

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
 Data File : BP001362.D
 Acq On : 23 Dec 2019 12:40
 Operator : JU
 Sample : PB125644BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125644BL

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-BP121919.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	5.628	596	602	617	rBV	234692	397537	12.92%	1.773%
2	6.222	696	703	707	rBV	1588075	2051510	66.65%	9.149%
3	7.763	957	965	968	rBV	1611412	2163953	70.31%	9.650%
4	7.946	989	996	1000	rBV	1759889	2213215	71.91%	9.870%
5	8.298	1051	1056	1063	rVB	348439	413138	13.42%	1.842%
6	8.545	1091	1098	1102	rBV	1465069	1737355	56.45%	7.748%
7	9.187	1199	1207	1211	rBV	1097312	1409127	45.78%	6.284%
8	10.334	1395	1402	1413	rBV	415089	489941	15.92%	2.185%
9	12.122	1698	1706	1722	rBV	2001049	2575220	83.67%	11.484%
10	13.198	1883	1889	1899	rBV	454213	582703	18.93%	2.599%
11	14.498	2102	2110	2131	rBV	1431107	1920971	62.41%	8.566%
12	15.598	2287	2297	2308	rBV	508251	634698	20.62%	2.830%
13	17.892	2680	2687	2691	rBV	2819913	3077939	100.00%	13.726%
14	19.251	2912	2918	2929	rBV	511133	599036	19.46%	2.671%
15	19.539	2962	2967	2973	rBV	72263	84658	2.75%	0.378%
16	19.933	3029	3034	3041	rBV	179886	211182	6.86%	0.942%
17	20.310	3093	3098	3104	rBV	287812	355261	11.54%	1.584%
18	20.704	3159	3165	3174	rBV	254538	345149	11.21%	1.539%
19	21.027	3213	3220	3232	rVB	396013	576241	18.72%	2.570%
20	21.145	3233	3240	3251	rBV	168764	267460	8.69%	1.193%
21	21.639	3318	3324	3332	rBV	96789	170526	5.54%	0.760%
22	22.209	3415	3421	3429	rBV2	48925	98099	3.19%	0.437%
23	22.880	3529	3535	3545	rVB3	19676	49463	1.61%	0.221%

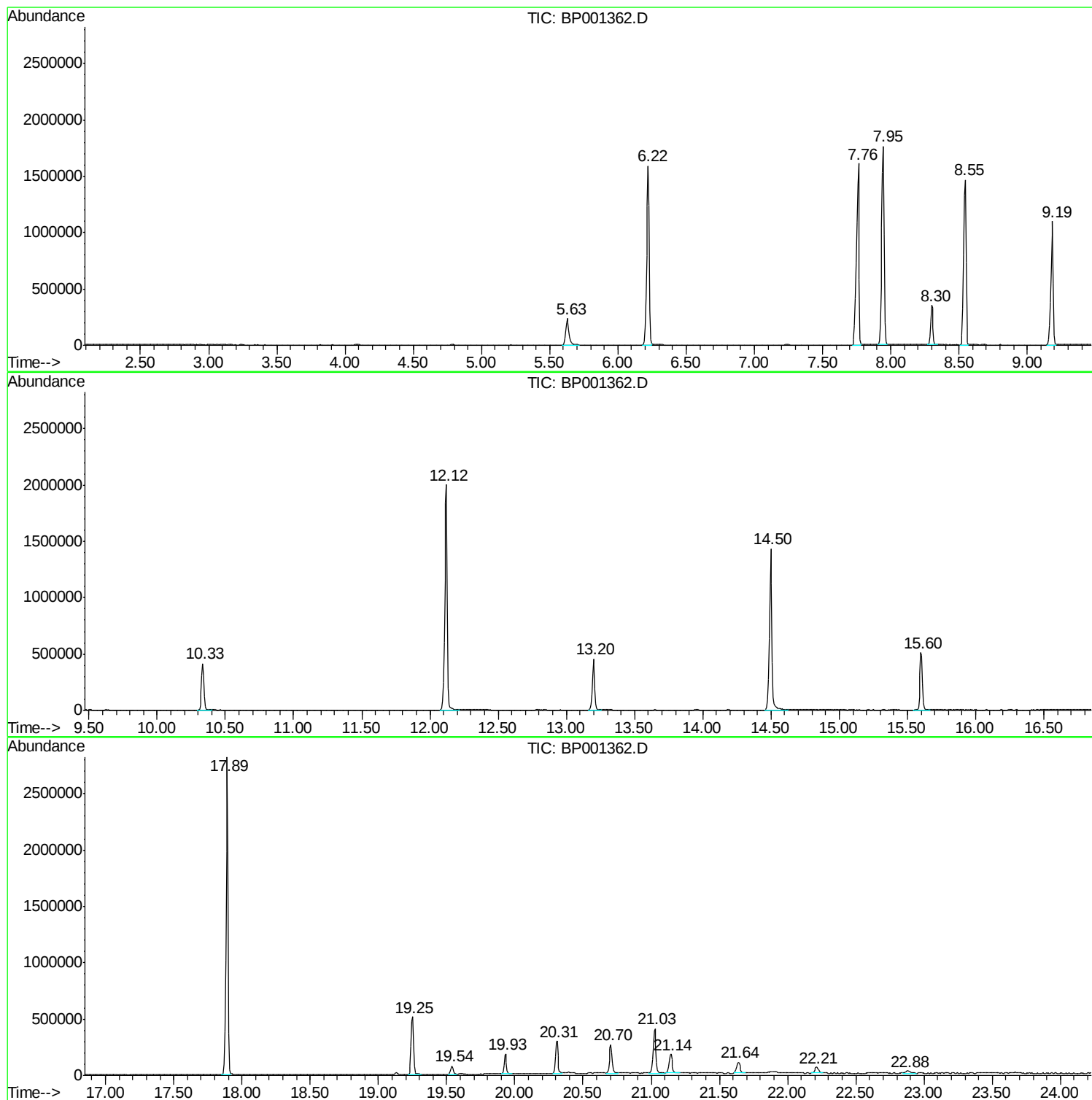
Sum of corrected areas: 22424382

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

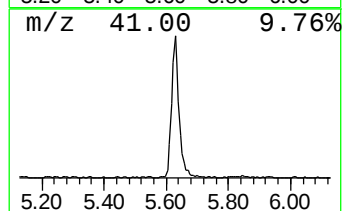
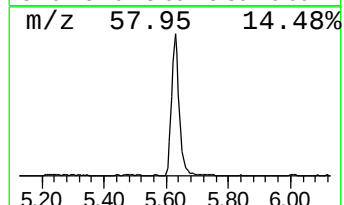
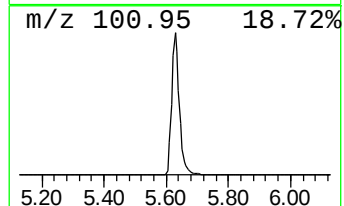
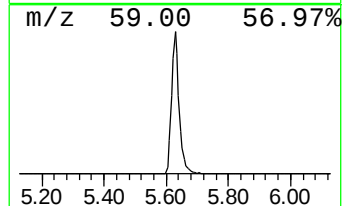
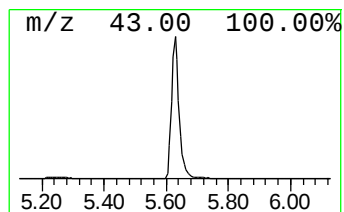
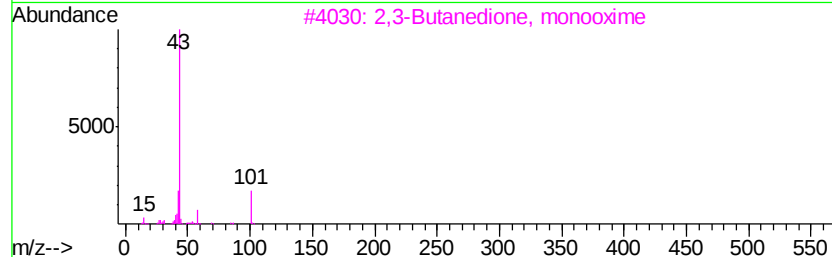
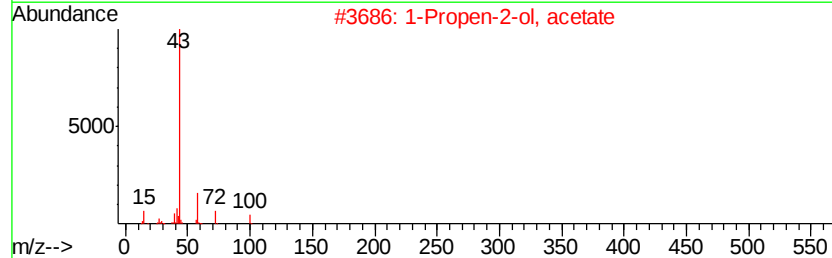
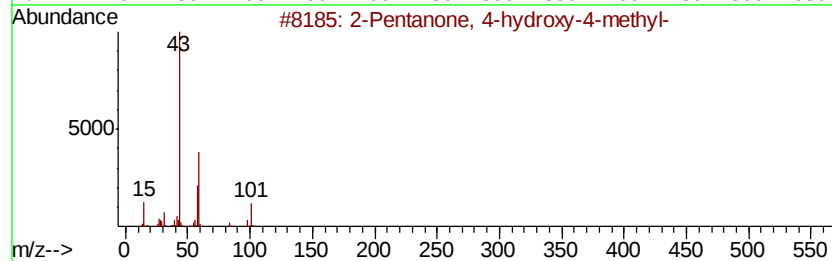
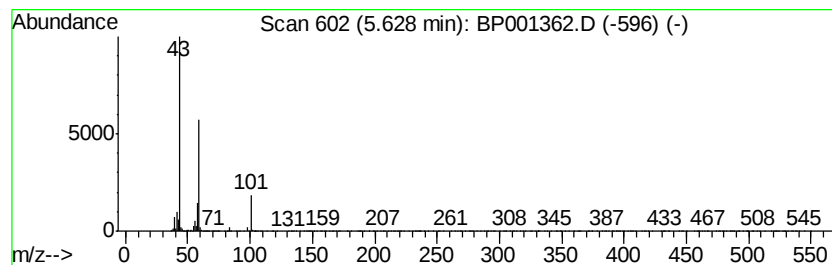
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	19.24 ng	397537	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

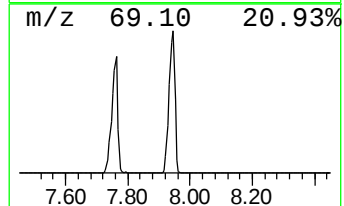
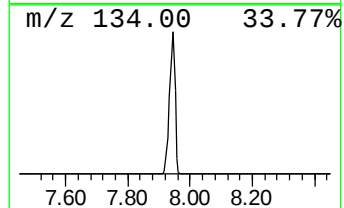
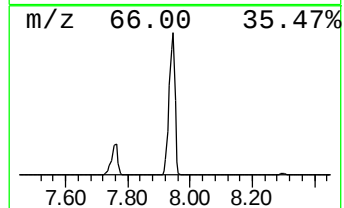
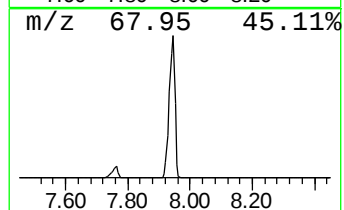
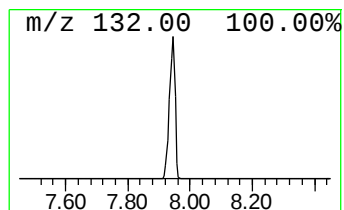
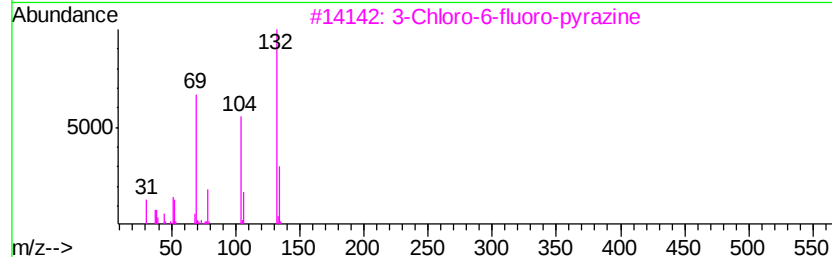
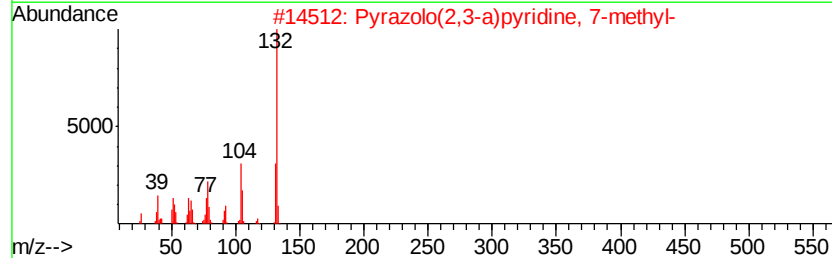
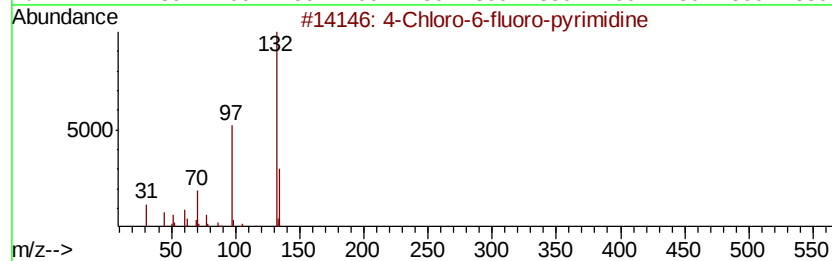
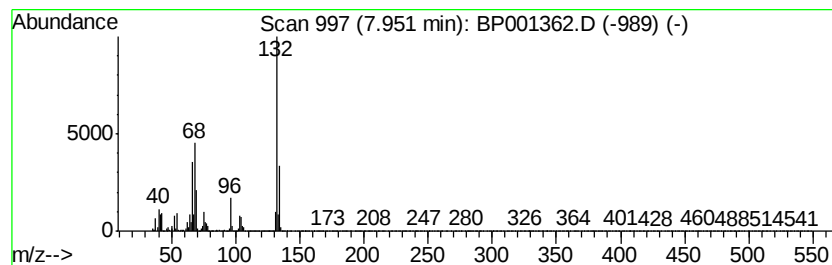
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown7.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	107.14 ng	2213220	1,4-Dichlorobenzene-d4	8.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	10
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

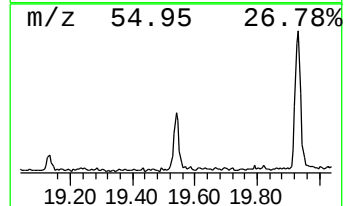
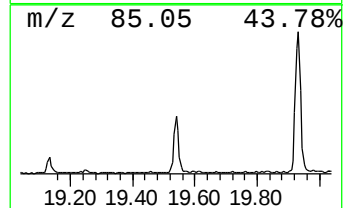
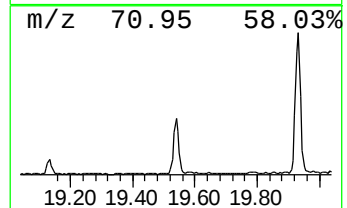
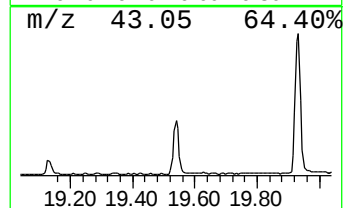
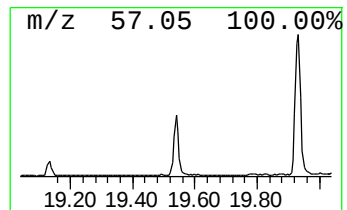
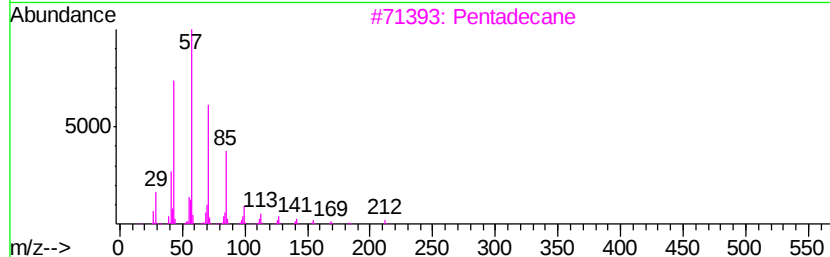
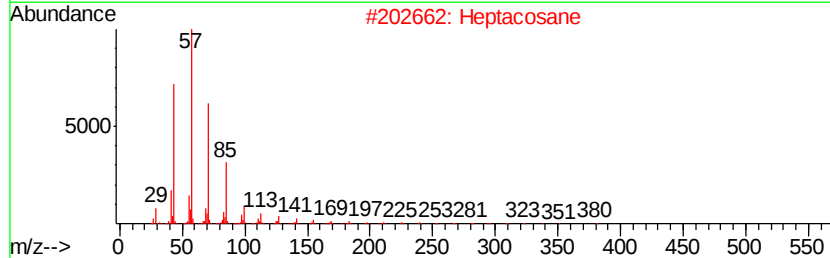
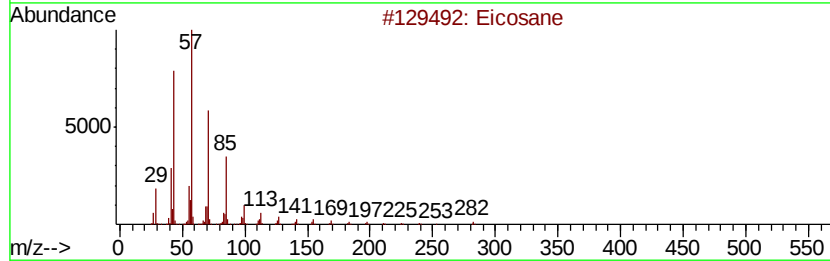
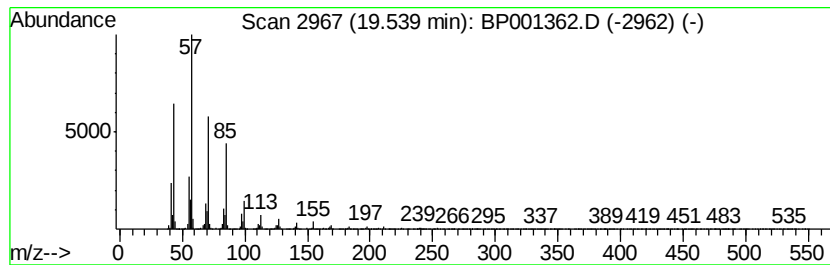
Instrument :
BNA_P
ClientSampleId :
PB125644BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.54	2.83 ng	84658	Chrysene-d12	19.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	94
2	Heptacosane	380 C27H56	000593-49-7	91
3	Pentadecane	212 C15H32	000629-62-9	91
4	Heneicosane	296 C21H44	000629-94-7	90
5	Tricosane	324 C23H48	000638-67-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

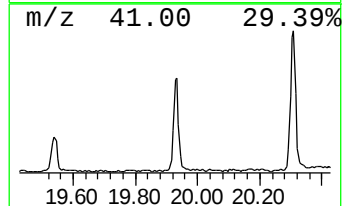
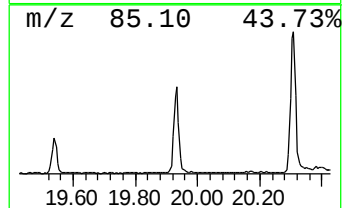
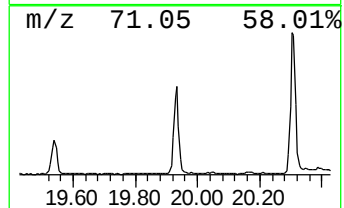
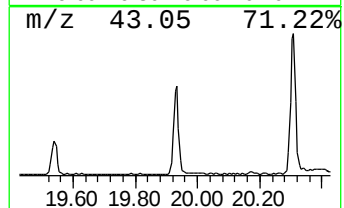
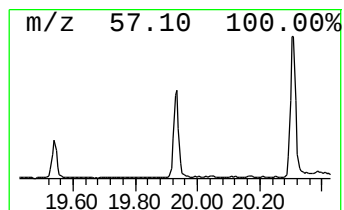
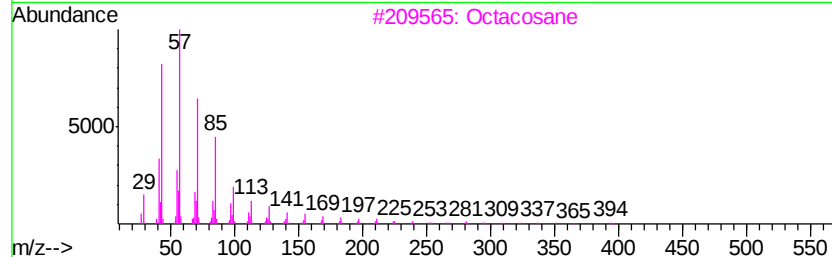
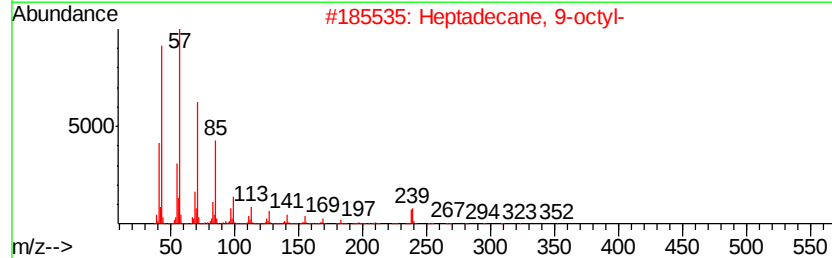
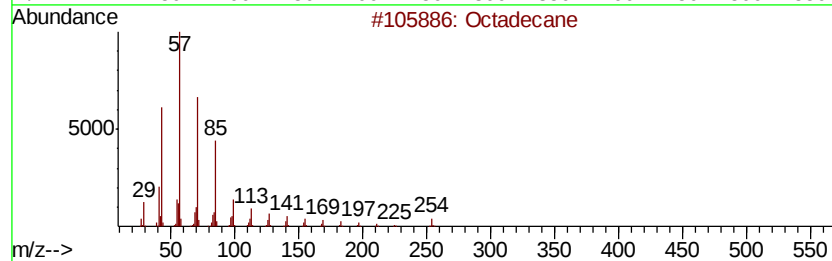
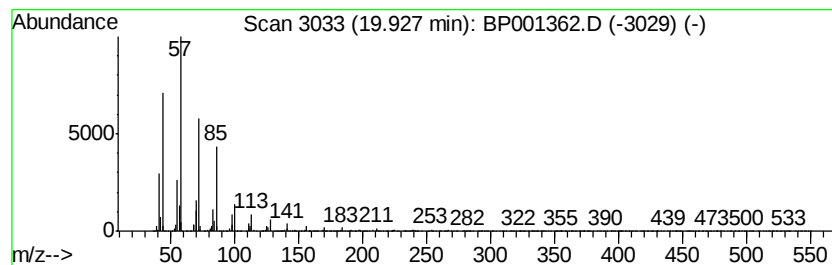
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.93	7.05 ng	211182	Chrysene-d12	19.25

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	95
2	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	95
3	Octacosane		394	C28H58	000630-02-4	94
4	Heptacosane		380	C27H56	000593-49-7	91
5	Heneicosane		296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

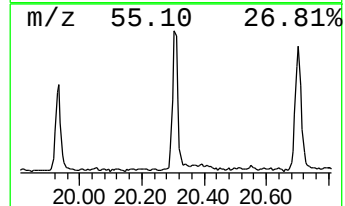
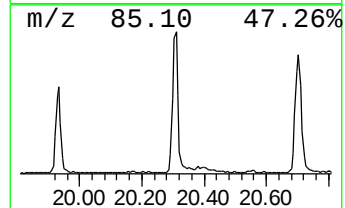
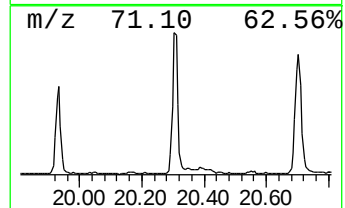
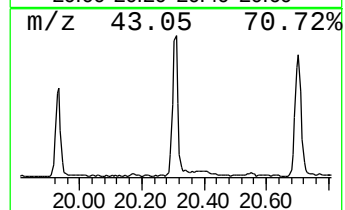
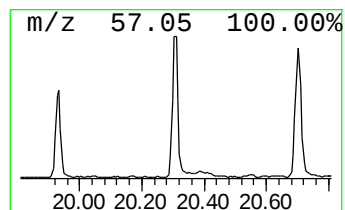
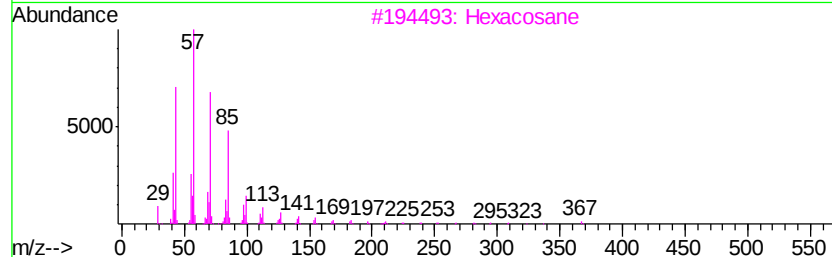
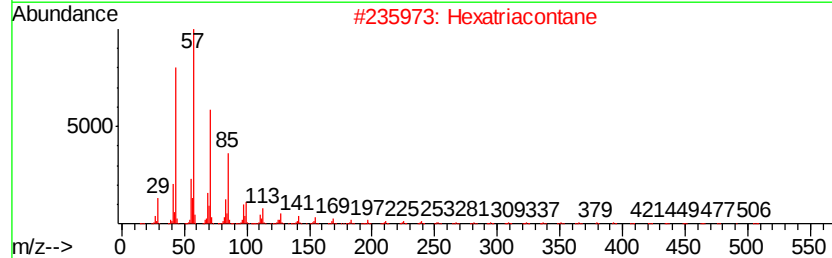
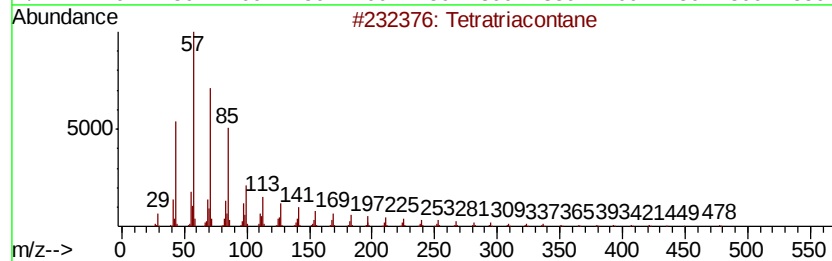
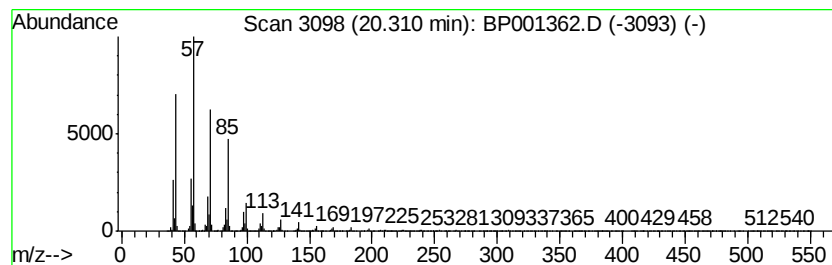
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetratriacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	12.33 ng	355261	Perylene-d12	21.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontane	479	C34H70	014167-59-0	97
2		Hexatriacontane	507	C36H74	000630-06-8	95
3		Hexacosane	366	C26H54	000630-01-3	91
4		Octacosane	394	C28H58	000630-02-4	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

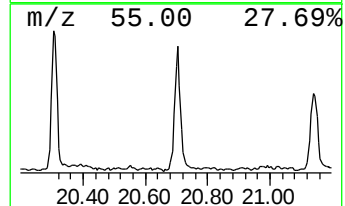
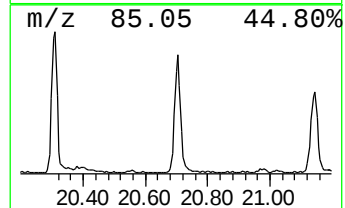
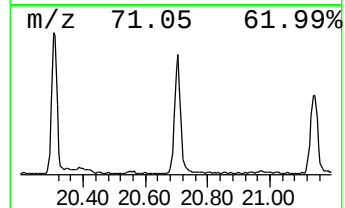
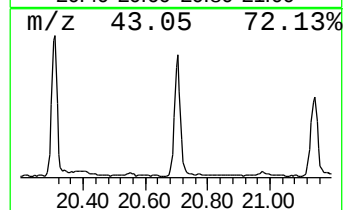
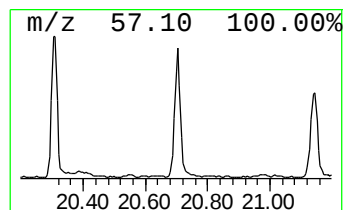
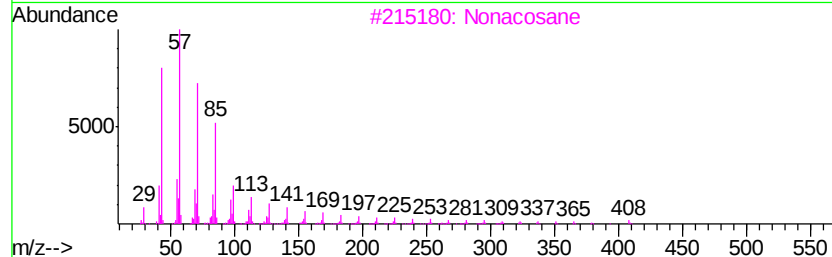
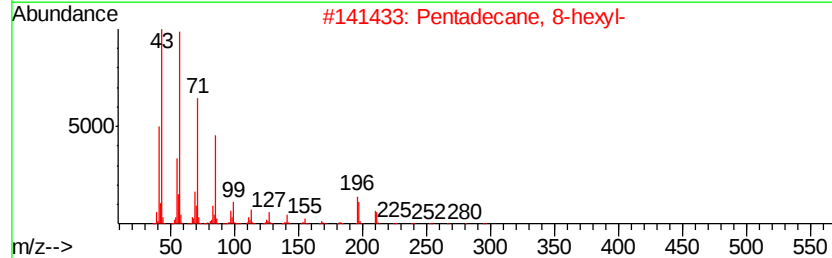
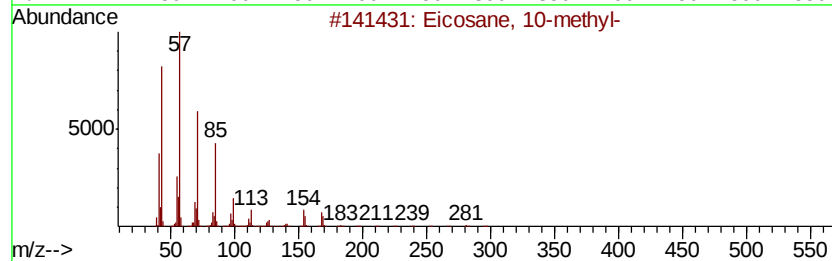
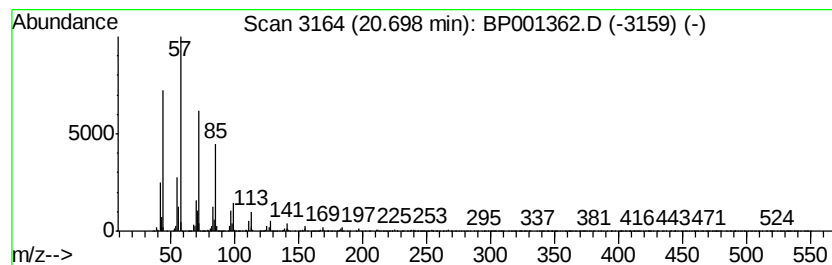
Instrument :
BNA_P
ClientSampleId :
PB125644BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Eicosane, 10-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.70	11.98 ng	345149	Perylene-d12	21.03
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane, 10-methyl-	296 C21H44	054833-23-7	94
2	Pentadecane, 8-hexyl-	296 C21H44	013475-75-7	94
3	Nonacosane	408 C29H60	000630-03-5	93
4	Hexacosane	366 C26H54	000630-01-3	91
5	1-Iodo-2-methylundecane	296 C12H25I	073105-67-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

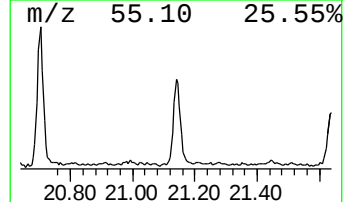
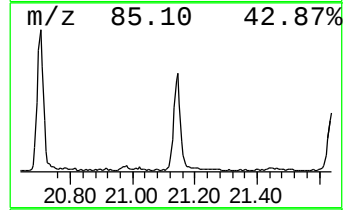
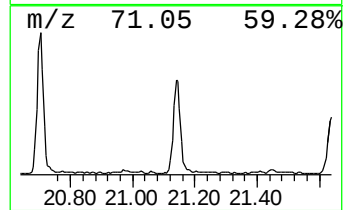
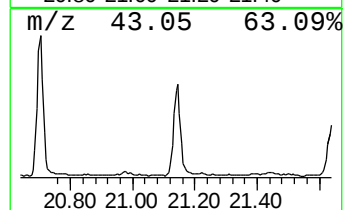
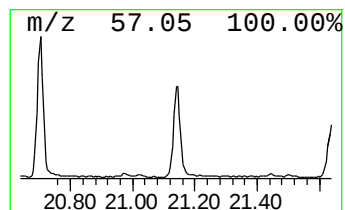
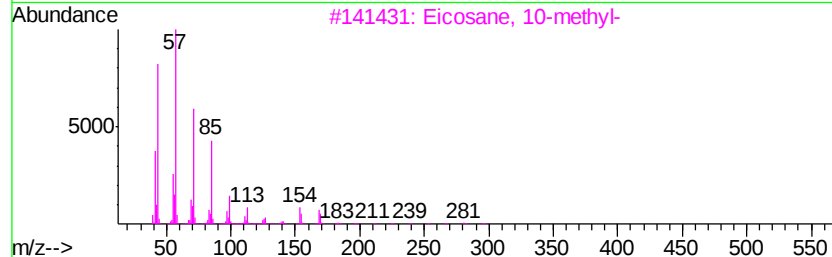
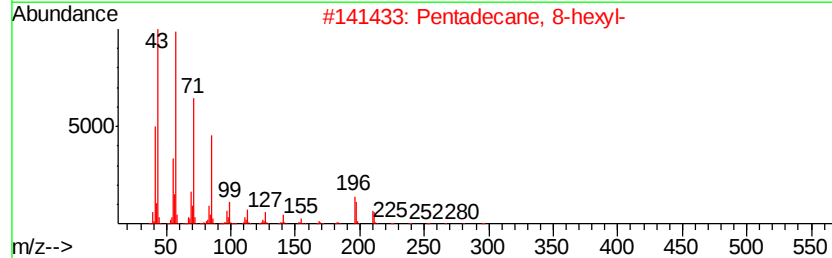
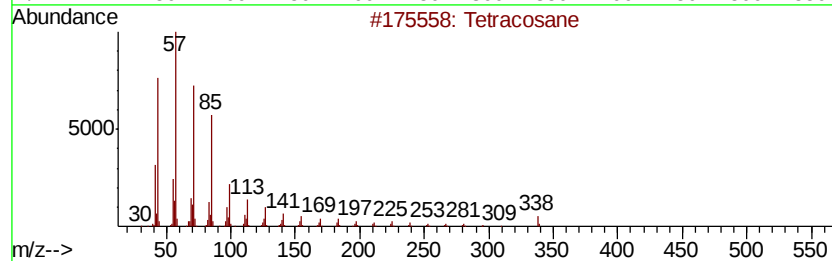
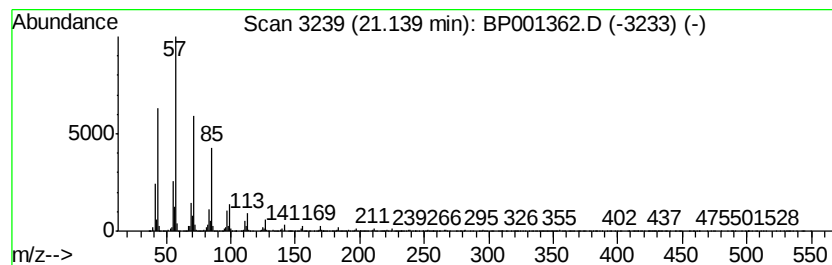
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.14	9.28 ng	267460	Perylene-d12	21.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

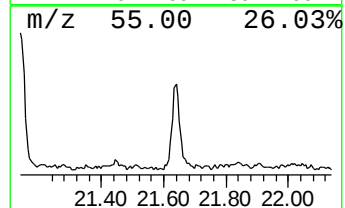
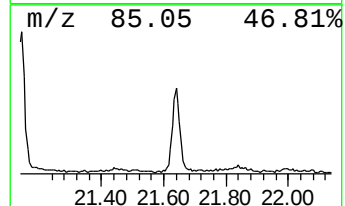
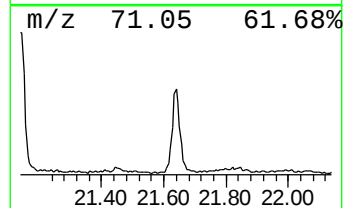
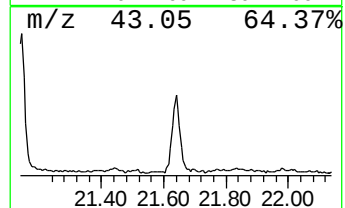
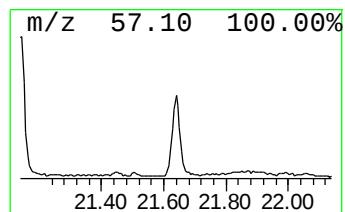
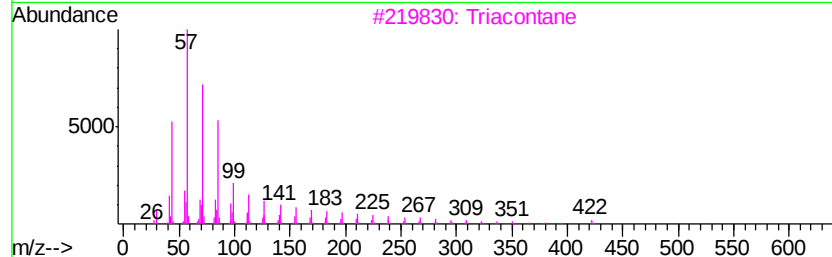
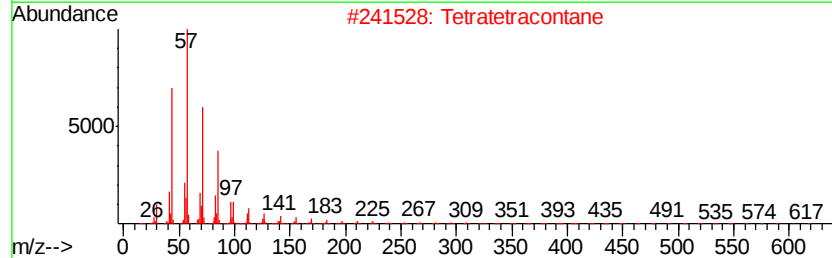
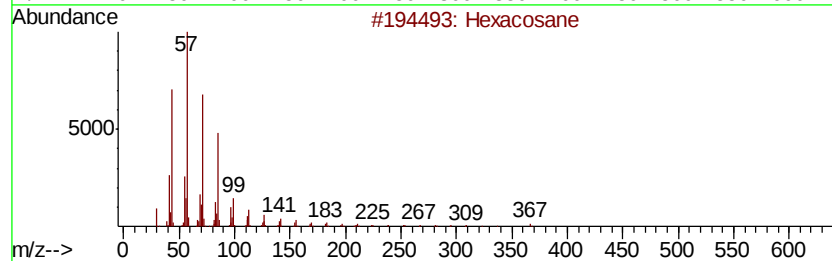
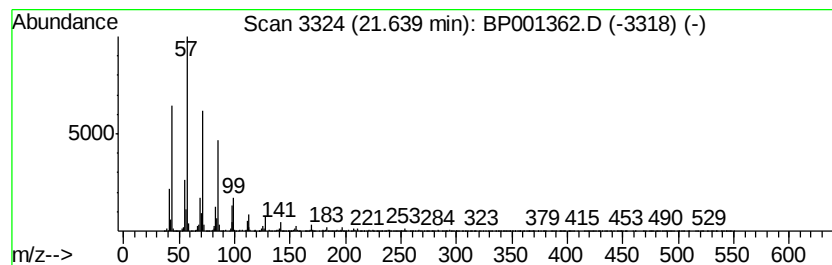
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.64	5.92 ng	170526	Perylene-d12	21.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

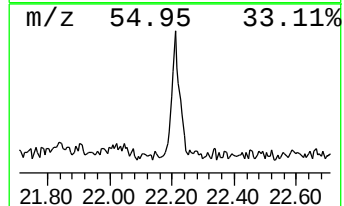
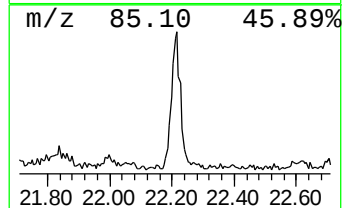
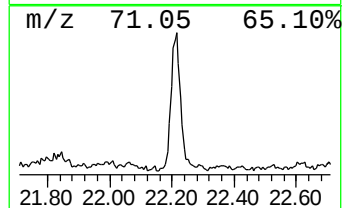
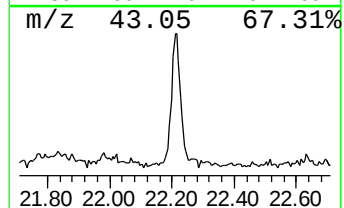
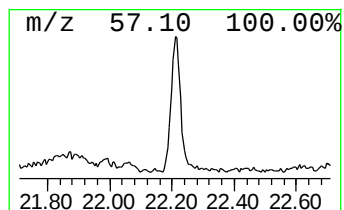
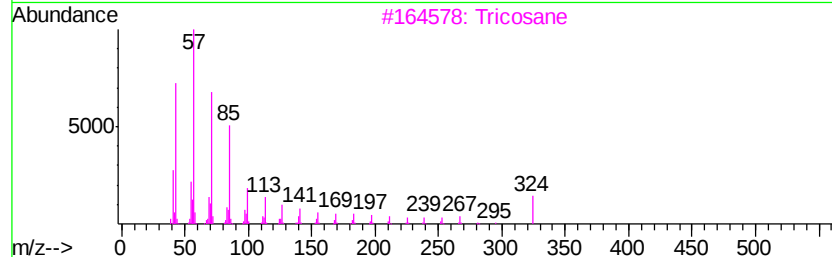
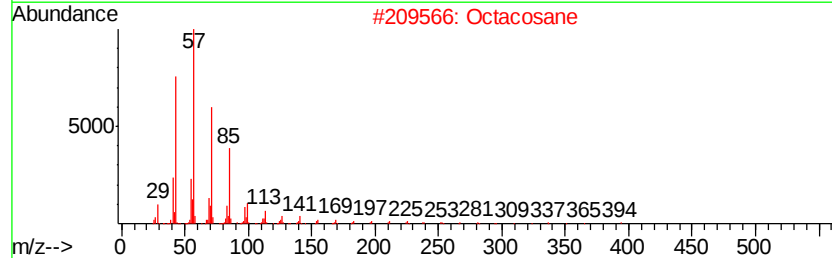
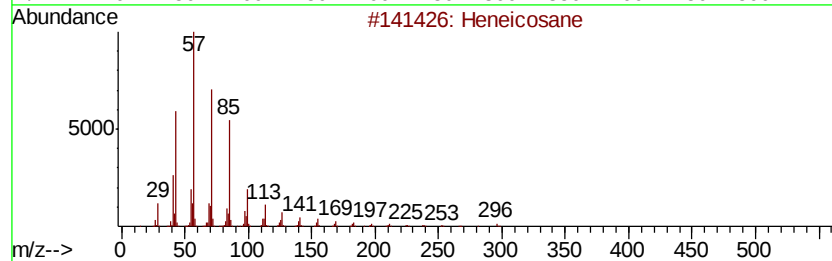
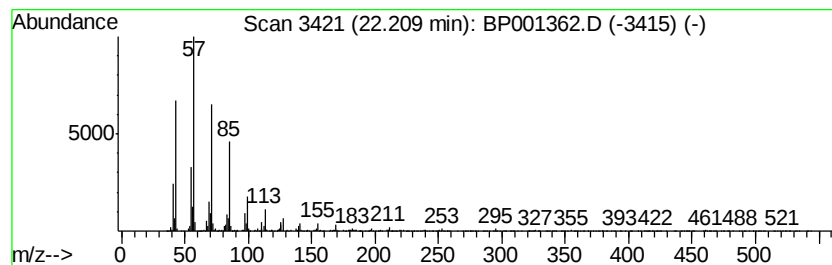
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.21	3.40 ng	98099	Perylene-d12	21.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Tricosane	324	C23H48	000638-67-5	91
4			Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP122319\
Data File : BP001362.D
Acq On : 23 Dec 2019 12:40
Operator : JU
Sample : PB125644BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB125644BL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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...	5.63	19.2	ng	397537	1	8.30	413138	20.0
unknown7.95	7.95	107.1	ng	2213220	1	8.30	413138	20.0
Eicosane	19.54	2.8	ng	84658	5	19.25	599036	20.0
Octadecane	19.93	7.0	ng	211182	5	19.25	599036	20.0
Tetratriacontane	20.31	12.3	ng	355261	6	21.03	576241	20.0
Eicosane, 10-methyl-	20.70	12.0	ng	345149	6	21.03	576241	20.0
Tetracosane	21.14	9.3	ng	267460	6	21.03	576241	20.0
Hexacosane	21.64	5.9	ng	170526	6	21.03	576241	20.0
Heneicosane	22.21	3.4	ng	98099	6	21.03	576241	20.0