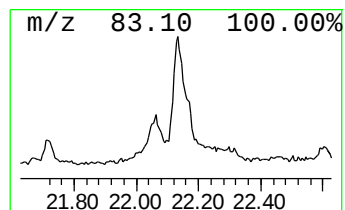
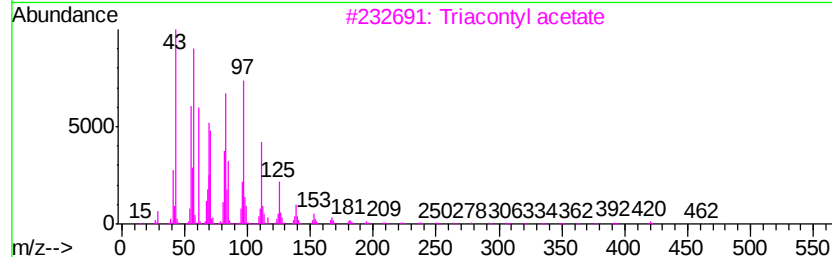
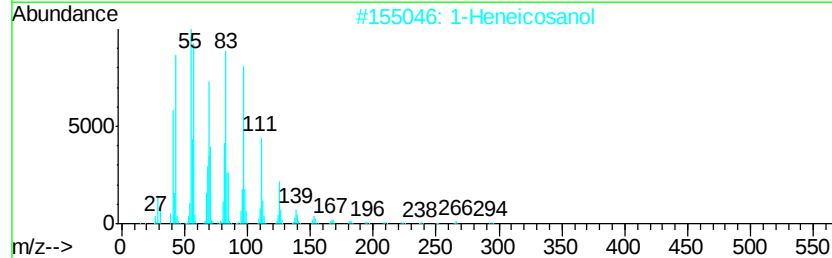
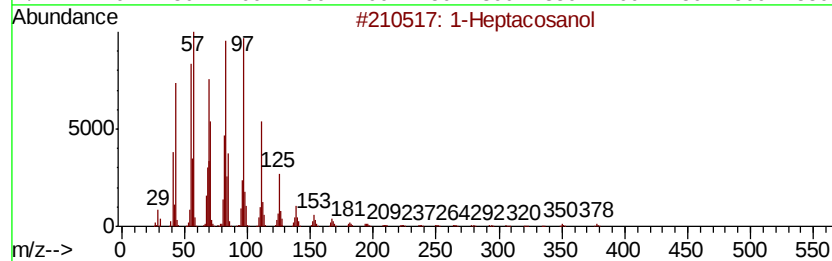
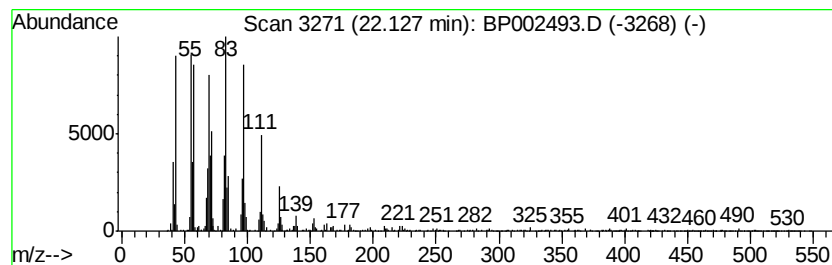
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Heptacosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.13	2.46 ng	266920	Chrysene-d12		21.10	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heptacosanol	396	C27H56O	002004-39-9	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	94
3		Triaccontyl acetate	480	C32H64O2	041755-58-2	93
4		17-Pentatriacontene	491	C35H70	006971-40-0	93
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	91



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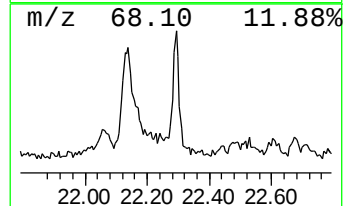
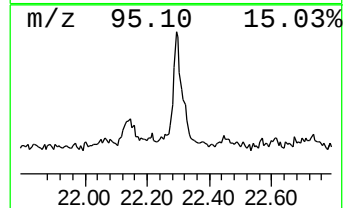
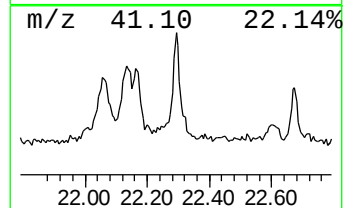
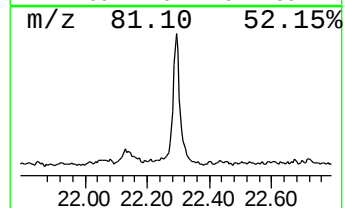
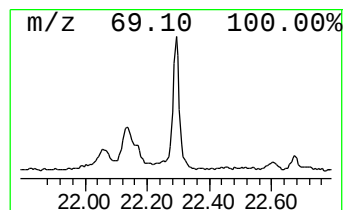
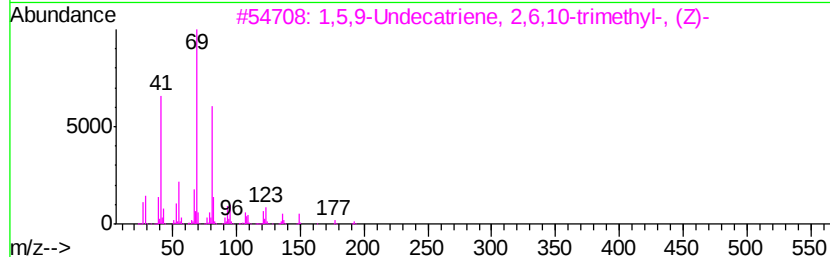
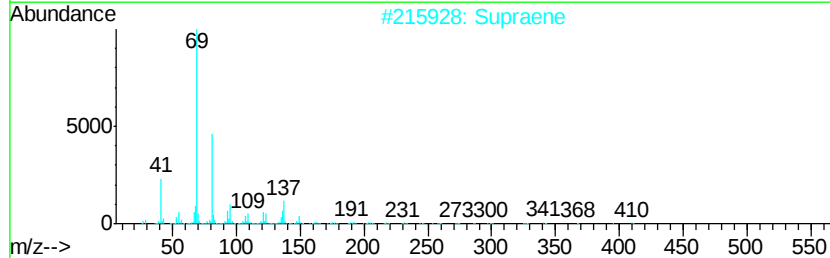
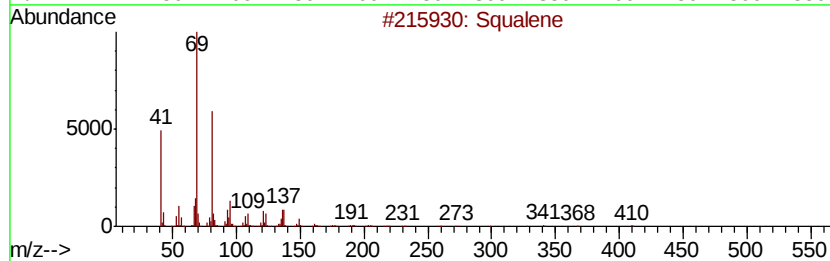
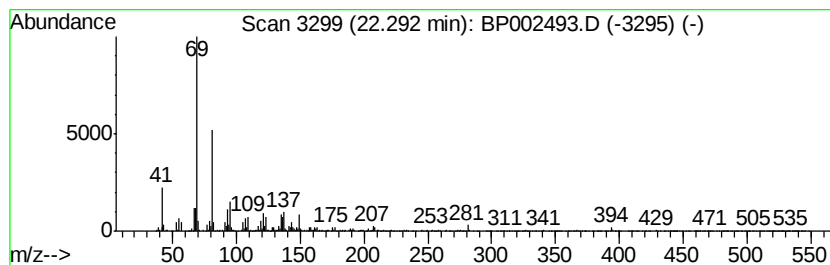
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.29	3.88 ng	475016	Perylene-d12	23.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene	410 C30H50	000111-02-4	83
2	Supraene	410 C30H50	007683-64-9	83
3	1,5,9-Undecatriene, 2,6,10-trime...	192 C14H24	062951-96-6	74
4	.psi.,.psi.-Carotene, 7,7',8,8',...	547 C40H66	000502-62-5	72
5	3,7,11-Tridecatrienoic acid, 4,8...	264 C17H28O2	036237-69-1	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062320\
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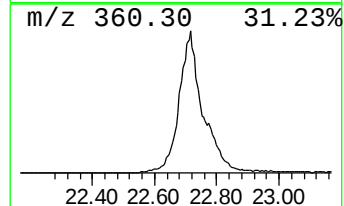
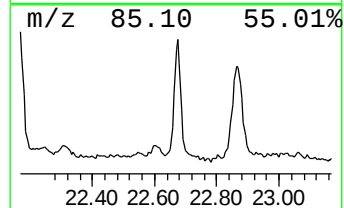
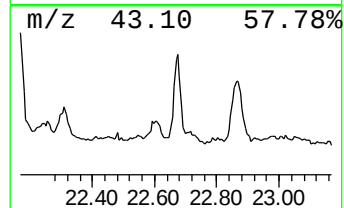
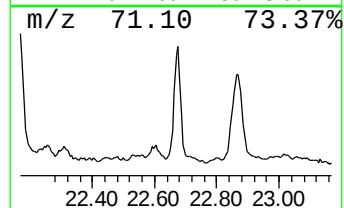
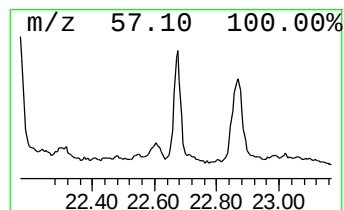
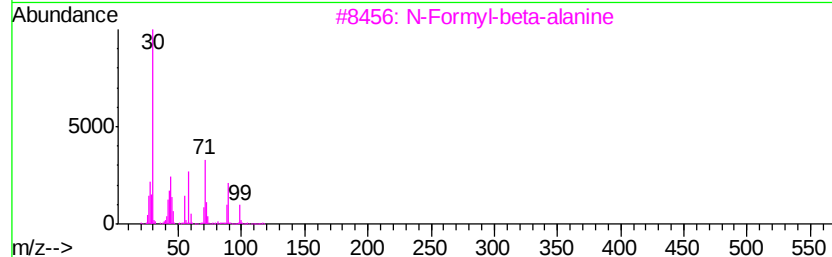
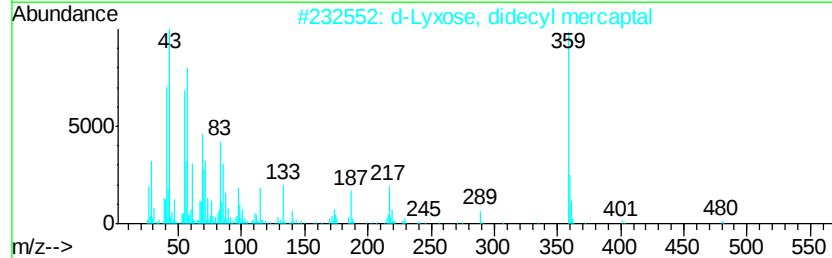
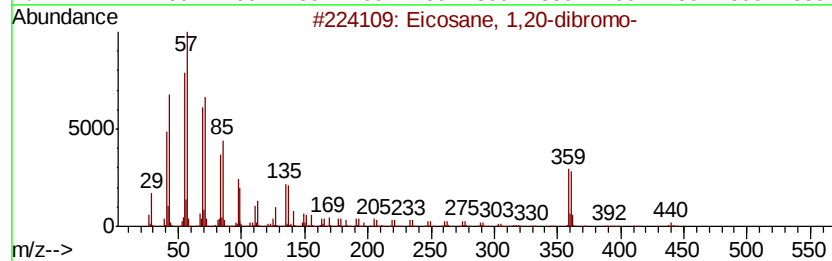
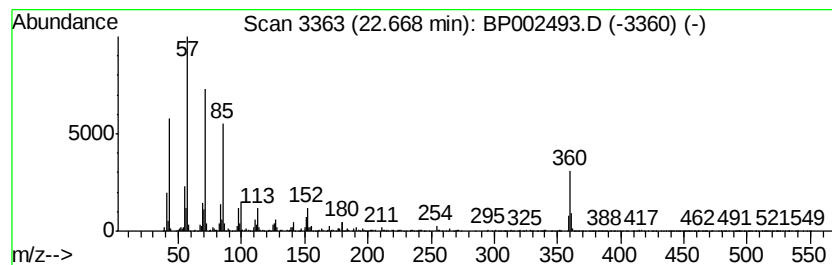
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown22.67 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.67	2.34 ng	286193	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 1,20-dibromo-	438	C20H40Br2	014296-16-3	9
2		d-Lyxose, didecyl mercaptal	480	C25H52O4S2	112159-64-5	7
3		N-Formyl-beta-alanine	117	C4H7NO3	014565-43-6	5
4		3.beta.-Acetoxy-17.alpha.-methyl...	360	C23H36O3	065731-78-4	4
5		Acetic acid, 1,2-diacetoxy-4,10,...	462	C26H38O7	1000192-72-1	4



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP062320\
Data File : BP002493.D
Acq On : 23 Jun 2020 09:24
Operator : CG/JU
Sample : L3082-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP060520.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-Butene, 1-chlor...	3.04	3.5 ng		118240	1	7.59	682854	20.0
2,4,7,9-Tetrameth...	13.15	4.0 ng		280509	3	14.23	1389330	20.0
n-Hexadecanoic acid	17.92	9.1 ng		770353	4	16.99	1686650	20.0
3,5-Dimethoxy-4-h...	18.17	2.4 ng		198454	4	16.99	1686650	20.0
unknown18.22	18.22	4.6 ng		387454	4	16.99	1686650	20.0
Heptadecane, 2,6,...	20.65	2.0 ng		221354	5	21.10	2169880	20.0
5-Eicosene, (E)-	20.85	6.8 ng		735138	5	21.10	2169880	20.0
Diethylene glycol...	20.89	7.0 ng		755272	5	21.10	2169880	20.0
Heneicosane, 11-d...	21.35	2.3 ng		245362	5	21.10	2169880	20.0
1-Heptacosanol	22.13	2.5 ng		266920	5	21.10	2169880	20.0
Squalene	22.29	3.9 ng		475016	6	23.30	2450440	20.0
unknown22.67	22.67	2.3 ng		286193	6	23.30	2450440	20.0