

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
 Data File : BP001284.D
 Acq On : 09 Dec 2019 22:14
 Operator : JU
 Sample : K6188-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SOIL-COMP-2

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-BP112719.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	589	602	616	rBV	339441	585119	15.80%	2.093%
2	6.222	696	703	707	rBV	1989348	2615967	70.66%	9.356%
3	7.763	957	965	976	rBV	2072499	2993934	80.87%	10.708%
4	7.951	990	997	1001	rBV	2169728	2780991	75.12%	9.946%
5	8.304	1053	1057	1065	rVB	397833	471078	12.72%	1.685%
6	8.551	1092	1099	1103	rBV	1861446	2154083	58.18%	7.704%
7	9.193	1200	1208	1212	rBV	1429388	1809364	48.87%	6.471%
8	10.339	1398	1403	1413	rVB	485777	555233	15.00%	1.986%
9	12.122	1696	1706	1710	rBV	2717402	3236381	87.42%	11.575%
10	12.786	1814	1819	1828	rBV	141132	162406	4.39%	0.581%
11	13.204	1883	1890	1902	rBV	528666	660074	17.83%	2.361%
12	14.498	2103	2110	2128	rBV	1817805	2311243	62.43%	8.266%
13	15.598	2292	2297	2309	rVB	621655	702770	18.98%	2.513%
14	16.492	2444	2449	2468	rBV2	108140	173894	4.70%	0.622%
15	17.886	2680	2686	2690	rBV	3386776	3702162	100.00%	13.241%
16	19.121	2891	2896	2911	rBV2	204341	463270	12.51%	1.657%
17	19.239	2911	2916	2921	rVV	592947	658509	17.79%	2.355%
18	19.533	2962	2966	2974	rBV	74601	85024	2.30%	0.304%
19	19.921	3028	3032	3035	rBV	104035	109151	2.95%	0.390%
20	20.298	3091	3096	3100	rBV	139644	154941	4.19%	0.554%
21	20.692	3158	3163	3168	rBV	152334	185034	5.00%	0.662%
22	21.010	3211	3217	3225	rVB	467310	654329	17.67%	2.340%
23	21.127	3232	3237	3246	rVB	127510	186406	5.04%	0.667%
24	21.627	3316	3322	3329	rBV	104962	174376	4.71%	0.624%
25	21.704	3329	3335	3341	rBV6	31449	93914	2.54%	0.336%
26	22.198	3413	3419	3433	rVB3	61471	144335	3.90%	0.516%
27	22.868	3526	3533	3544	rVB3	33752	95740	2.59%	0.342%
28	23.651	3661	3666	3675	rVB4	17387	40584	1.10%	0.145%

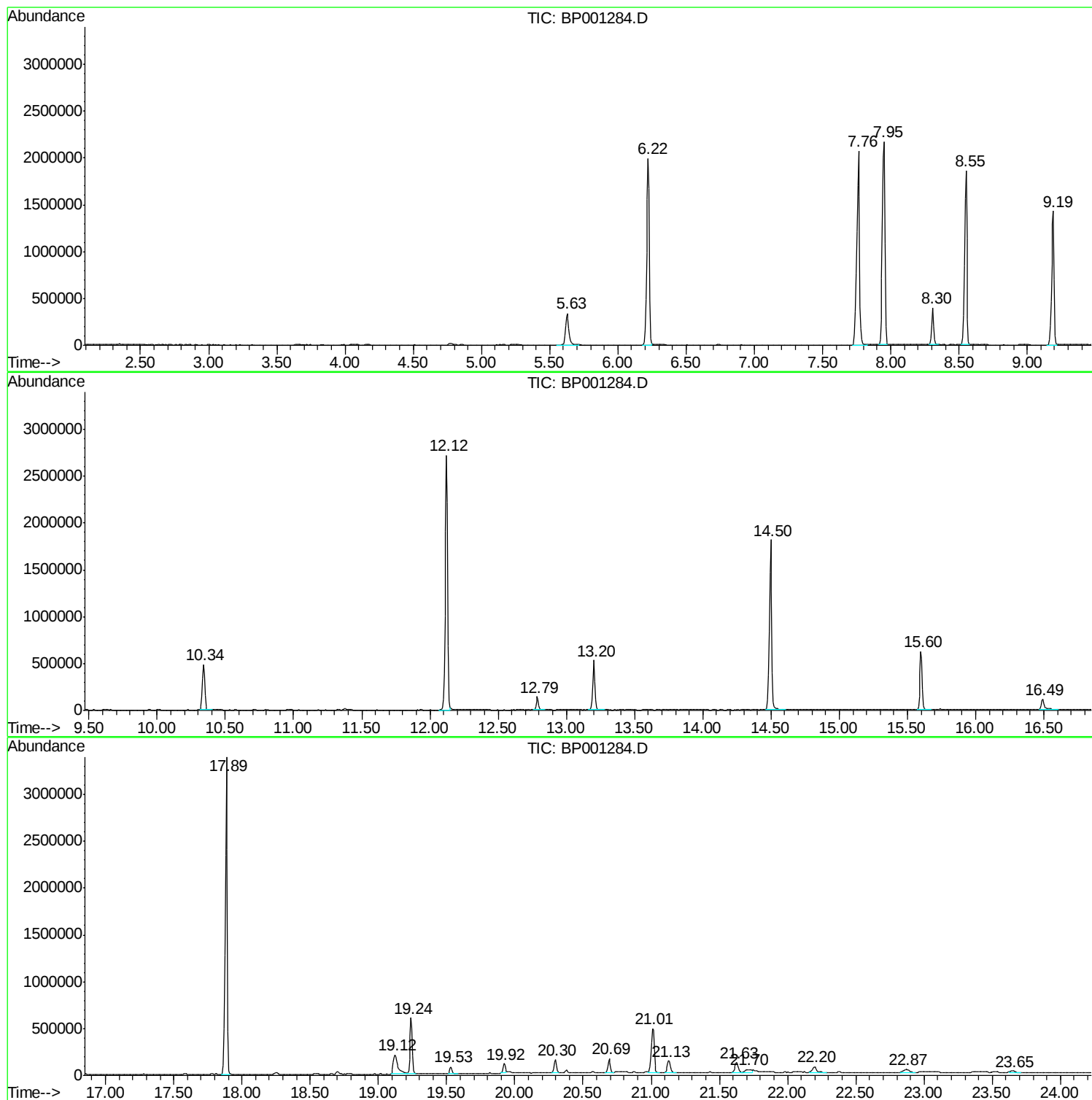
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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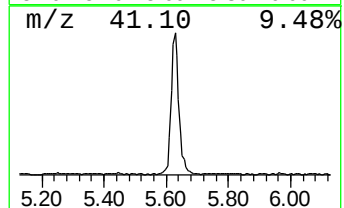
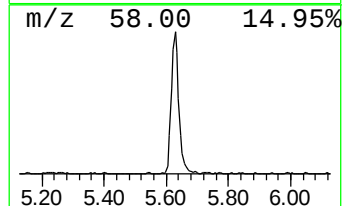
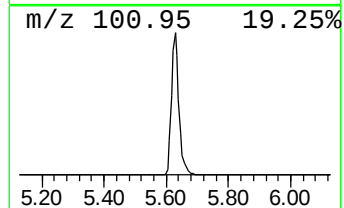
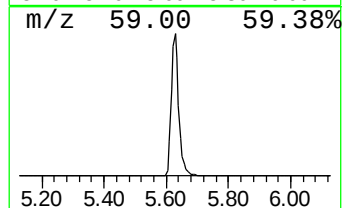
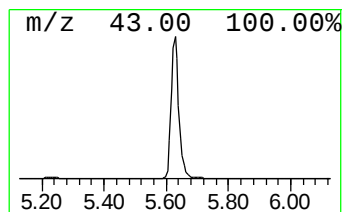
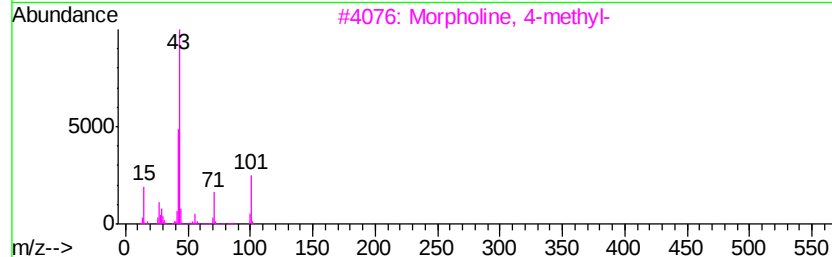
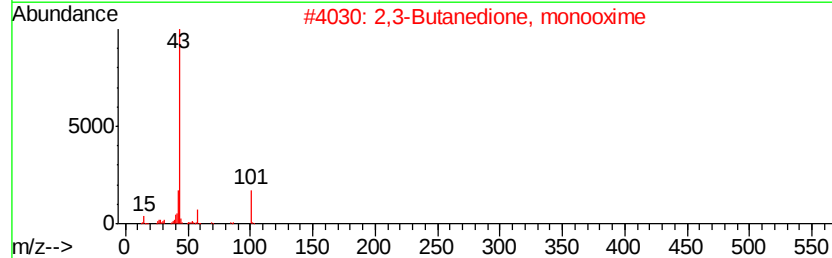
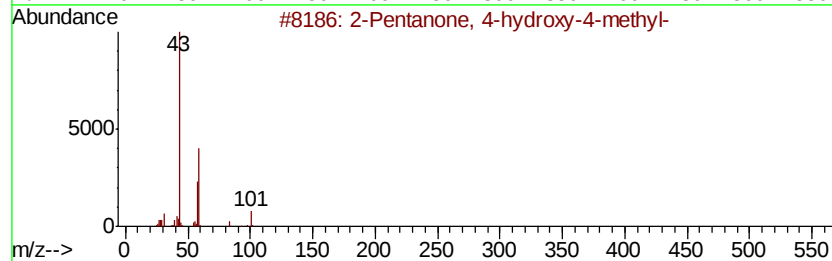
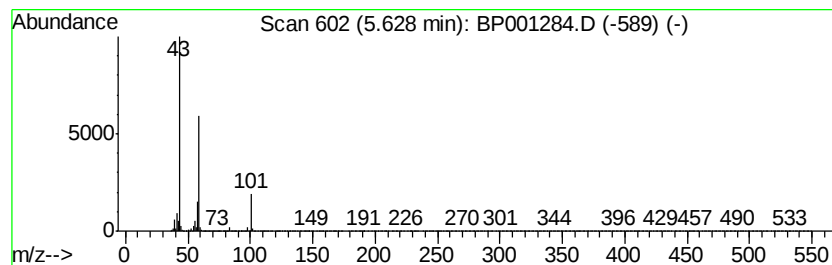
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	24.84 ng	585119	1,4-Dichlorobenzene-d4	8.31

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Butane	58	C4H10	000106-97-8	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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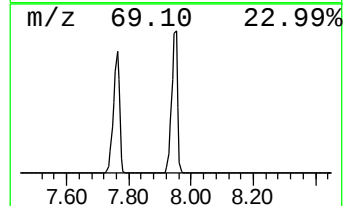
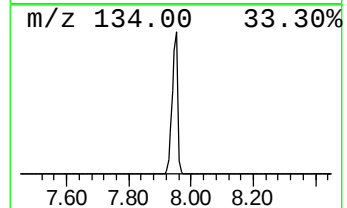
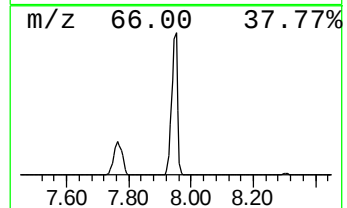
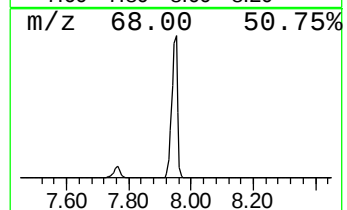
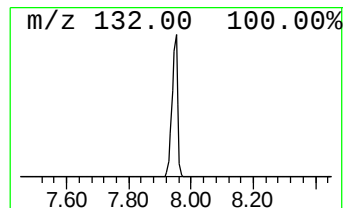
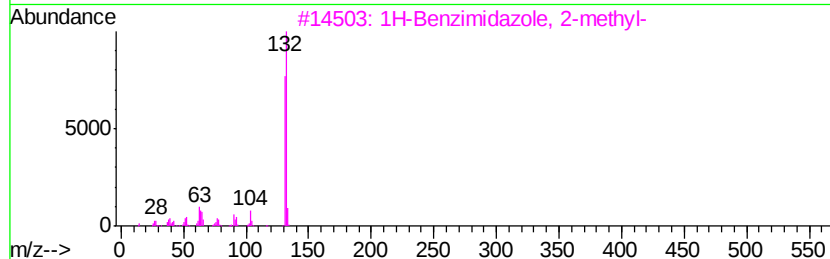
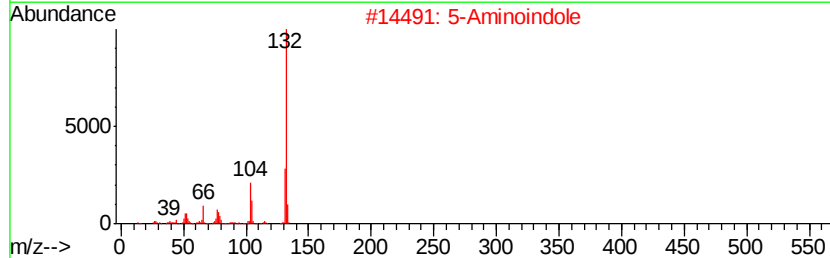
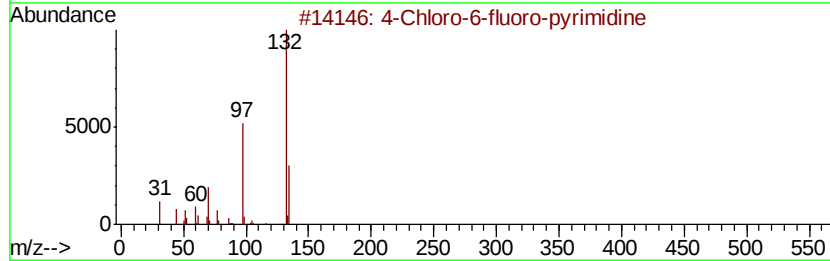
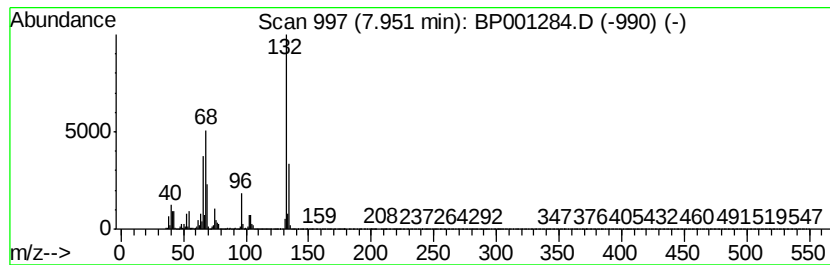
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Peak Number 2 unknown7.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	118.07 ng	2780990	1,4-Dichlorobenzene-d4	8.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14



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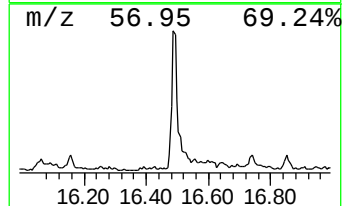
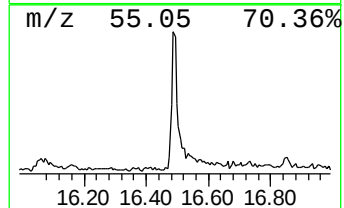
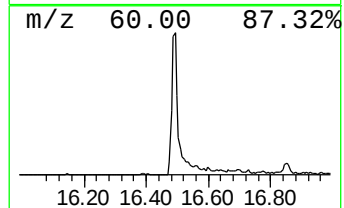
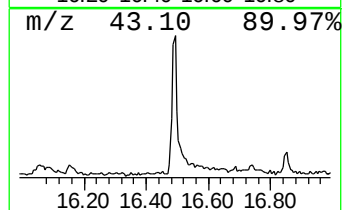
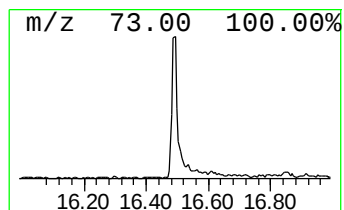
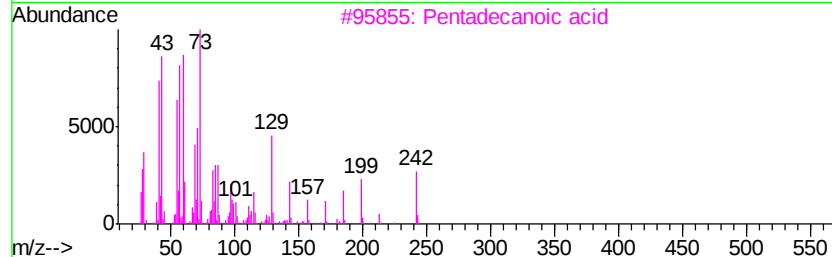
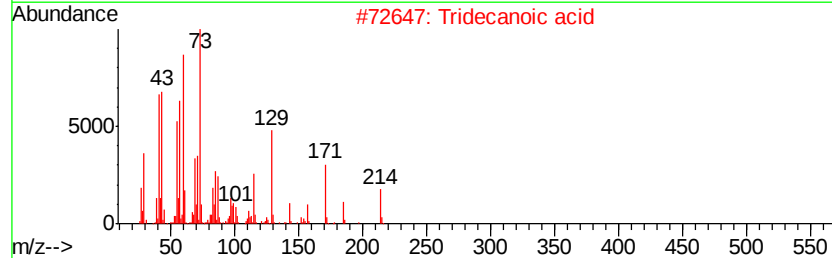
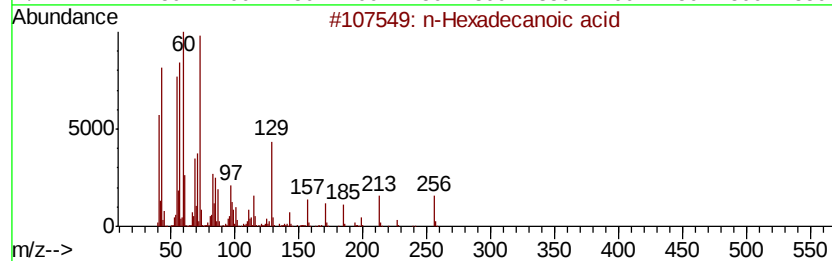
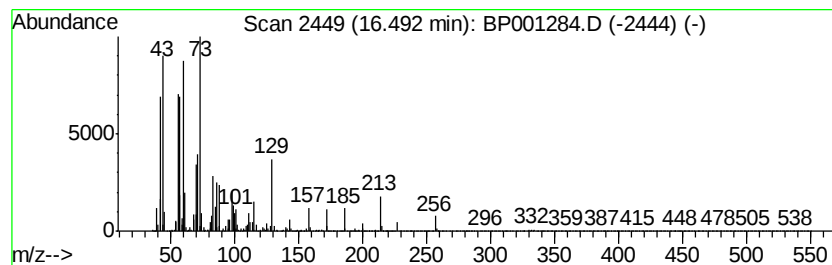
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	4.95 ng	173894	Phenanthrene-d10	15.60

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		Octadecanoic acid	284	C18H36O2	000057-11-4	87
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	83



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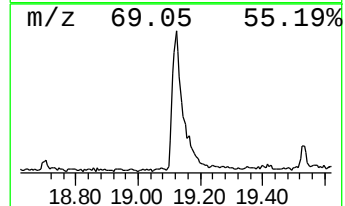
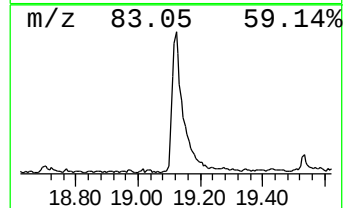
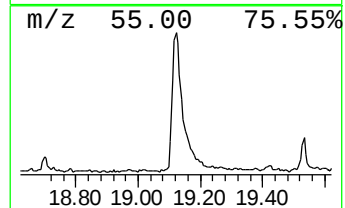
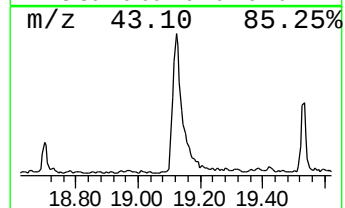
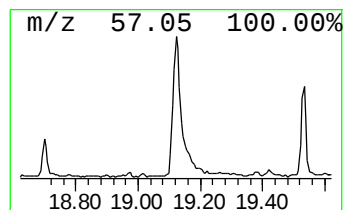
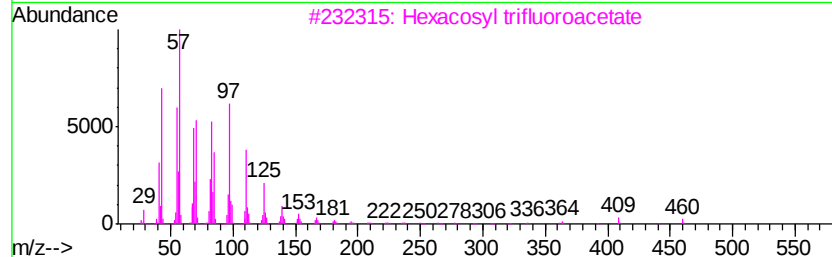
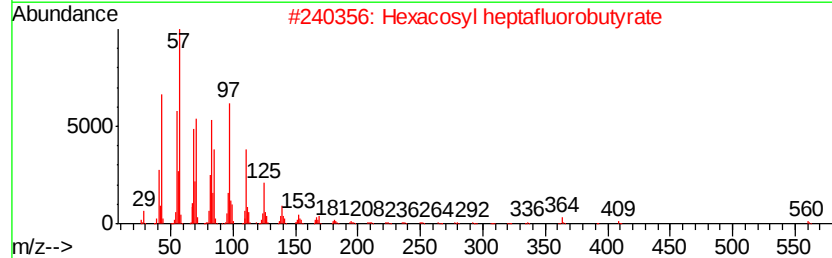
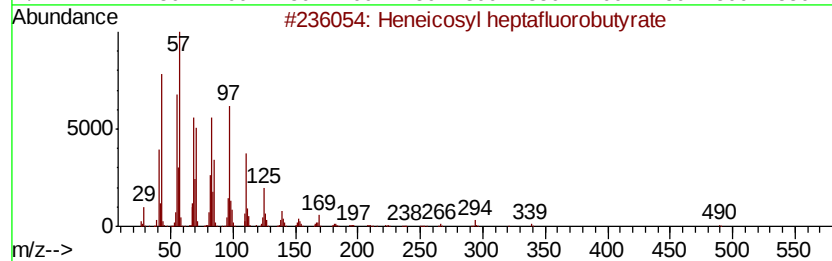
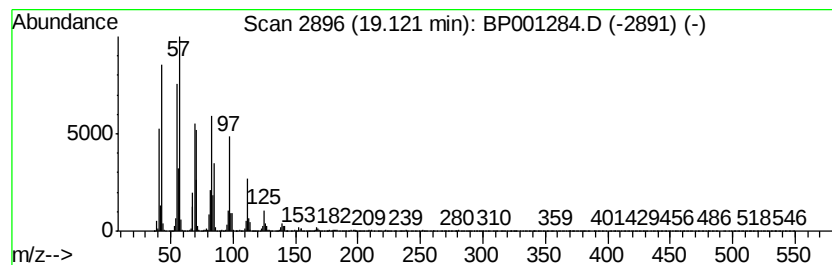
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Peak Number 5 Heneicosyl heptafluorobutyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.12	14.07 ng	463270	Chrysene-d12	19.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	91
2		Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	91
3		Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	91
4		Heptafluorobutyric acid, n-octadecyl	466	C22H37F7O2	000400-57-7	91
5		Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	91



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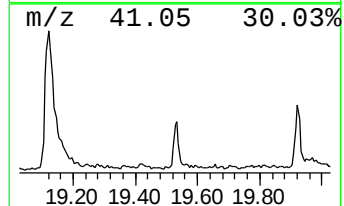
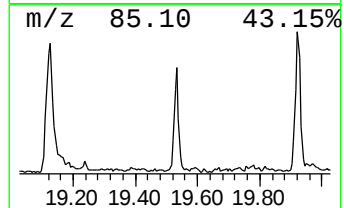
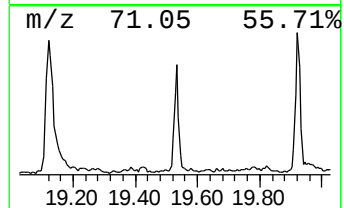
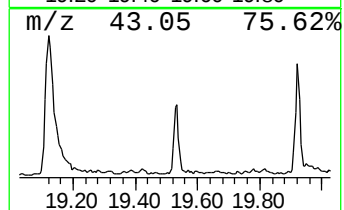
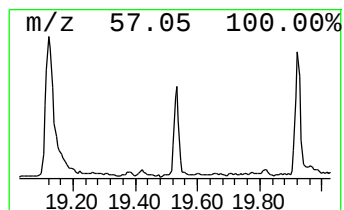
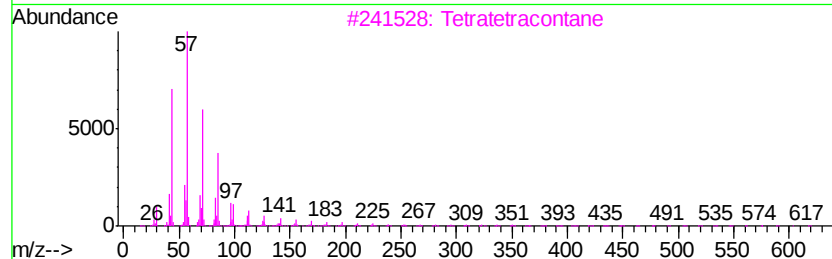
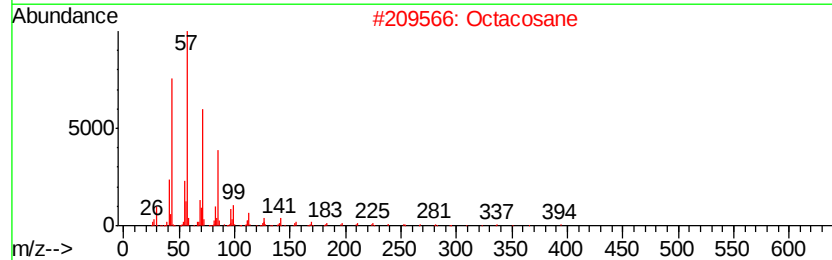
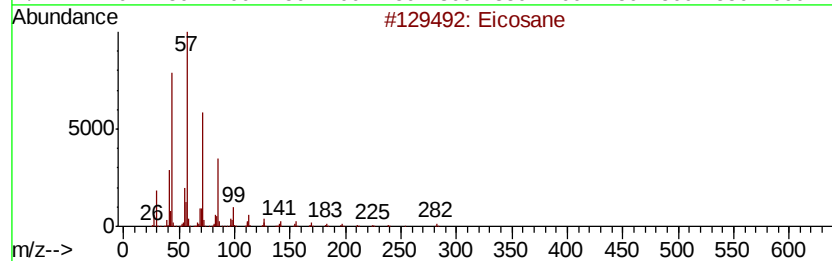
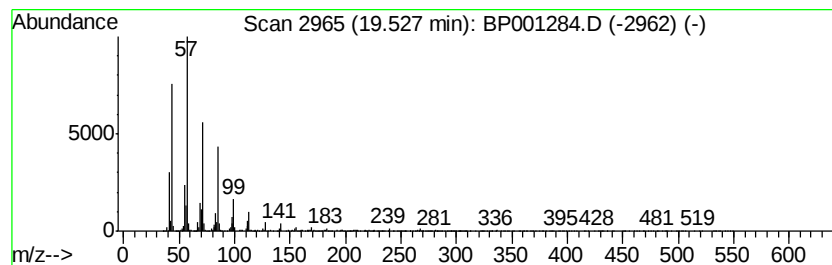
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	2.58 ng	85024	Chrysene-d12	19.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	93
2	Octacosane	394 C28H58	000630-02-4	91
3	Tetratetracontane	619 C44H90	007098-22-8	90
4	Octadecane	254 C18H38	000593-45-3	90
5	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	87



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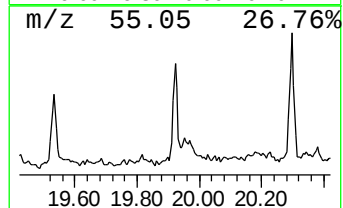
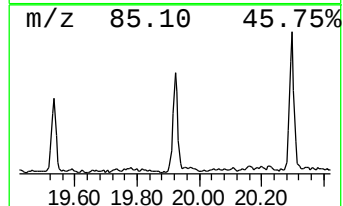
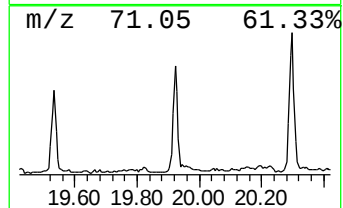
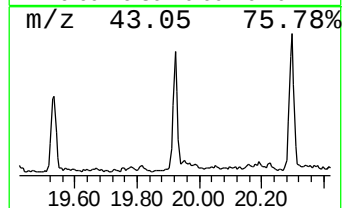
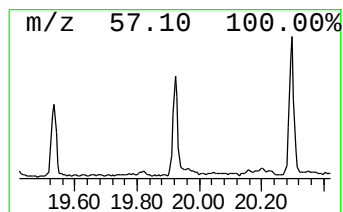
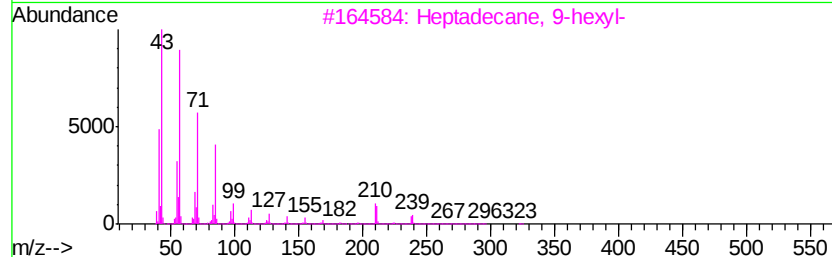
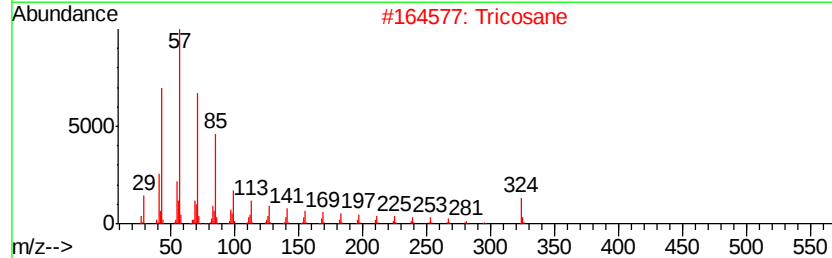
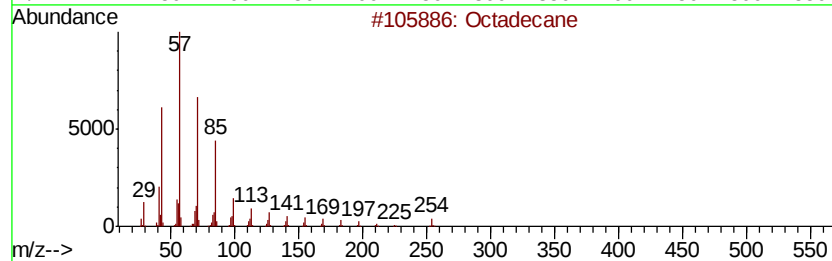
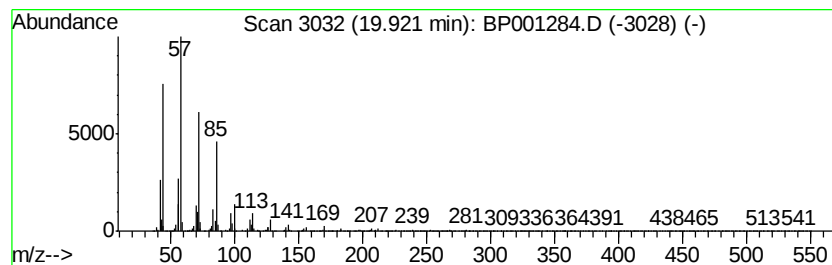
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.92	3.32 ng	109151	Chrysene-d12	19.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	95
2		Tricosane	324	C23H48	000638-67-5	93
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001284.D
Acq On : 09 Dec 2019 22:14
Operator : JU
Sample : K6188-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SOIL-COMP-2

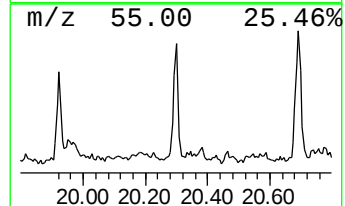
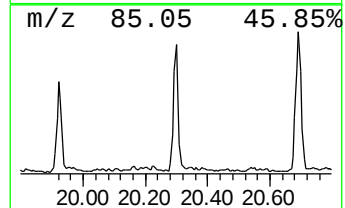
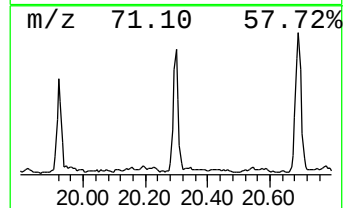
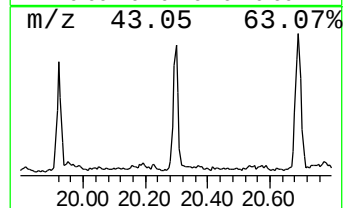
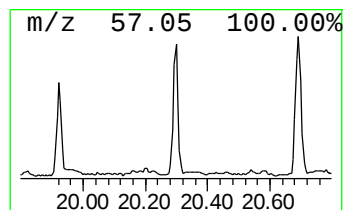
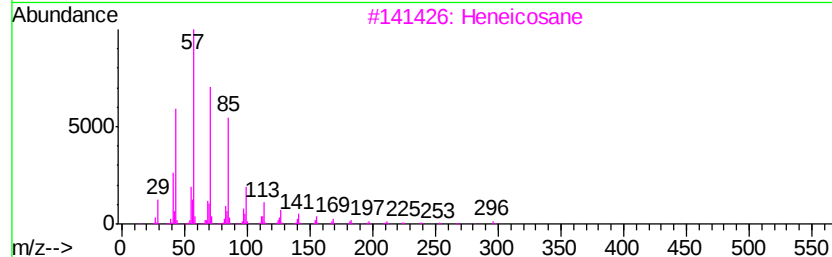
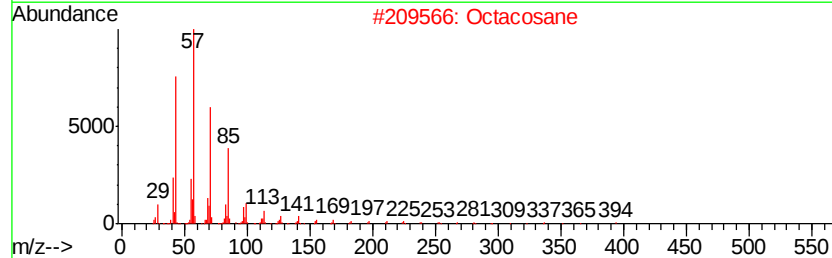
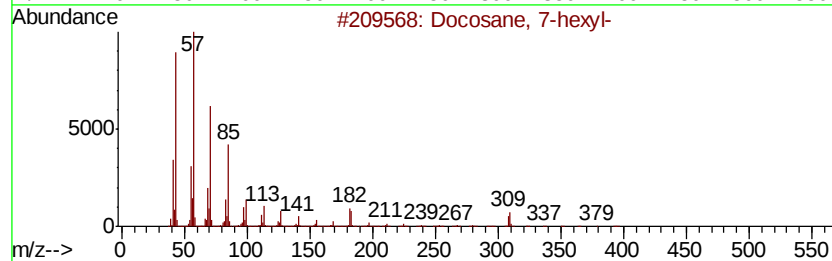
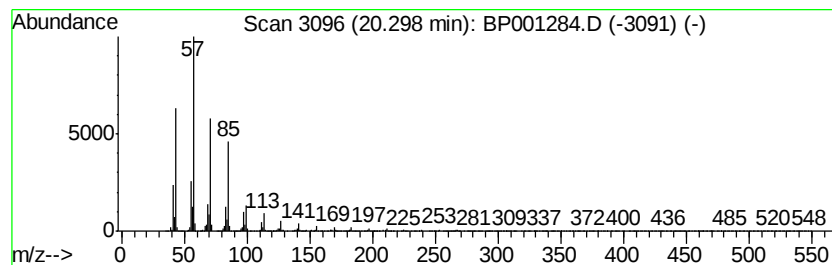
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Docosane, 7-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	4.74 ng	154941	Perylene-d12	21.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 7-hexyl-	394	C28H58	055373-86-9	96
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001284.D
Acq On : 09 Dec 2019 22:14
Operator : JU
Sample : K6188-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SOIL-COMP-2

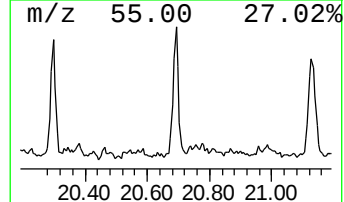
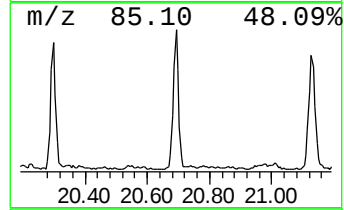
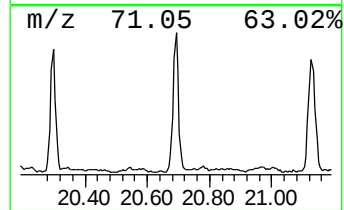
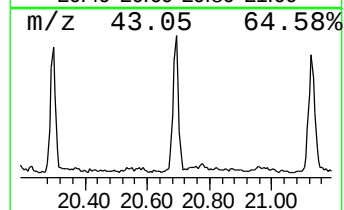
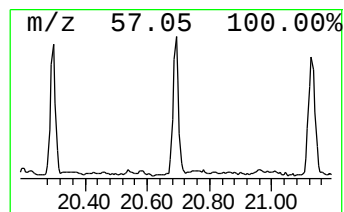
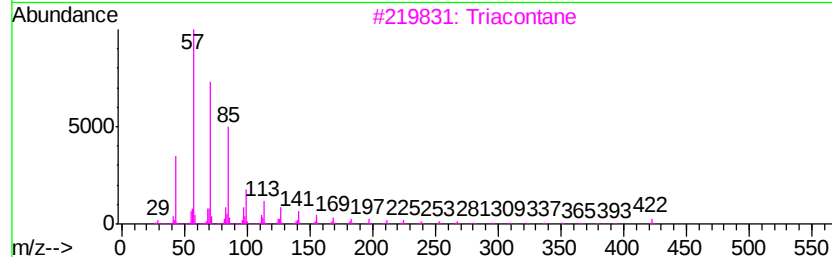
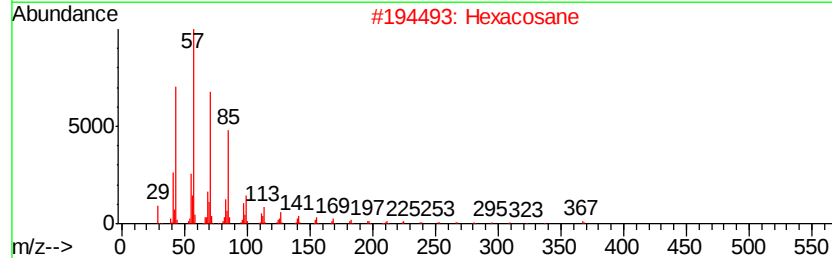
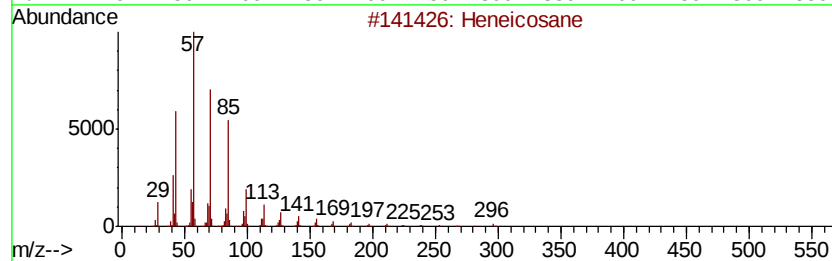
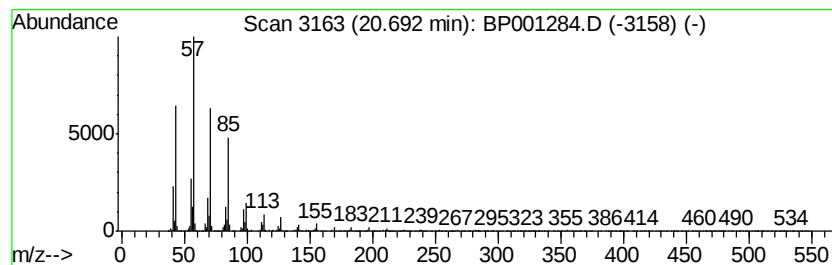
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	5.66 ng	185034	Perylene-d12	21.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	triacontane		422	C30H62	000638-68-6	91
4	Nonadecane		268	C19H40	000629-92-5	87
5	Eicosane		282	C20H42	000112-95-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001284.D
Acq On : 09 Dec 2019 22:14
Operator : JU
Sample : K6188-05
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SOIL-COMP-2

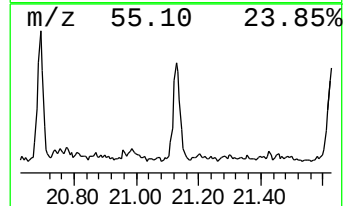
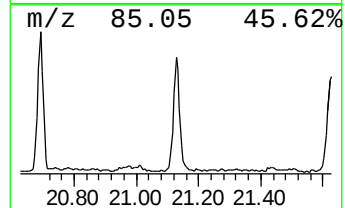
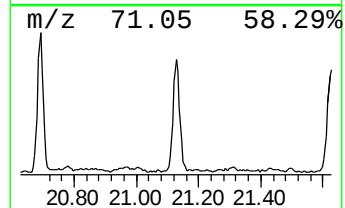
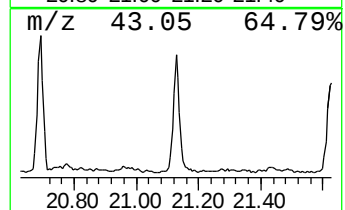
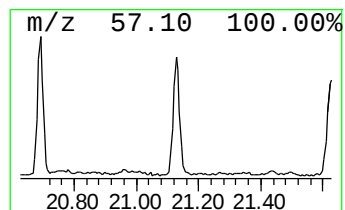
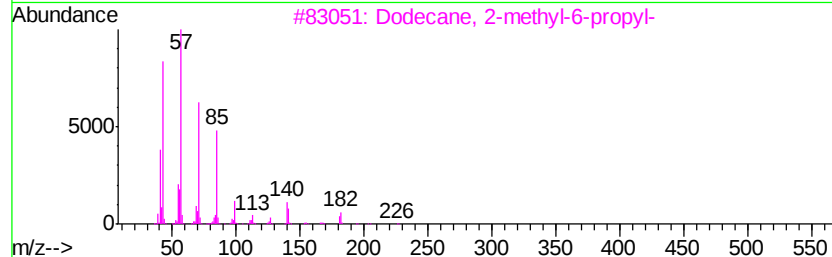
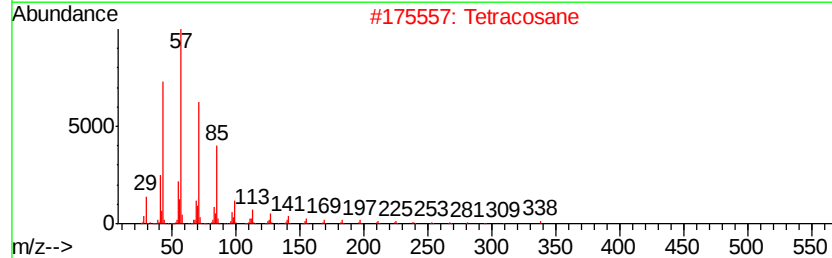
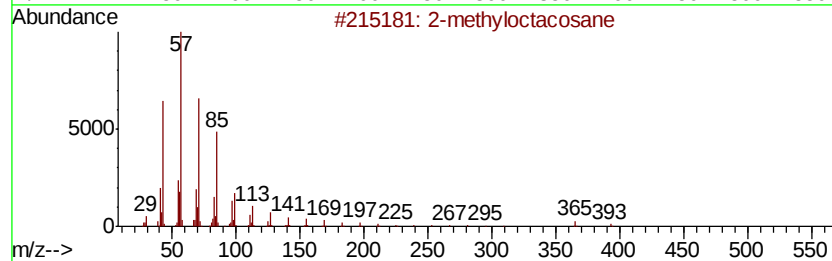
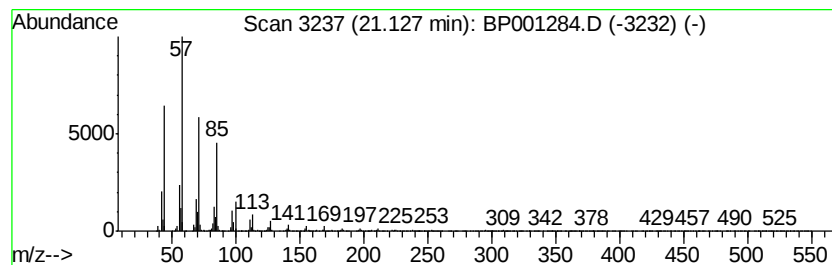
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-methyloctacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.13	5.70 ng	186406	Perylene-d12	21.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	90
2			Tetracosane	338	C24H50	000646-31-1	90
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4			Pentadecane	212	C15H32	000629-62-9	87
5			Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
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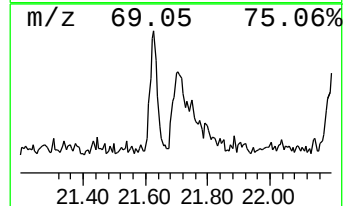
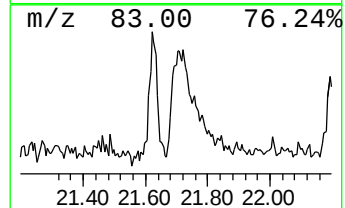
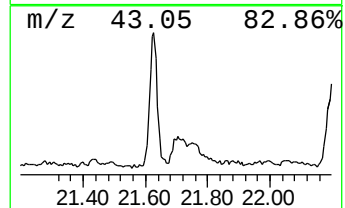
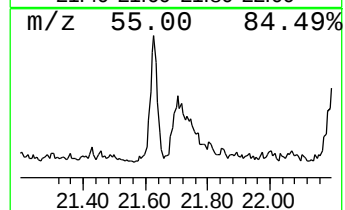
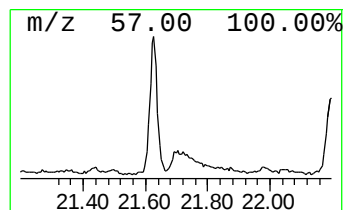
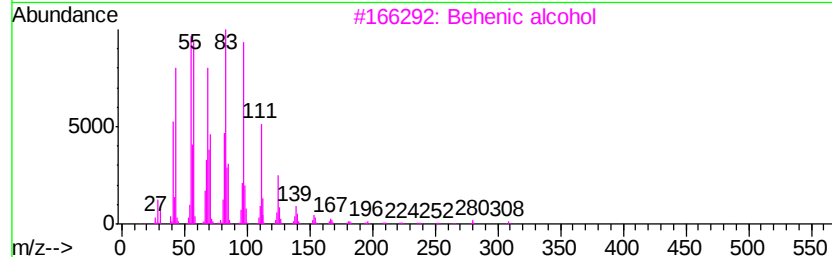
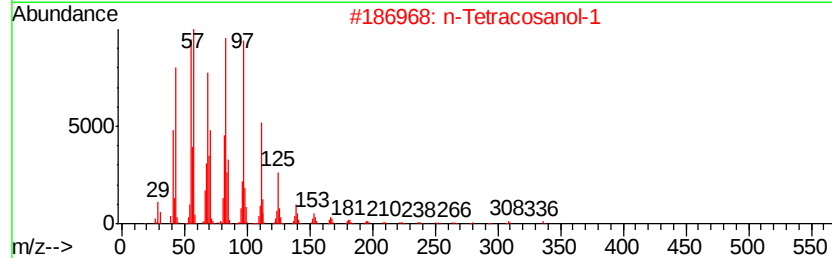
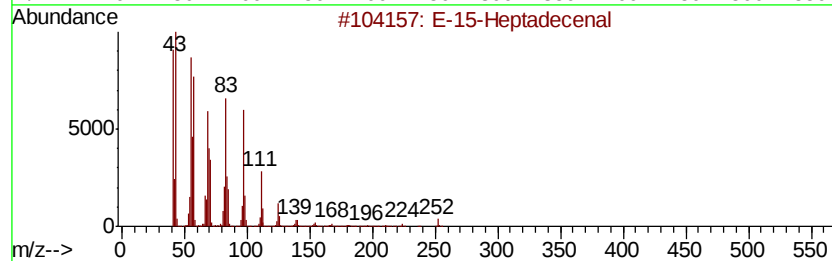
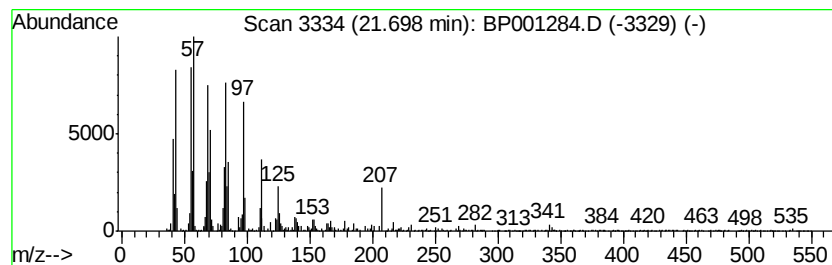
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 E-15-Heptadecenal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.70	2.87 ng	93914	Perylene-d12		21.01		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	E-15-Heptadecenal			252	C17H32O	1000130-97-9	95
2	n-Tetracosanol-1			354	C24H50O	000506-51-4	90
3	Behenic alcohol			326	C22H46O	000661-19-8	90
4	1-Docosene			308	C22H44	001599-67-3	86
5	5-Eicosene, (E)-			280	C20H40	074685-30-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001284.D
Acq On : 09 Dec 2019 22:14
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Misc :
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Instrument :
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ClientSampleId :
SOIL-COMP-2

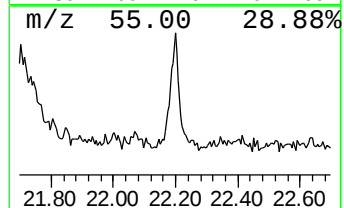
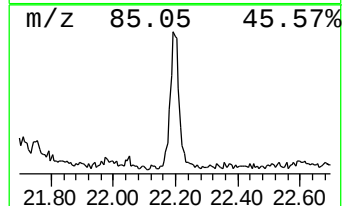
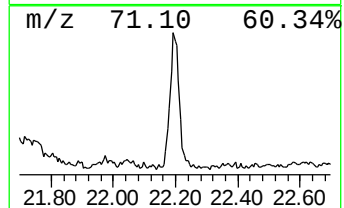
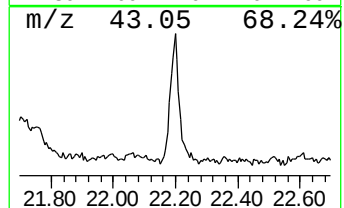
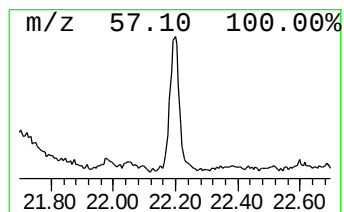
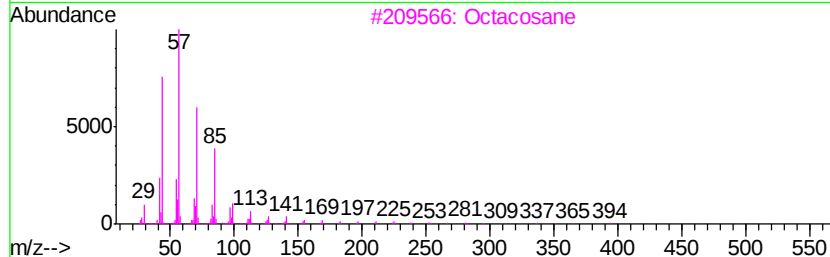
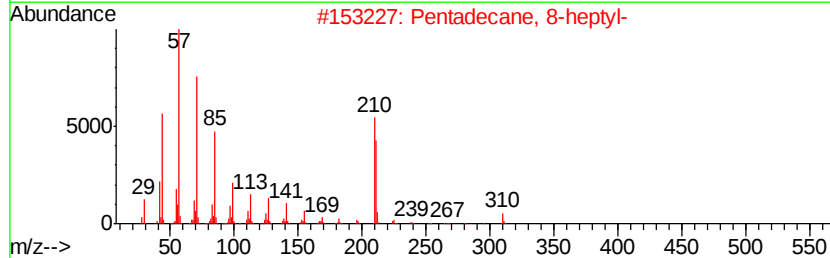
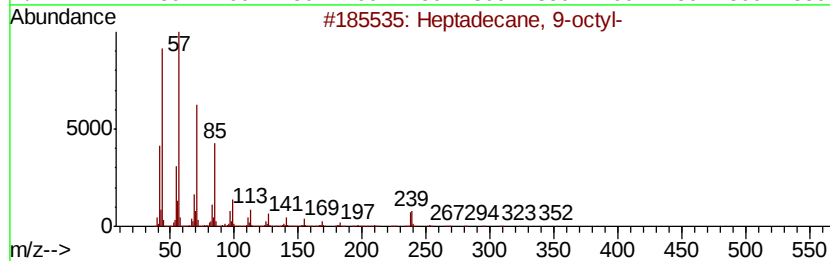
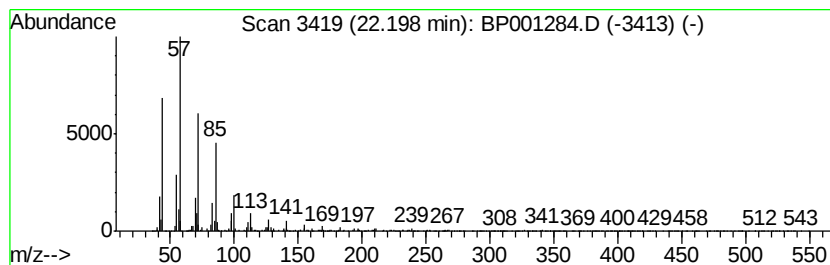
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 9-octyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.20	4.41 ng	144335	Perylene-d12	21.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	94
3			Octacosane	394	C28H58	000630-02-4	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
Data File : BP001284.D
Acq On : 09 Dec 2019 22:14
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
SOIL-COMP-2

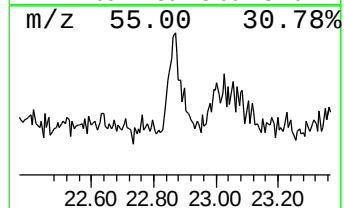
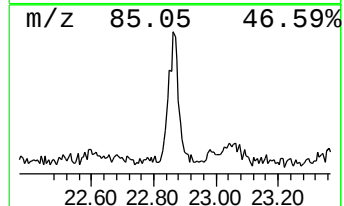
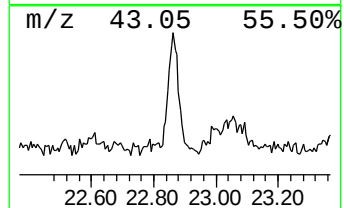
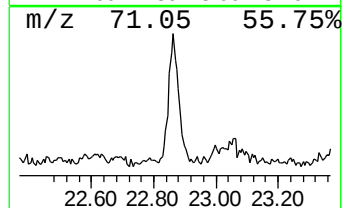
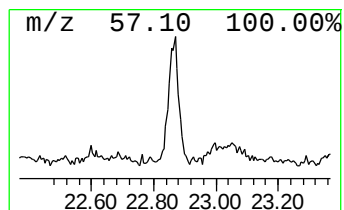
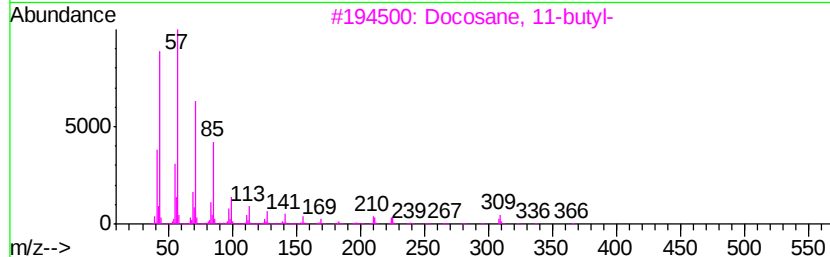
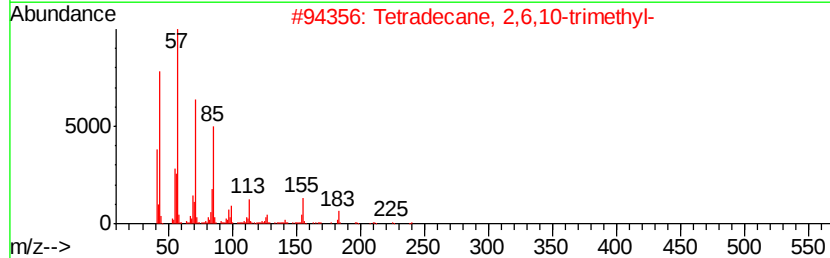
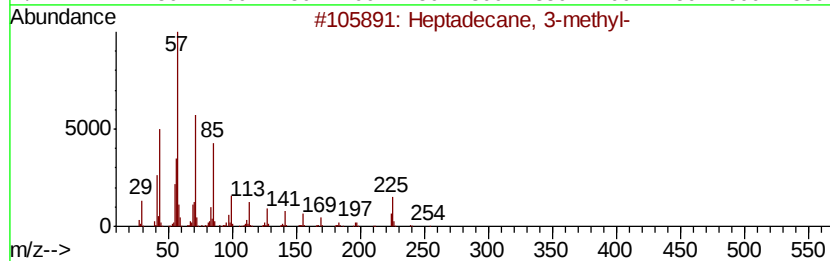
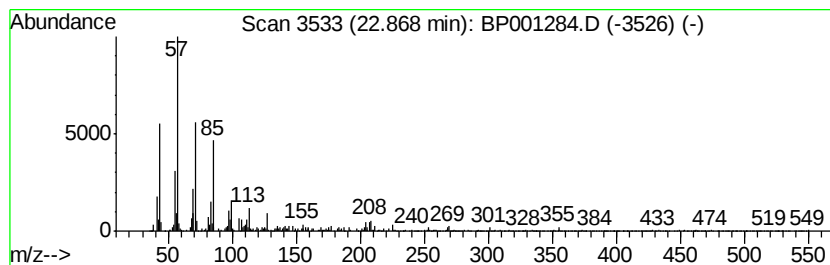
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecane, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.87	2.93 ng	95740	Perylene-d12	21.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	91
2		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	90
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
5		Docosane, 9-butyl-	366	C26H54	055282-14-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120919\
 Data File : BP001284.D
 Acq On : 09 Dec 2019 22:14
 Operator : JU
 Sample : K6188-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SOIL-COMP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.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	24.8	ng	585119	1	8.31	471078	20.0
unknown7.95	7.95	118.1	ng	2780990	1	8.31	471078	20.0
n-Hexadecanoic acid	16.49	5.0	ng	173894	4	15.60	702770	20.0
Heneicosyl heptaf...	19.12	14.1	ng	463270	5	19.24	658509	20.0
Eicosane	19.53	2.6	ng	85024	5	19.24	658509	20.0
Octadecane	19.92	3.3	ng	109151	5	19.24	658509	20.0
Docosane, 7-hexyl-	20.30	4.7	ng	154941	6	21.01	654329	20.0
Heneicosane	20.69	5.7	ng	185034	6	21.01	654329	20.0
2-methyloctacosane	21.13	5.7	ng	186406	6	21.01	654329	20.0
E-15-Heptadecenal	21.70	2.9	ng	93914	6	21.01	654329	20.0
Heptadecane, 9-oc...	22.20	4.4	ng	144335	6	21.01	654329	20.0
Heptadecane, 3-me...	22.87	2.9	ng	95740	6	21.01	654329	20.0