

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
 Data File : BF113644.D
 Acq On : 8 Apr 2019 17:28
 Operator : JU/SJ
 Sample : K2264-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OPEN-WALL-EXCAVATION

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-BF040319.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	3.652	253	266	269	rBV	4846485	10483360	100.00%	19.138%
2	5.299	542	546	565	rBV	2165547	2366472	22.57%	4.320%
3	6.334	718	722	739	rBV	2266799	2537060	24.20%	4.632%
4	6.457	739	743	753	rVB	2578178	2509390	23.94%	4.581%
5	6.681	777	781	790	rBV	772827	689360	6.58%	1.258%
6	6.834	803	807	825	rBV	2050360	1941096	18.52%	3.544%
7	7.245	873	877	890	rBV	1643328	1588223	15.15%	2.899%
8	7.963	995	999	1017	rBV	947945	920791	8.78%	1.681%
9	9.039	1177	1182	1185	rBV	3175786	2935630	28.00%	5.359%
10	9.434	1245	1249	1257	rVB	328598	286133	2.73%	0.522%
11	9.710	1292	1296	1300	rBV2	1198497	1058715	10.10%	1.933%
12	9.934	1328	1334	1345	rBV3	48860	106484	1.02%	0.194%
13	10.504	1427	1431	1449	rBV	2090010	2105662	20.09%	3.844%
14	11.192	1545	1548	1550	rBV	1338947	1168943	11.15%	2.134%
15	11.216	1550	1552	1557	rVV	792861	682945	6.51%	1.247%
16	11.269	1557	1561	1565	rVB	106797	120238	1.15%	0.220%
17	11.733	1635	1640	1645	rVV2	922348	1080684	10.31%	1.973%
18	11.804	1650	1652	1655	rVV	138436	169334	1.62%	0.309%
19	12.404	1750	1754	1758	rBV	948411	854540	8.15%	1.560%
20	12.516	1767	1773	1778	rBV	968141	1313132	12.53%	2.397%
21	12.633	1787	1793	1796	rVB	756309	748434	7.14%	1.366%
22	12.775	1811	1817	1821	rBV	3289031	3370782	32.15%	6.154%
23	13.069	1862	1867	1870	rBV2	80989	114761	1.09%	0.210%
24	13.133	1870	1878	1880	rBV4	143861	308361	2.94%	0.563%
25	13.680	1967	1971	1977	rVB2	211297	285751	2.73%	0.522%
26	13.827	1982	1996	1999	rVV2	1337621	2138267	20.40%	3.904%
27	13.851	1999	2000	2009	rVB	354576	322358	3.07%	0.588%
28	13.992	2019	2024	2029	rVB	602861	635165	6.06%	1.160%
29	14.292	2070	2075	2081	rVB	1290820	1541858	14.71%	2.815%
30	14.592	2120	2126	2129	rBV	1841520	1918525	18.30%	3.502%
31	14.839	2165	2168	2174	rBV	226496	417011	3.98%	0.761%
32	14.898	2174	2178	2183	rVB	1746803	2094586	19.98%	3.824%
33	15.121	2212	2216	2220	rVB2	150461	194065	1.85%	0.354%
34	15.180	2222	2226	2230	rVB	192628	204651	1.95%	0.374%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.245	2231	2237	2250	rVB2	1785271	2772207	26.44%	5.061%
36	15.639	2298	2304	2310	rBV	909769	1375460	13.12%	2.511%
37	16.098	2375	2382	2392	rVB	404031	744016	7.10%	1.358%
38	16.621	2463	2471	2482	rVB2	145351	420012	4.01%	0.767%
39	17.015	2533	2538	2543	rBV2	60268	113316	1.08%	0.207%
40	17.251	2574	2578	2587	rVB3	59669	139421	1.33%	0.255%

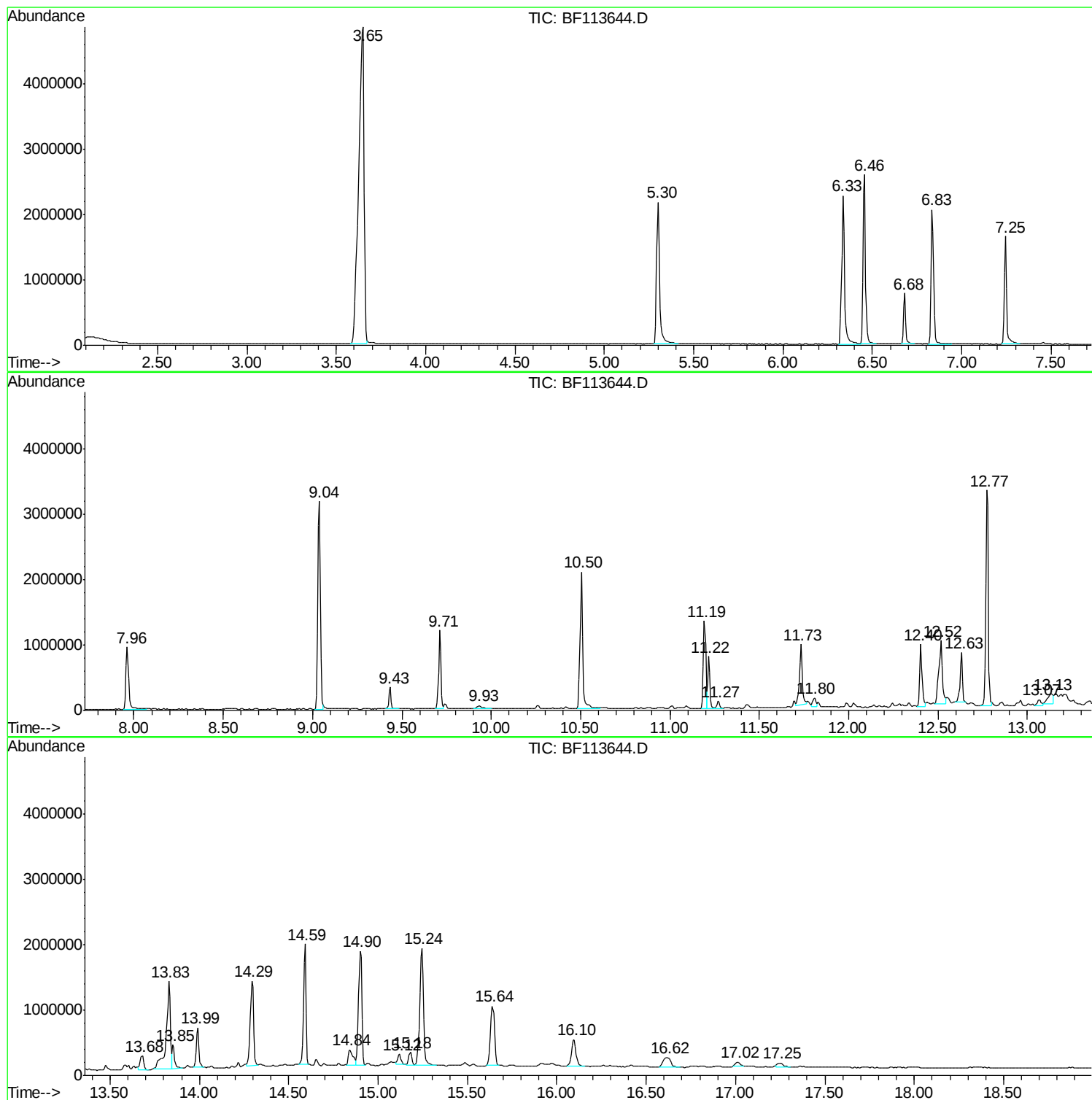
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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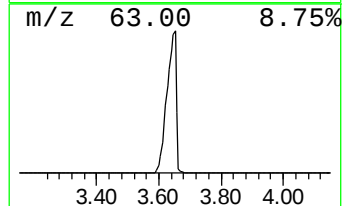
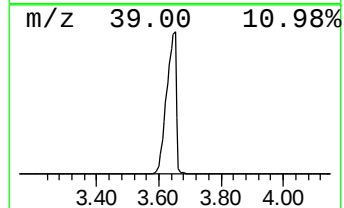
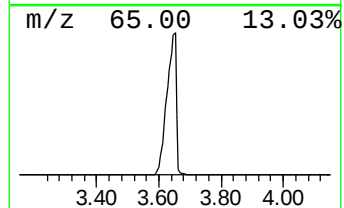
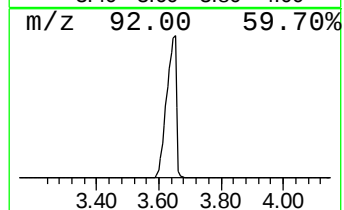
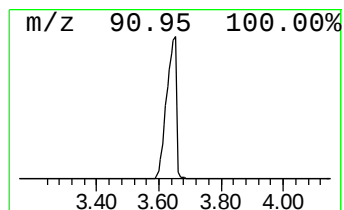
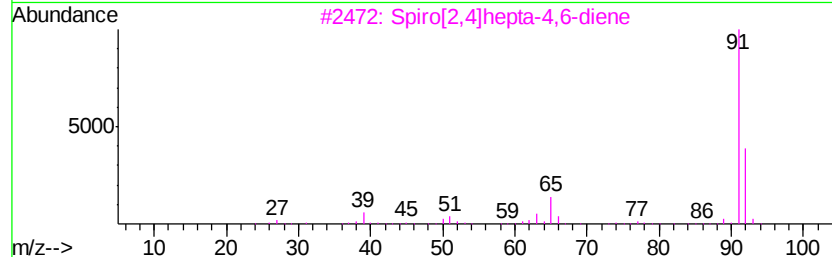
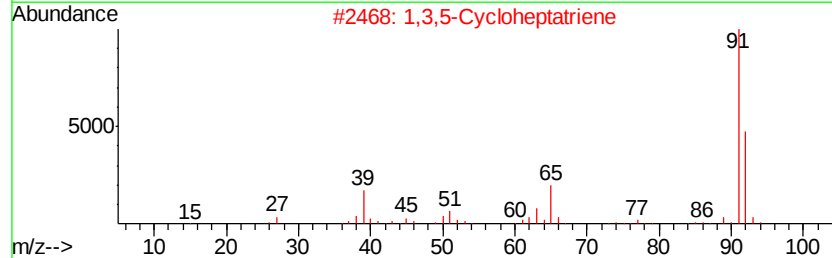
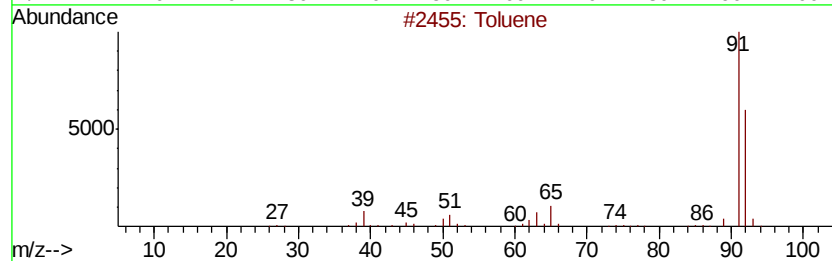
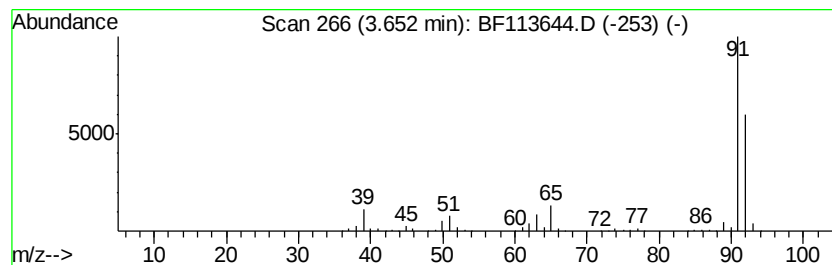
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1,3,5-Cycloheptatriene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.65	304.15 ng	10483400	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Toluene	92	C7H8	000108-88-3	95
2		1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	90
3		Spiro[2.4]hepta-4,6-diene	92	C7H8	000765-46-8	87
4		2,5-Norbornadiene	92	C7H8	000121-46-0	83
5		Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	59



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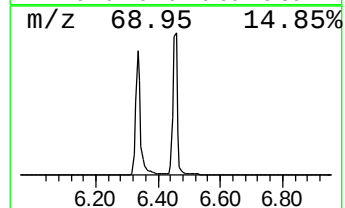
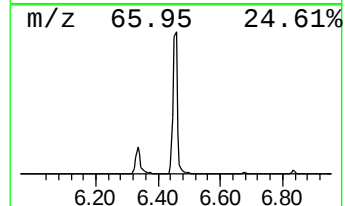
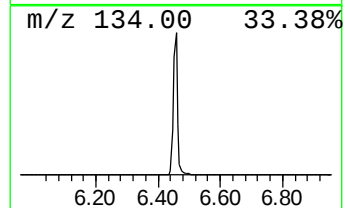
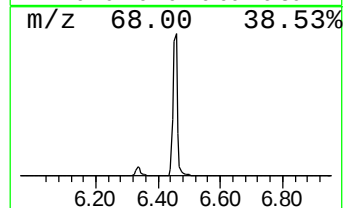
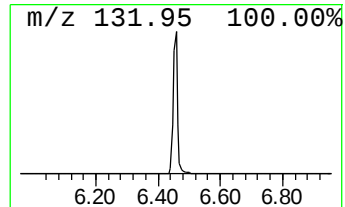
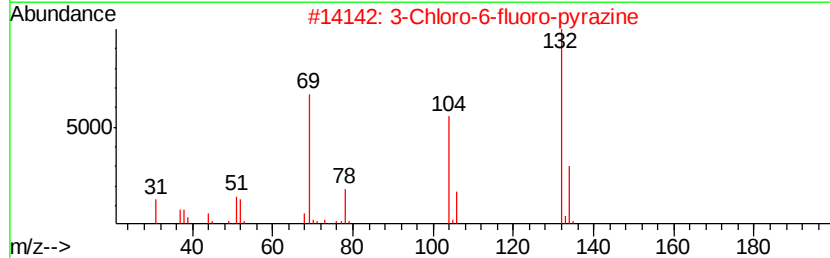
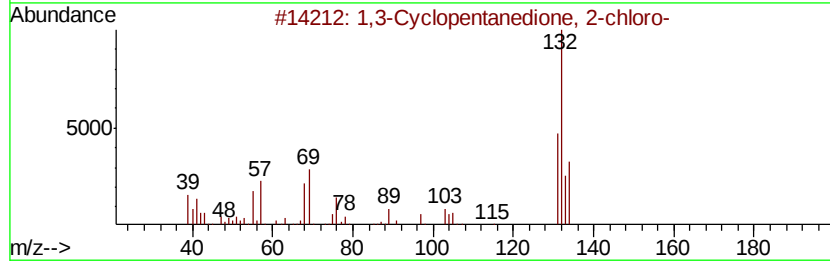
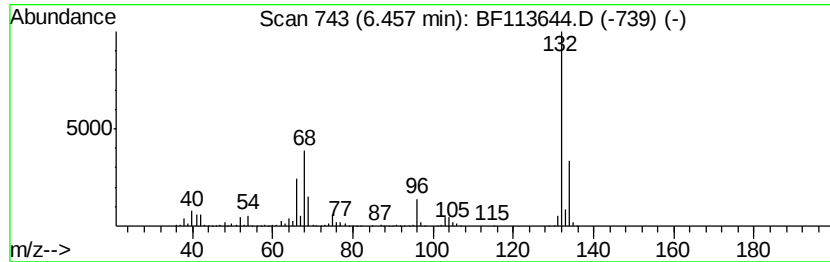
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.46 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	72.80 ng	2509390	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	37
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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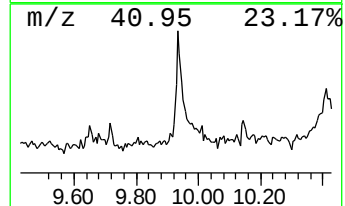
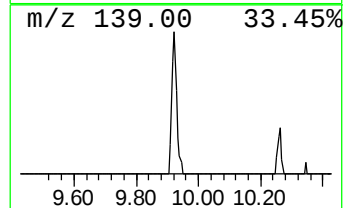
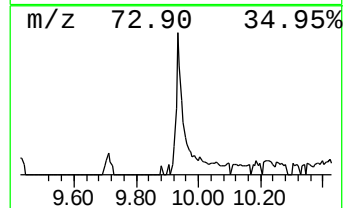
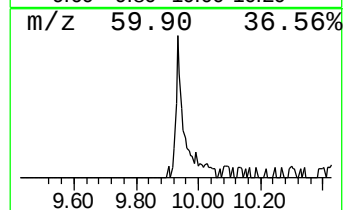
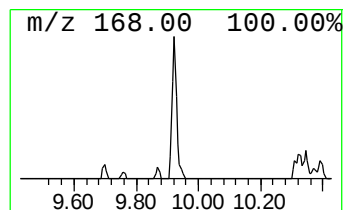
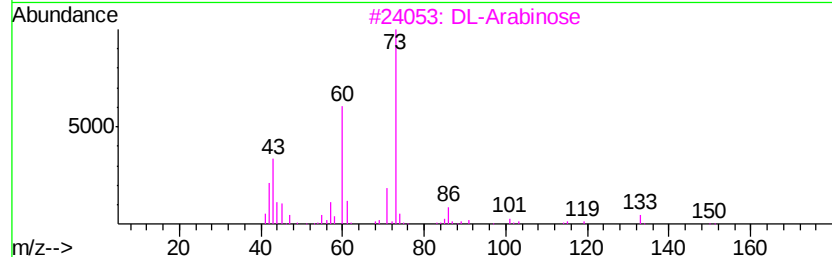
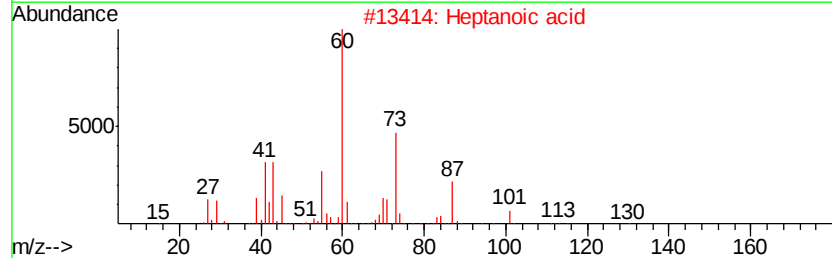
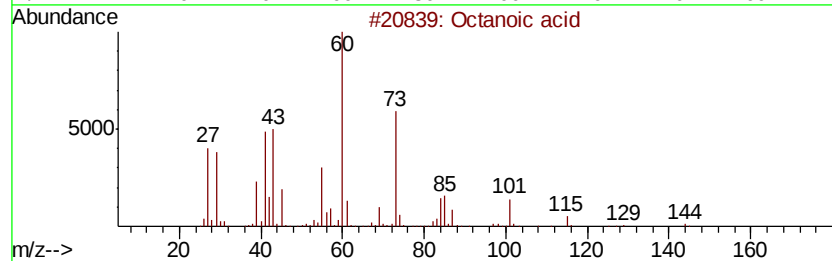
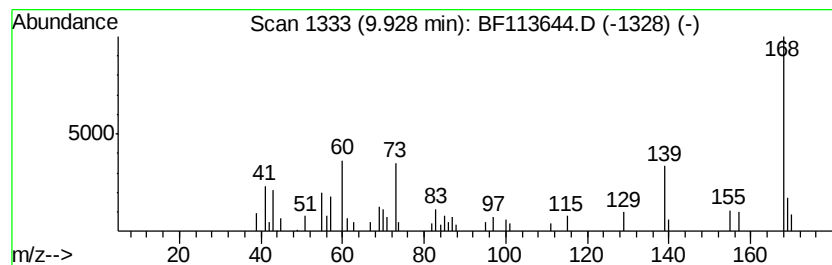
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown9.93 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	2.01 ng	106484	Acenaphthene-d10	9.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octanoic acid	144	C8H16O2	000124-07-2	27
2		Heptanoic acid	130	C7H14O2	000111-14-8	27
3		DL-Arabinose	150	C5H10O5	020235-19-2	27
4		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	27
5		d-Glucohexodialdose	178	C6H10O6	1000149-19-1	25



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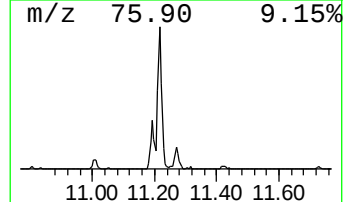
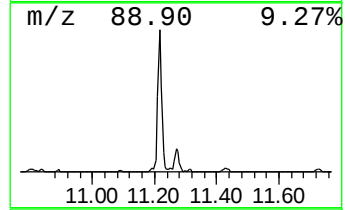
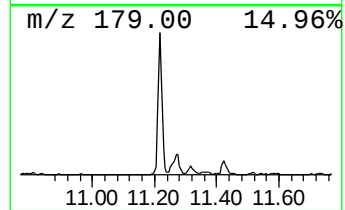
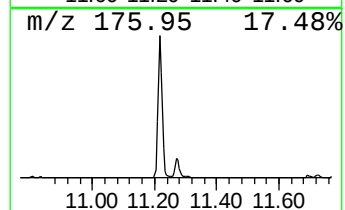
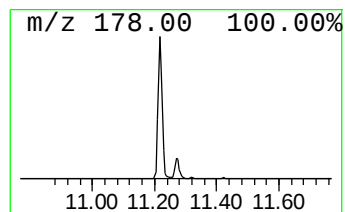
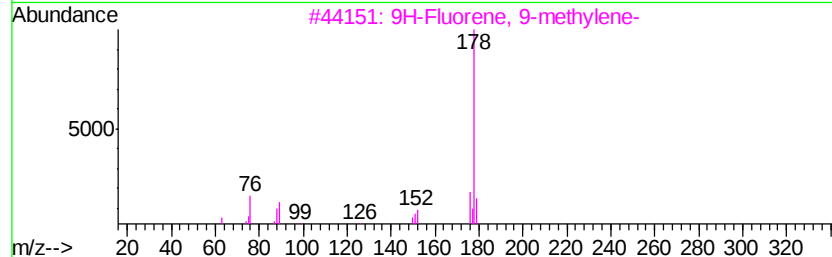
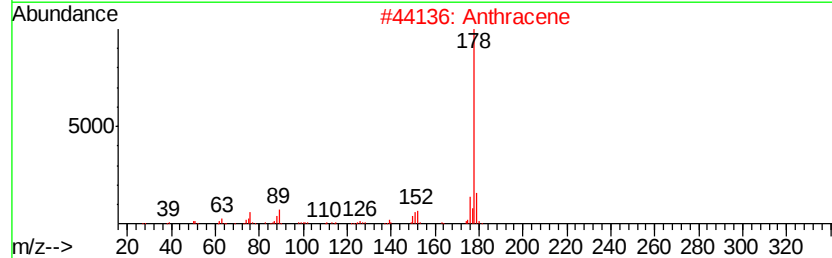
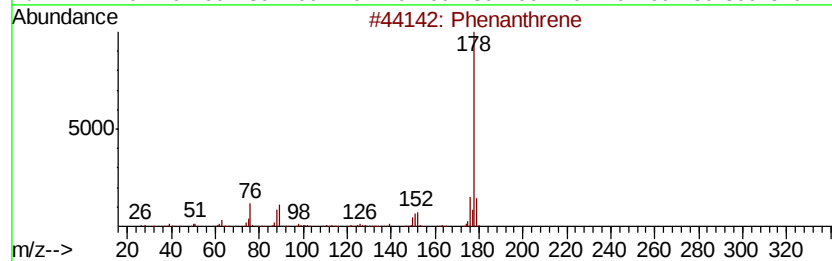
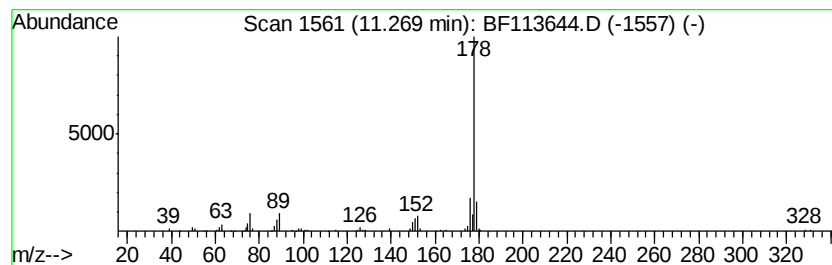
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9H-Fluorene, 9-methylene- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	2.06 ng	120238	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene	178	C14H10	000085-01-8	96
2		Anthracene	178	C14H10	000120-12-7	95
3		9H-Fluorene, 9-methvlene-	178	C14H10	004425-82-5	93
4		Diphenylacetylene	178	C14H10	000501-65-5	90
5		2-Cyclopropen-1-one, 2,3-diphenyl-	206	C15H10O	000886-38-4	83



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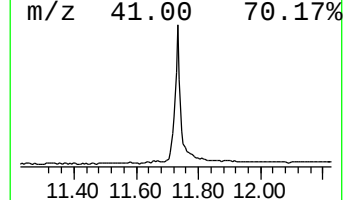
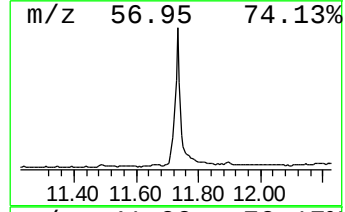
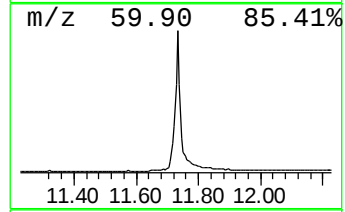
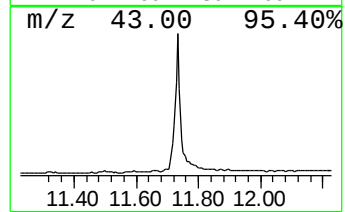
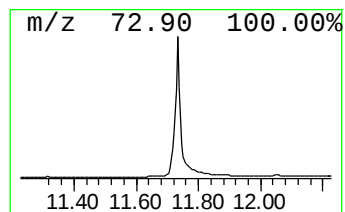
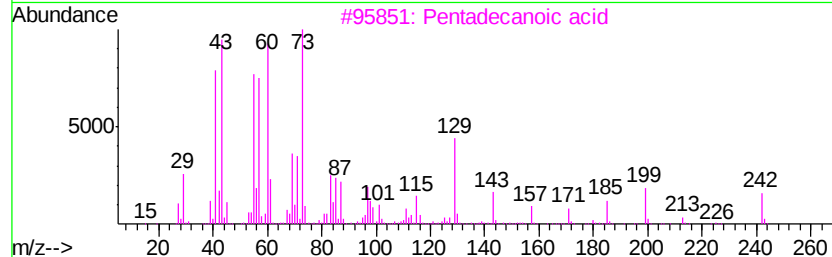
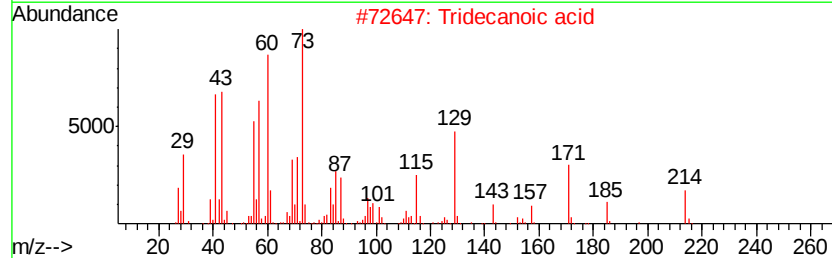
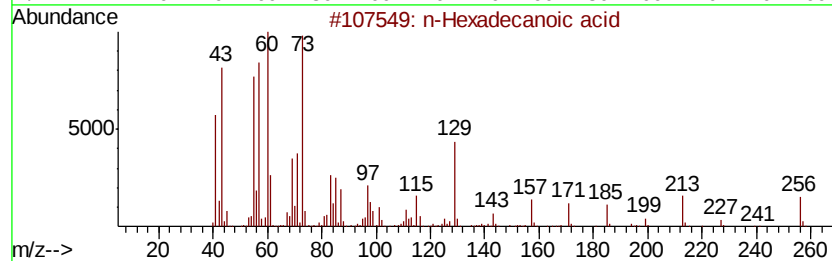
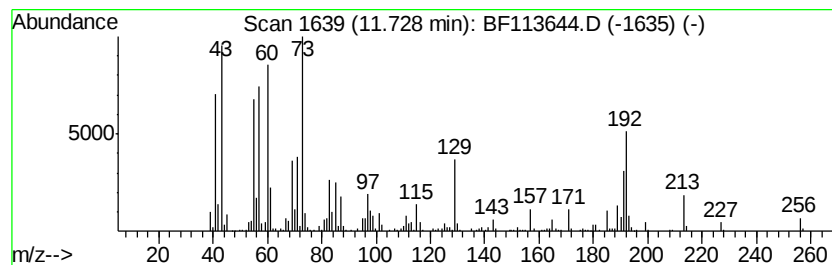
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.73	18.49 ng	1080680	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	93
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	76
5		n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
Operator : JU/SJ
Sample : K2264-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

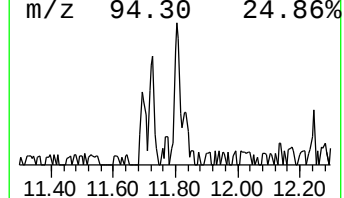
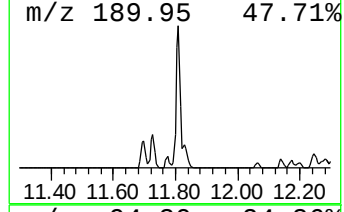
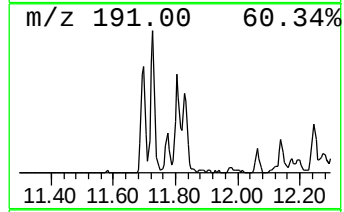
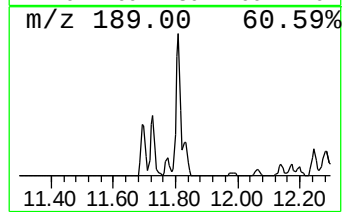
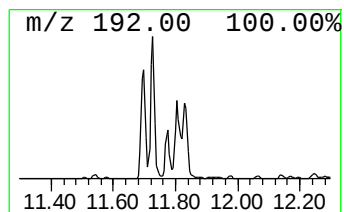
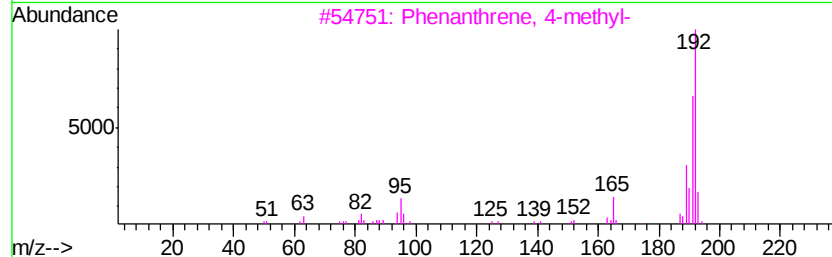
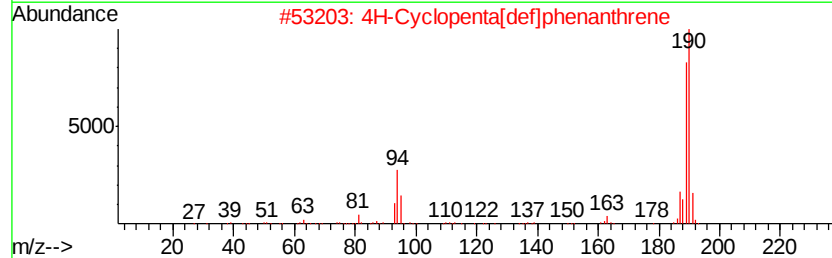
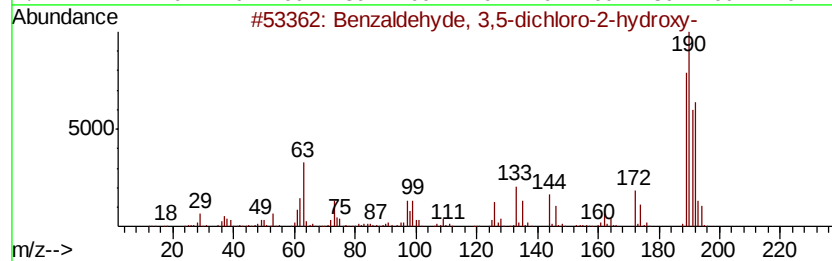
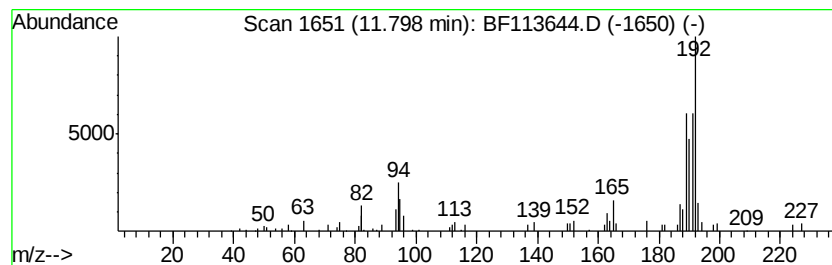
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ClientSampleId :
OPEN-WALL-EXCAVATION

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzaldehyde, 3,5-dichloro-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.80	2.90 ng	169334	Phenanthrene-d10	11.19	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
5	8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	42



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
Operator : JU/SJ
Sample : K2264-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

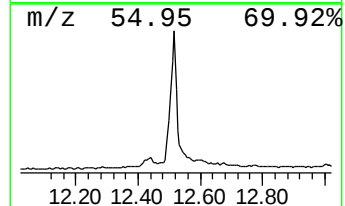
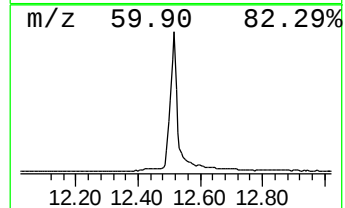
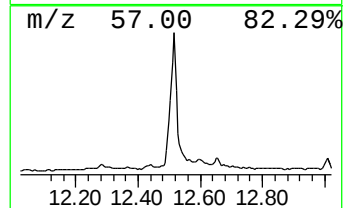
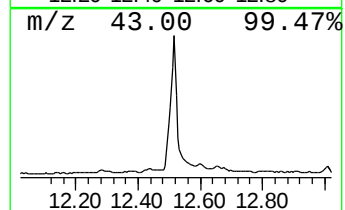
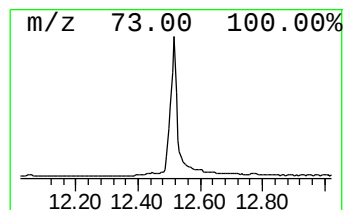
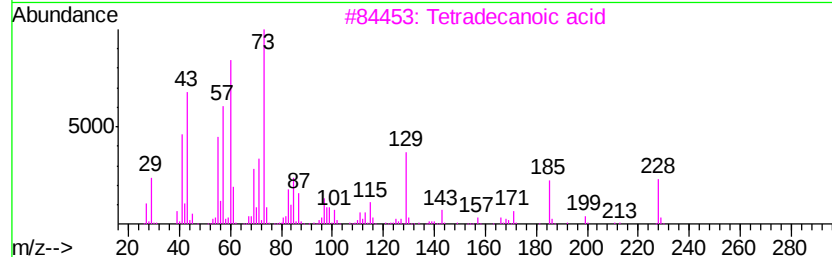
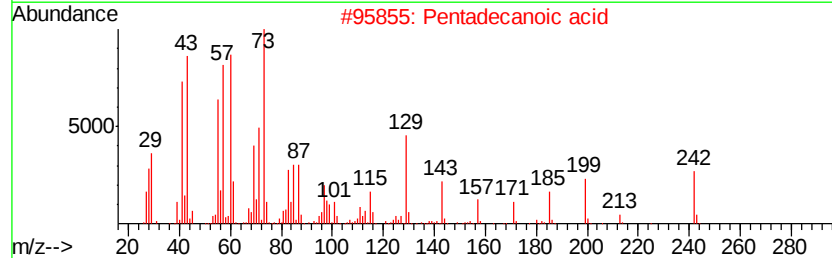
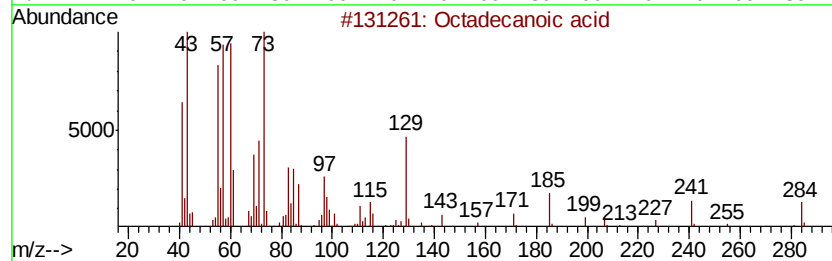
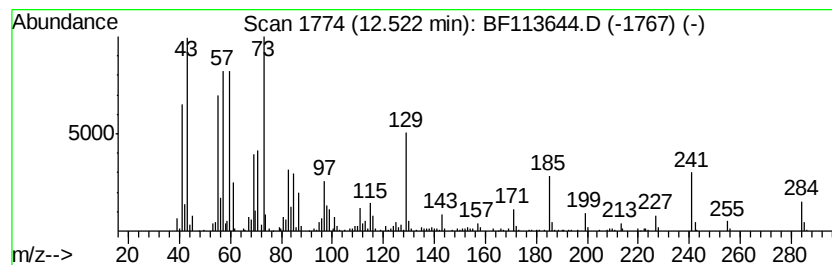
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ClientSampleId :
OPEN-WALL-EXCAVATION

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	12.28 ng	1313130	Chrysene-d12	13.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octadecanoic acid	284 C18H36O2	000057-11-4 99
2		Pentadecanoic acid	242 C15H30O2	001002-84-2 91
3		Tetradecanoic acid	228 C14H28O2	000544-63-8 87
4		n-Hexadecanoic acid	256 C16H32O2	000057-10-3 86
5		n-Decanoic acid	172 C10H20O2	000334-48-5 64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
Operator : JU/SJ
Sample : K2264-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

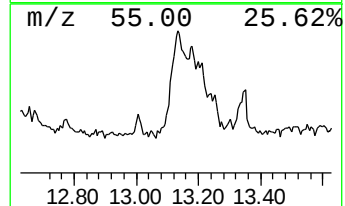
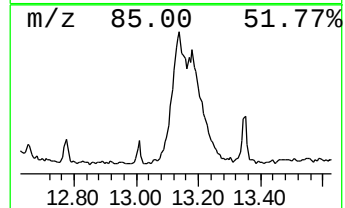
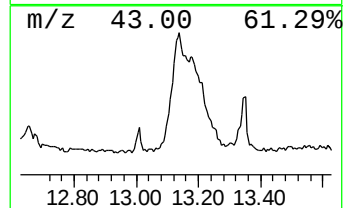
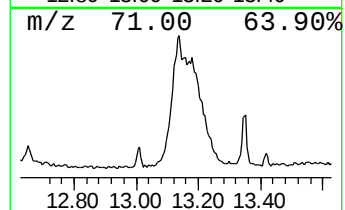
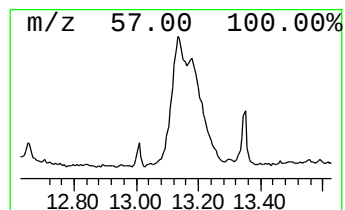
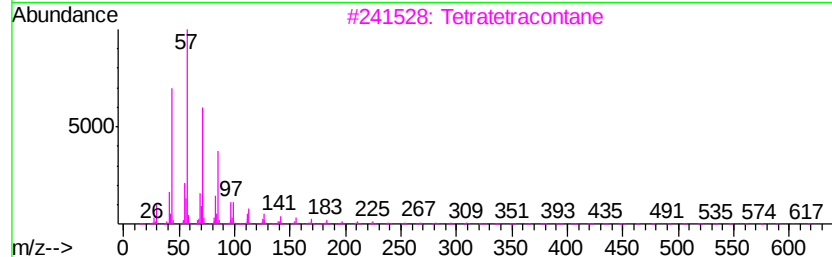
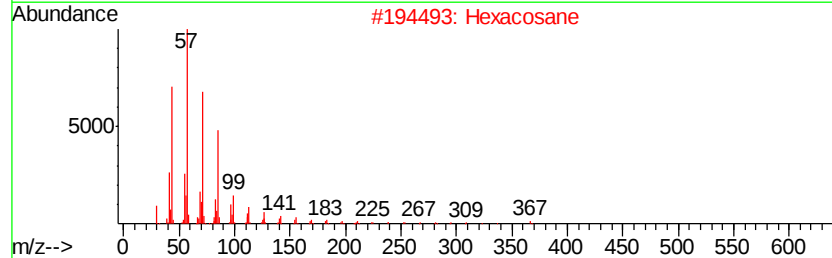
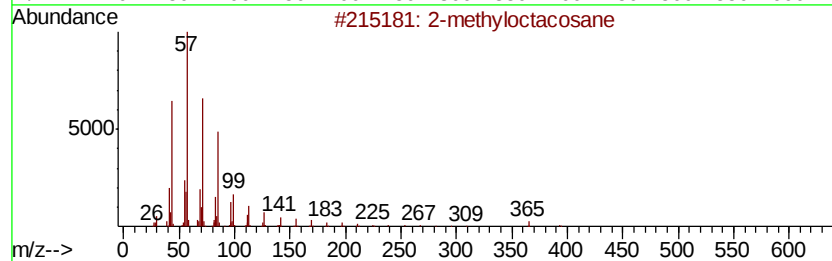
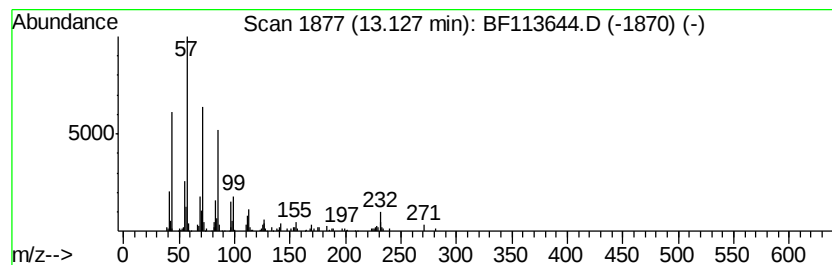
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-methyloctacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	2.88 ng	308361	Chrysene-d12	13.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-methyloctacosane	408 C29H60	1000376-72-8	90
2	Hexacosane	366 C26H54	000630-01-3	90
3	Tetratetracontane	619 C44H90	007098-22-8	90
4	Heptacosane, 1-chloro-	414 C27H55Cl	062016-79-9	87
5	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
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Sample : K2264-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

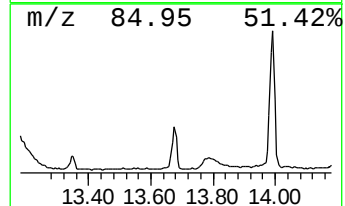
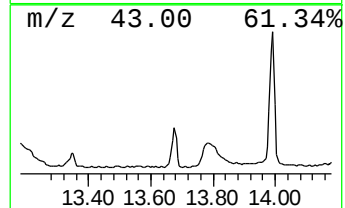
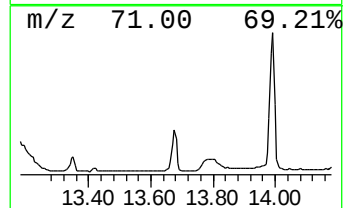
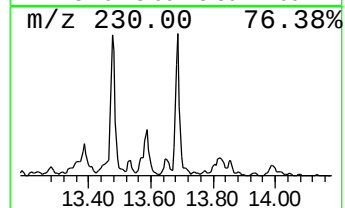
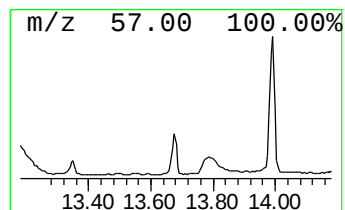
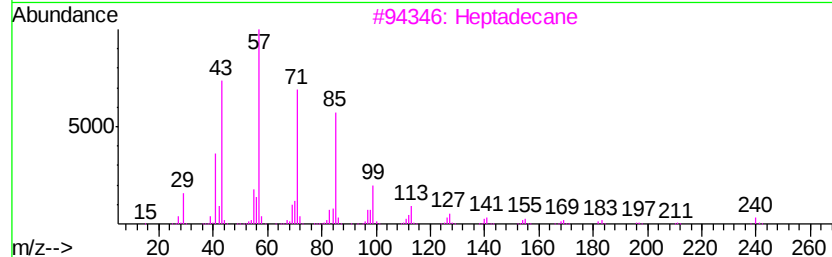
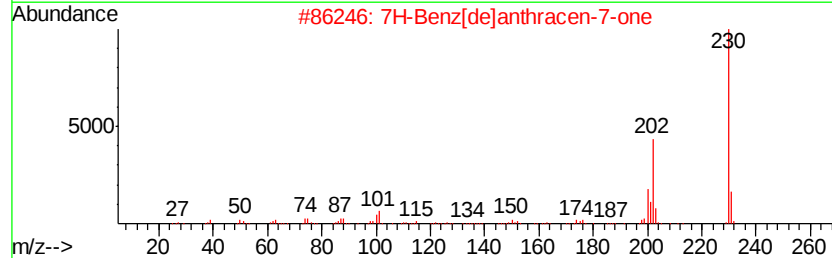
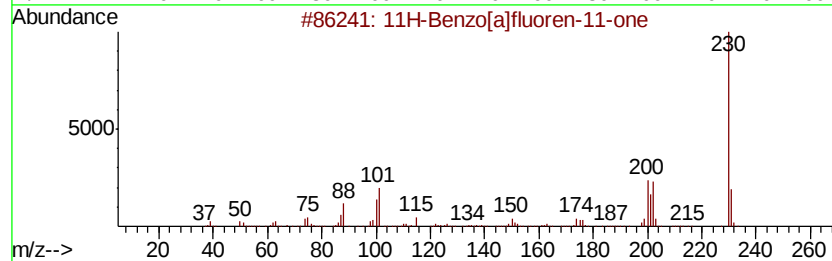
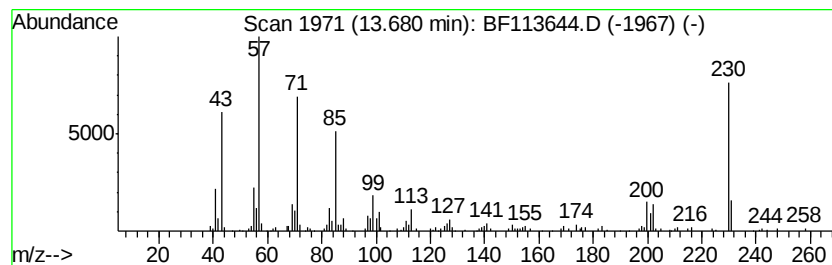
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluoren-11-one Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.68	2.67 ng	285751	Chrysene-d12		13.83	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	70
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	50
3		Heptadecane	240	C17H36	000629-78-7	42
4		Heneicosane	296	C21H44	000629-94-7	42
5		Hexadecane	226	C16H34	000544-76-3	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
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Sample : K2264-01
Misc :
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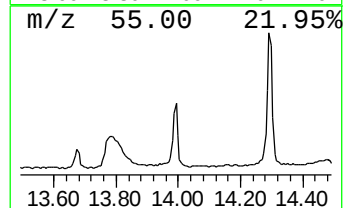
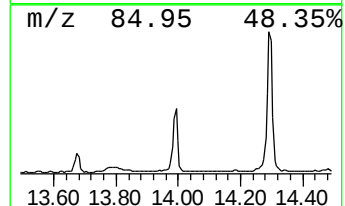
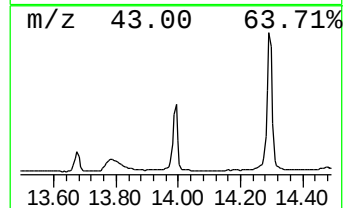
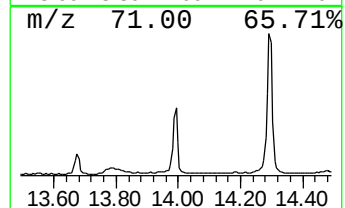
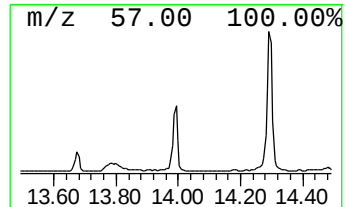
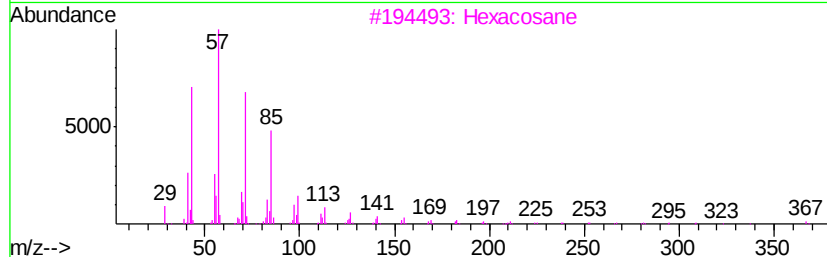
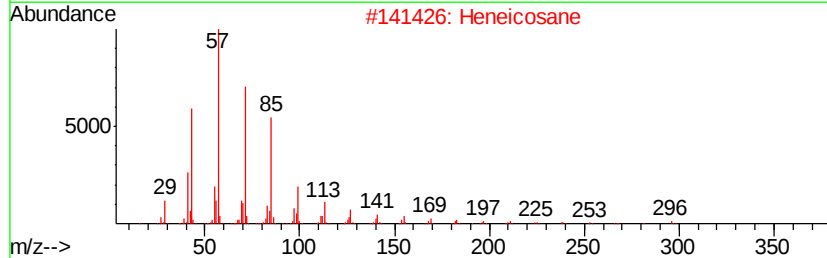
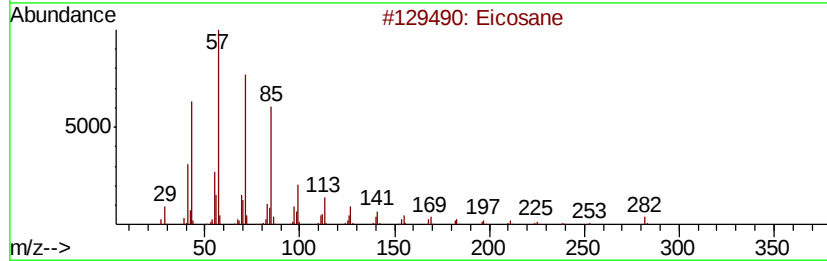
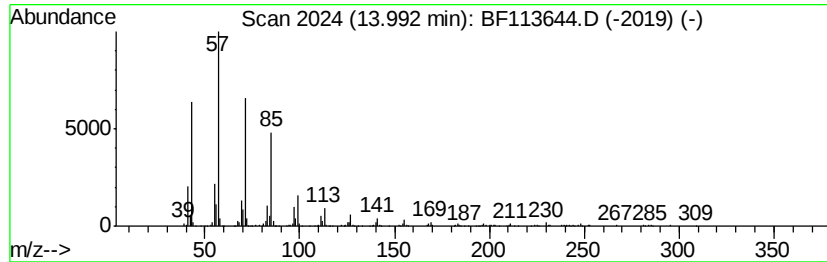
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	5.94 ng	635165	Chrysene-d12	13.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	97
2	Heneicosane	296 C21H44	000629-94-7	91
3	Hexacosane	366 C26H54	000630-01-3	91
4	Octacosane	394 C28H58	000630-02-4	91
5	Nonadecane, 2-methyl-	282 C20H42	001560-86-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
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Operator : JU/SJ
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OPEN-WALL-EXCAVATION

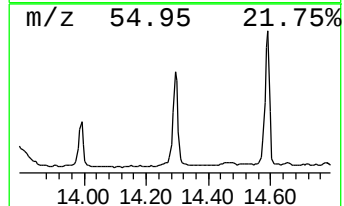
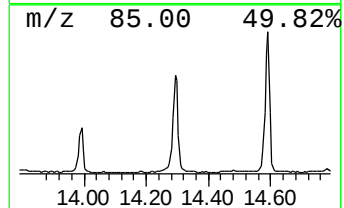
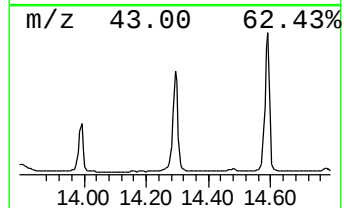
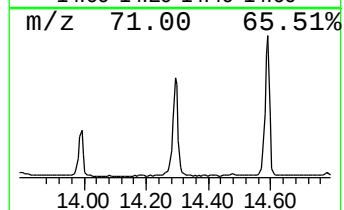
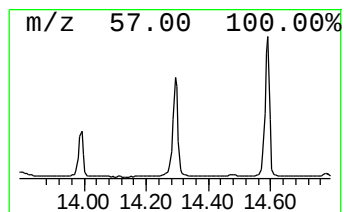
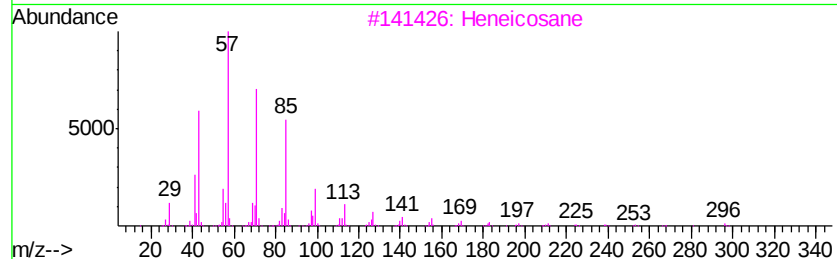
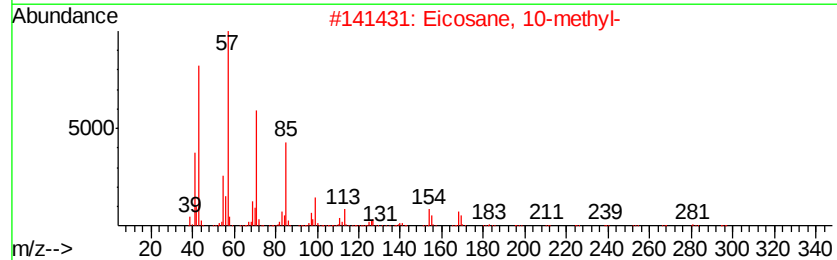
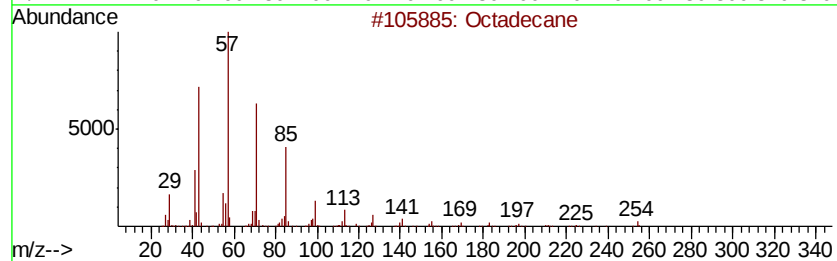
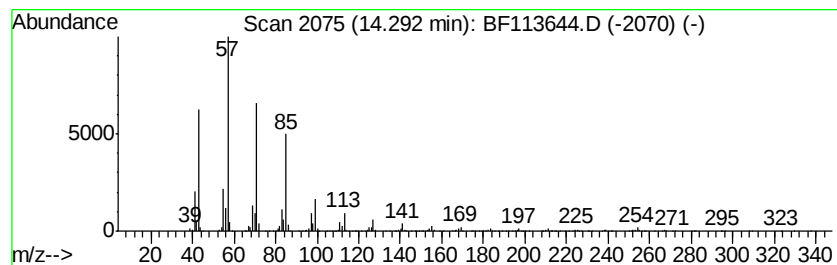
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	14.42 ng	1541860	Chrysene-d12	13.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	96
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
Operator : JU/SJ
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ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OPEN-WALL-EXCAVATION

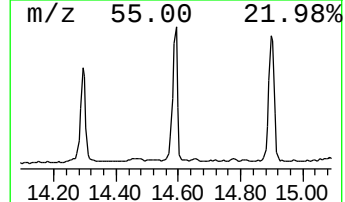
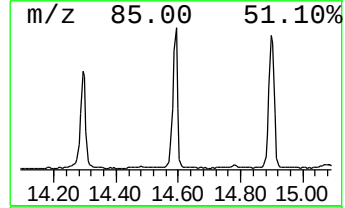
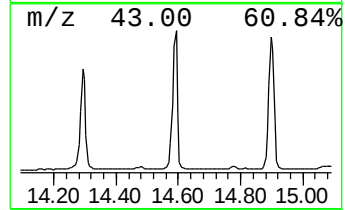
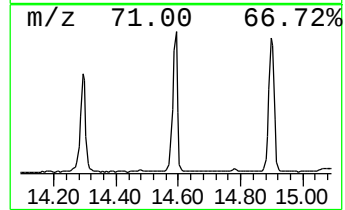
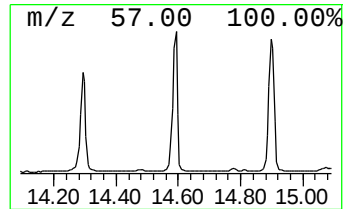
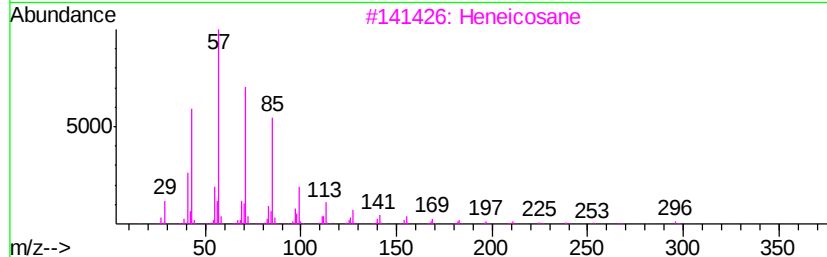
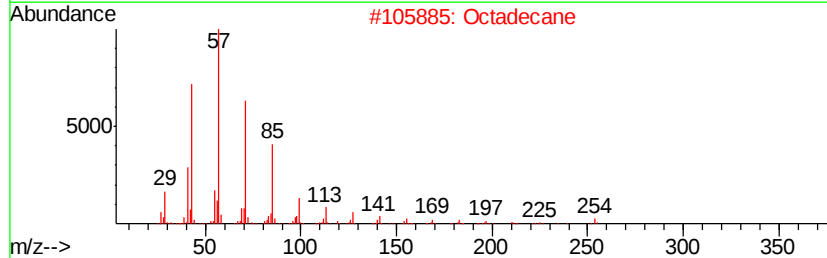
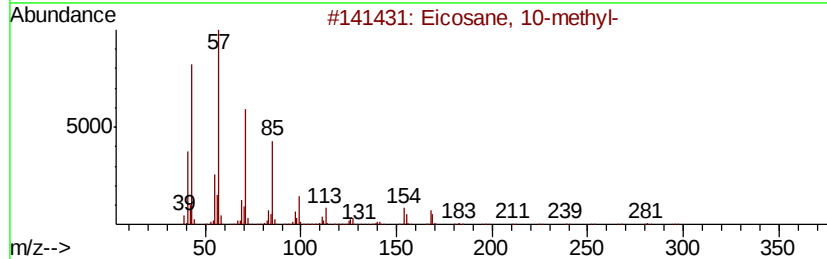
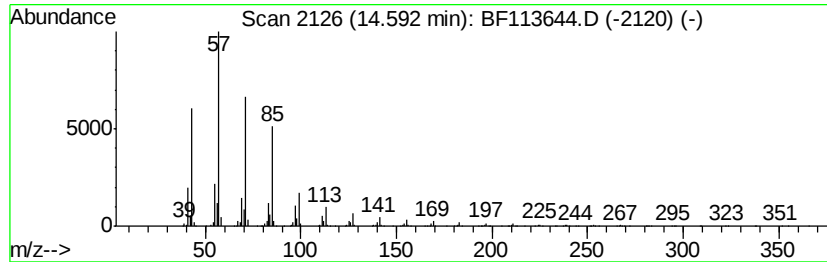
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane, 10-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	13.84 ng	1918530	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2		Octadecane	254	C18H38	000593-45-3	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
Operator : JU/SJ
Sample : K2264-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OPEN-WALL-EXCAVATION

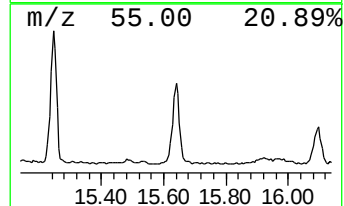
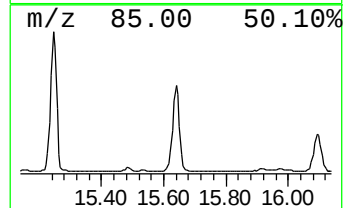
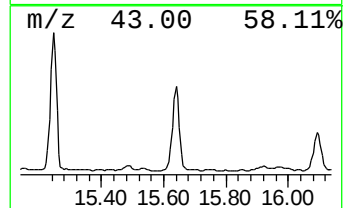
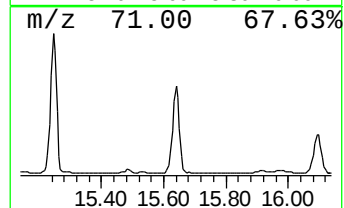
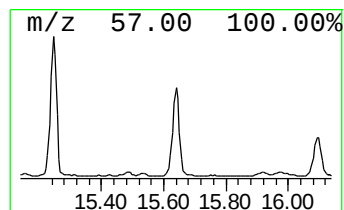
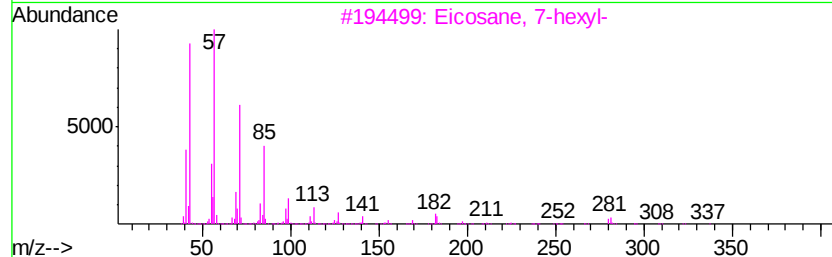
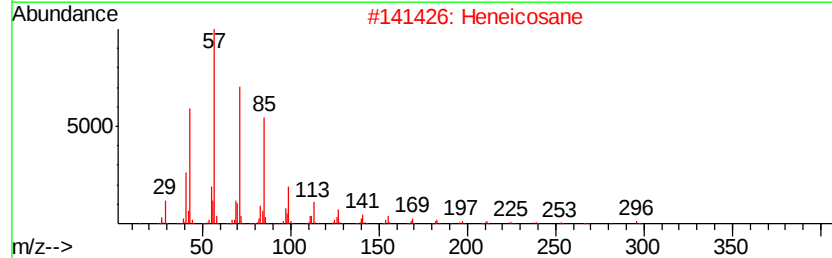
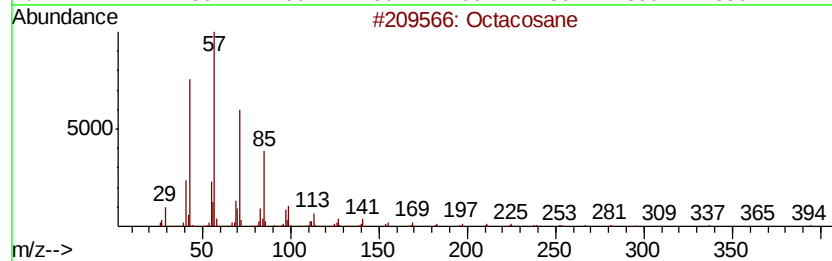
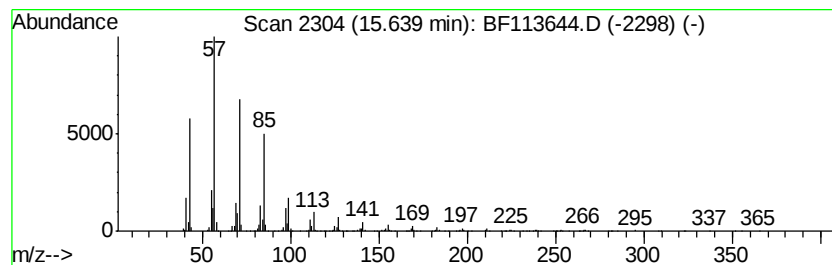
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	9.92 ng	1375460	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
4		Docosane	310	C22H46	000629-97-0	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040819\
Data File : BF113644.D
Acq On : 8 Apr 2019 17:28
Operator : JU/SJ
Sample : K2264-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OPEN-WALL-EXCAVATION

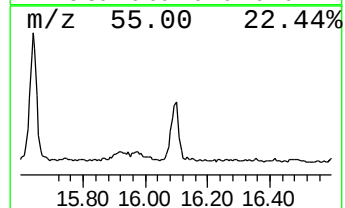
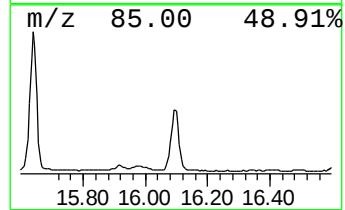
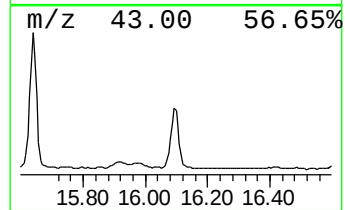
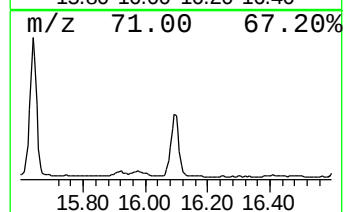
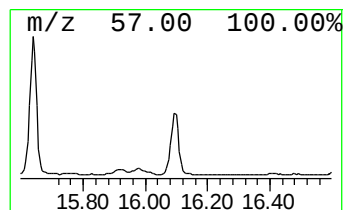
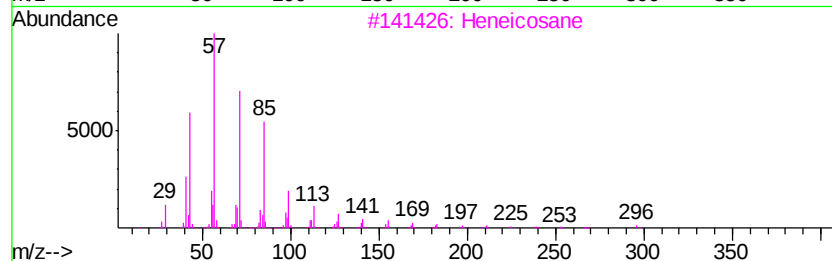
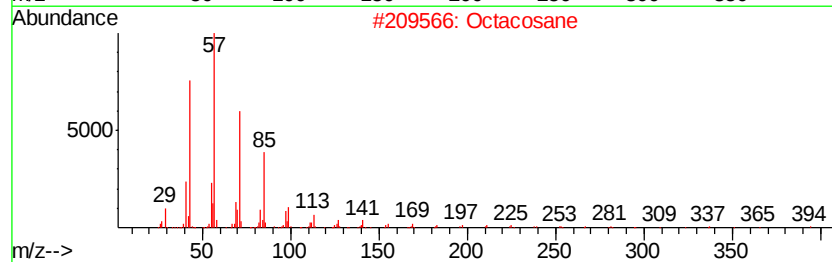
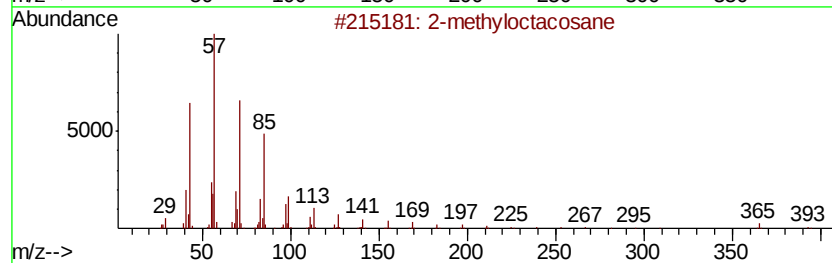
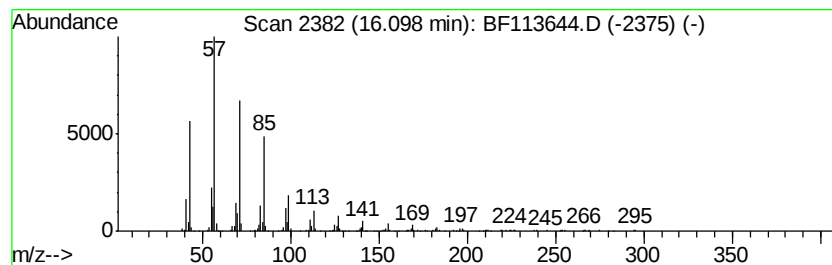
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2-Bromotetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	5.37 ng	744016	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5		2-Bromotetradecane	276	C14H29Br	074036-95-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040819\
 Data File : BF113644.D
 Acq On : 8 Apr 2019 17:28
 Operator : JU/SJ
 Sample : K2264-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OPEN-WALL-EXCAVATION

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040319.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
1,3,5-Cycloheptat...	3.65	304.1	ng	10483400	1	6.68	689360	20.0
unknown6.46	6.46	72.8	ng	2509390	1	6.68	689360	20.0
unknown9.93	9.93	2.0	ng	106484	3	9.71	1058720	20.0
9H-Fluorene, 9-me...	11.27	2.1	ng	120238	4	11.19	1168940	20.0
n-Hexadecanoic acid	11.73	18.5	ng	1080680	4	11.19	1168940	20.0
Benzaldehyde, 3,5...	11.80	2.9	ng	169334	4	11.19	1168940	20.0
Octadecanoic acid	12.52	12.3	ng	1313130	5	13.83	2138270	20.0
2-methyloctacosane	13.13	2.9	ng	308361	5	13.83	2138270	20.0
11H-Benzo[a]fluor...	13.68	2.7	ng	285751	5	13.83	2138270	20.0
Eicosane	13.99	5.9	ng	635165	5	13.83	2138270	20.0
Octadecane	14.29	14.4	ng	1541860	5	13.83	2138270	20.0
Eicosane, 10-methyl-	14.59	13.8	ng	1918530	6	15.24	2772210	20.0
Octacosane	15.64	9.9	ng	1375460	6	15.24	2772210	20.0
2-Bromotetradecane	16.10	5.4	ng	744016	6	15.24	2772210	20.0