

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030920\
 Data File : BP002136.D
 Acq On : 10 Mar 2020 00:02
 Operator : CG/JU
 Sample : L1834-16
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0DK9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % 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\SOM-EPA-BP022720MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	23	rVB	972029	1338422	13.29%	1.314%
2	3.164	43	47	54	rBV	59519	84676	0.84%	0.083%
3	3.440	89	94	101	rBV6	10140	26294	0.26%	0.026%
4	3.628	122	126	129	rBV2	13145	18706	0.19%	0.018%
5	3.805	153	156	162	rVB4	8891	12111	0.12%	0.012%
6	3.899	167	172	176	rVB2	11186	18674	0.19%	0.018%
7	4.111	205	208	212	rBV5	8379	11809	0.12%	0.012%
8	4.793	320	324	329	rBV4	5990	12127	0.12%	0.012%
9	4.993	350	358	370	rBV	3788818	5873924	58.32%	5.767%
10	5.334	412	416	421	rBV2	5643	12286	0.12%	0.012%
11	6.817	661	668	682	rVB	254466	466318	4.63%	0.458%
12	6.975	686	695	704	rBV	945050	1496599	14.86%	1.469%
13	7.175	722	729	740	rBV	1007665	1699139	16.87%	1.668%
14	7.528	783	789	800	rVB	384768	601897	5.98%	0.591%
15	7.634	800	807	810	rBV	982762	1662472	16.51%	1.632%
16	7.669	810	813	829	rVB	1984288	3031754	30.10%	2.976%
17	7.987	859	867	879	rVB	1483044	2403456	23.86%	2.360%
18	8.340	917	927	934	rBV	794706	1292143	12.83%	1.269%
19	8.393	934	936	945	rVB2	29160	57006	0.57%	0.056%
20	8.775	993	1001	1010	rBV	1073020	1820921	18.08%	1.788%
21	9.263	1079	1084	1089	rBV2	32924	54802	0.54%	0.054%
22	9.499	1117	1124	1133	rBV	875373	1458523	14.48%	1.432%
23	10.034	1207	1215	1224	rBV	1561866	2661215	26.42%	2.613%
24	10.193	1238	1242	1248	rVB4	10859	19432	0.19%	0.019%
25	10.316	1254	1263	1272	rBV	71085	163118	1.62%	0.160%
26	10.405	1272	1278	1286	rVB	1354150	2311104	22.95%	2.269%
27	10.552	1297	1303	1309	rBV2	44422	82750	0.82%	0.081%
28	11.757	1503	1508	1512	rBV2	19340	27460	0.27%	0.027%
29	11.910	1530	1534	1540	rVB2	15764	27106	0.27%	0.027%
30	12.005	1545	1550	1556	rBV2	30246	58038	0.58%	0.057%
31	12.781	1680	1682	1687	rBV2	7594	11572	0.11%	0.011%
32	12.840	1688	1692	1698	rVB4	13828	23491	0.23%	0.023%
33	13.075	1728	1732	1736	rBV4	8645	13842	0.14%	0.014%
34	13.293	1766	1769	1777	rVB	16970	27132	0.27%	0.027%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	13.687	1829	1836	1841	rBV	3038324	4386917	43.56%	4.307%
36	13.728	1841	1843	1851	rVB	298718	411919	4.09%	0.404%
37	13.957	1875	1882	1898	rBV	3527603	5364904	53.27%	5.267%
38	14.269	1929	1935	1942	rVV	2197154	3317664	32.94%	3.257%
39	14.481	1965	1971	1988	rBV2	132104	307600	3.05%	0.302%
40	14.798	2022	2025	2033	rVB2	20129	34230	0.34%	0.034%
41	15.269	2098	2105	2118	rBV	4751579	6981291	69.32%	6.854%
42	15.387	2119	2125	2138	rVV	1110656	1608740	15.97%	1.579%
43	15.510	2142	2146	2151	rVB3	56624	83563	0.83%	0.082%
44	16.057	2236	2239	2245	rVB3	22343	31575	0.31%	0.031%
45	17.022	2397	2403	2413	rBV	2832363	3986252	39.58%	3.913%
46	17.122	2413	2420	2432	rBV	5352288	7910045	78.54%	7.766%
47	17.287	2444	2448	2453	rBV3	31491	42526	0.42%	0.042%
48	17.428	2469	2472	2478	rVB	66514	83024	0.82%	0.082%
49	17.945	2555	2560	2575	rBV3	242735	456259	4.53%	0.448%
50	18.898	2719	2722	2726	rBV2	48530	68069	0.68%	0.067%
51	19.010	2738	2741	2746	rVB	75162	97259	0.97%	0.095%
52	19.110	2755	2758	2764	rVB2	40270	54218	0.54%	0.053%
53	19.181	2767	2770	2777	rVV	76677	110389	1.10%	0.108%
54	19.257	2779	2783	2787	rVB2	76261	110669	1.10%	0.109%
55	19.369	2796	2802	2808	rBV	6804104	9358845	92.92%	9.188%
56	19.928	2894	2897	2900	rVB2	47013	47123	0.47%	0.046%
57	20.410	2977	2979	2983	rVB2	47457	45753	0.45%	0.045%
58	20.857	3051	3055	3062	rBV	499493	836760	8.31%	0.821%
59	21.116	3094	3099	3109	rBV	3669691	4732913	46.99%	4.646%
60	21.292	3125	3129	3133	rBV	242229	273918	2.72%	0.269%
61	21.722	3198	3202	3216	rBV	479445	678340	6.74%	0.666%
62	22.180	3275	3280	3286	rBV	584284	769693	7.64%	0.756%
63	22.686	3361	3366	3380	rVB	838429	1266257	12.57%	1.243%
64	23.180	3442	3450	3457	rBV	5445346	10071511	100.00%	9.888%
65	23.251	3457	3462	3467	rVV	675619	1153029	11.45%	1.132%
66	23.322	3467	3474	3490	rVB	2547631	4972930	49.38%	4.882%
67	23.898	3566	3572	3580	rBV	545736	1062034	10.54%	1.043%
68	24.351	3644	3649	3665	rVB	146453	419243	4.16%	0.412%
69	24.639	3688	3698	3705	rBV2	411540	1448098	14.38%	1.422%
70	26.368	3987	3992	4001	rVB7	161114	396072	3.93%	0.389%

LSC Area Percent Report

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Title : SVOA CALIBRATION

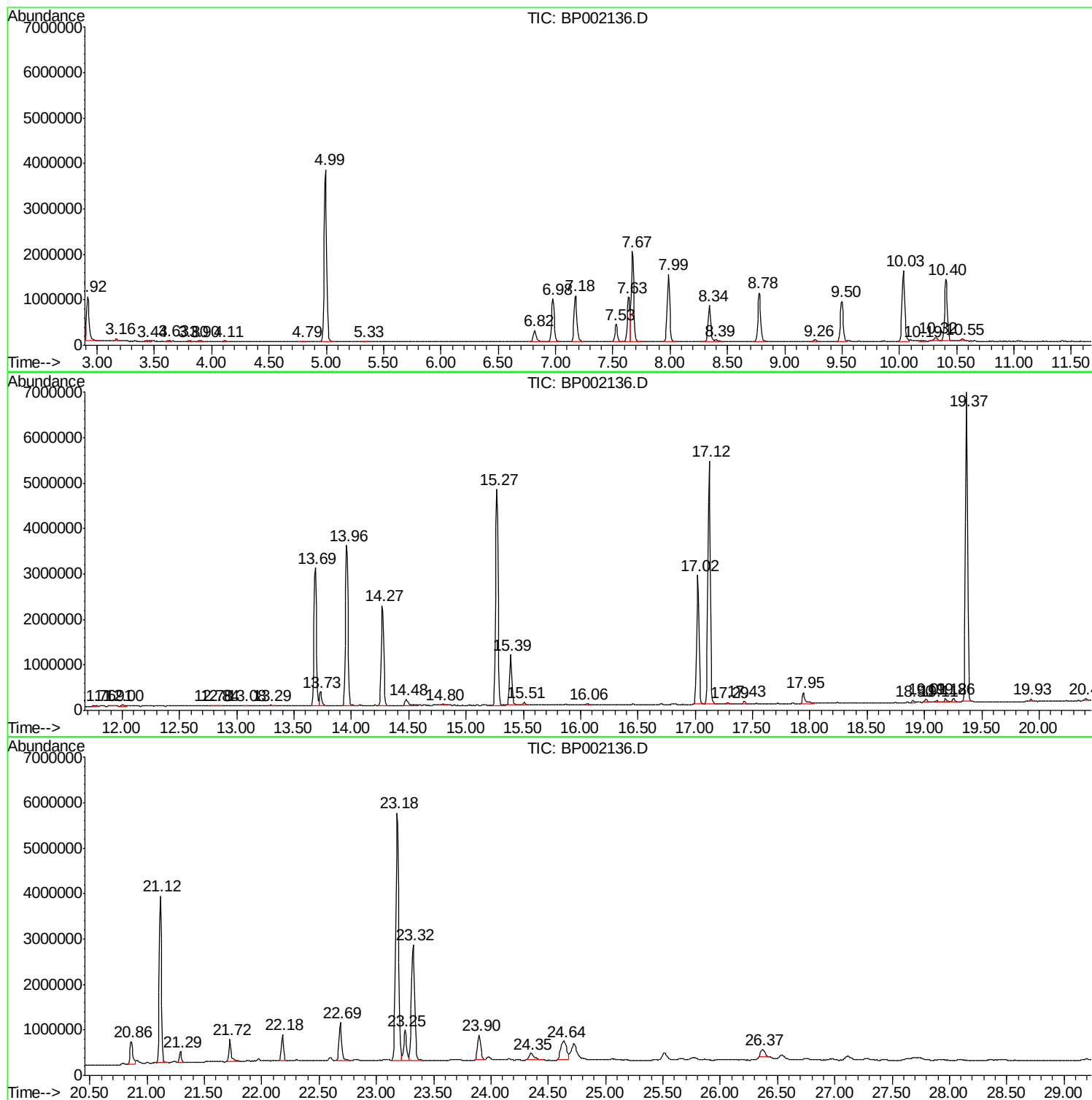
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
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Sample : L1834-16
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Instrument :
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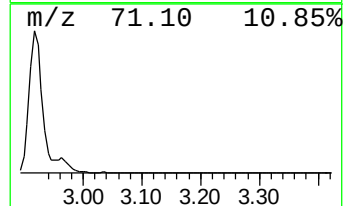
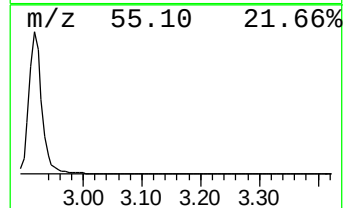
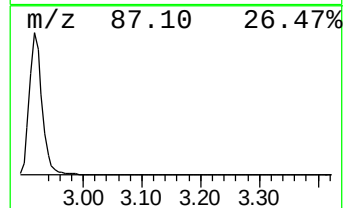
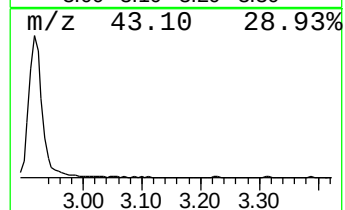
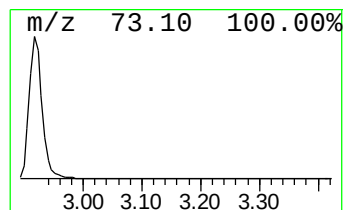
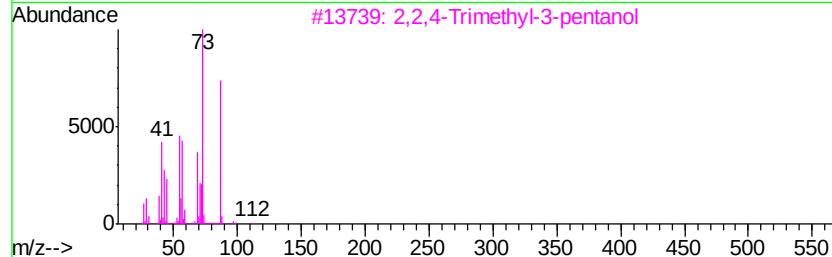
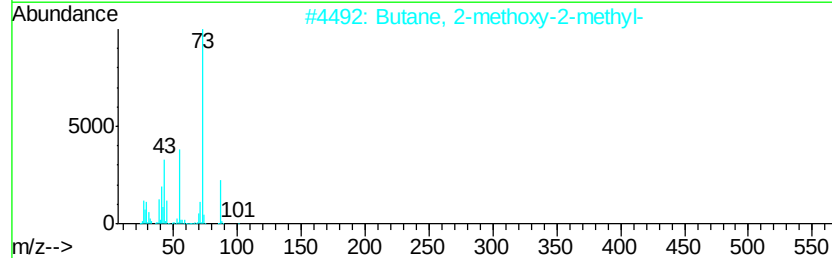
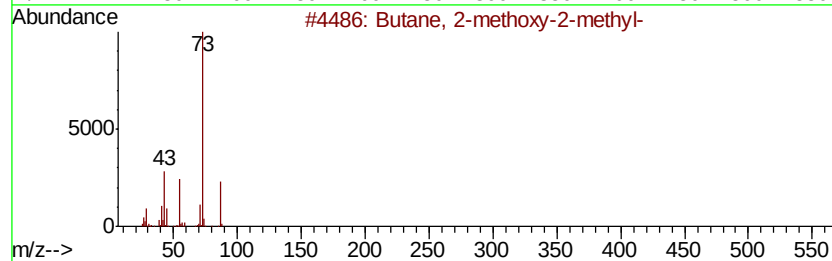
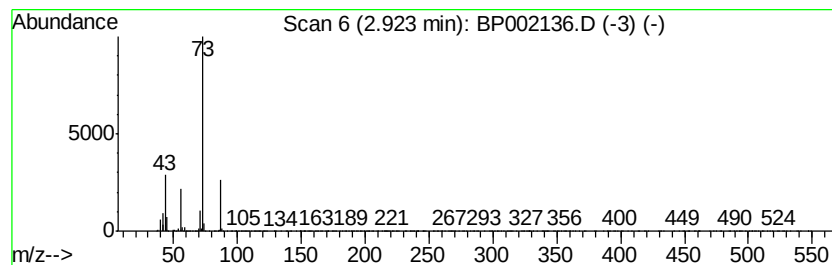
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	16.10 ng/ul	1338420	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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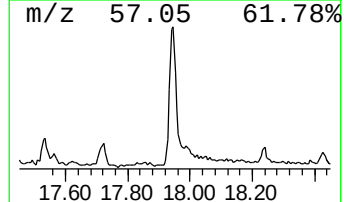
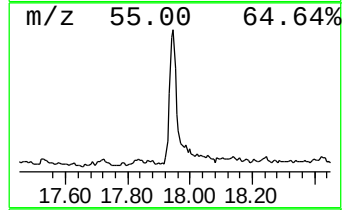
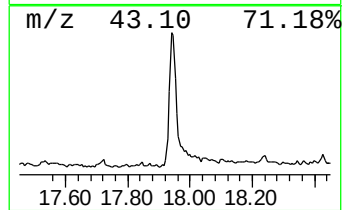
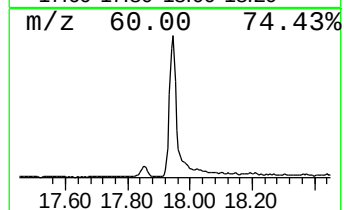
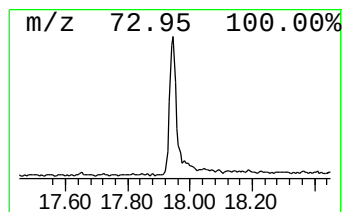
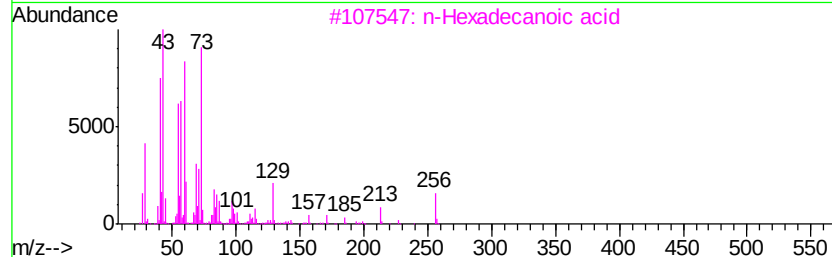
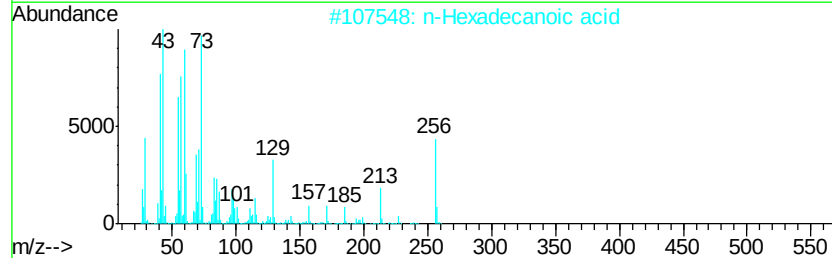
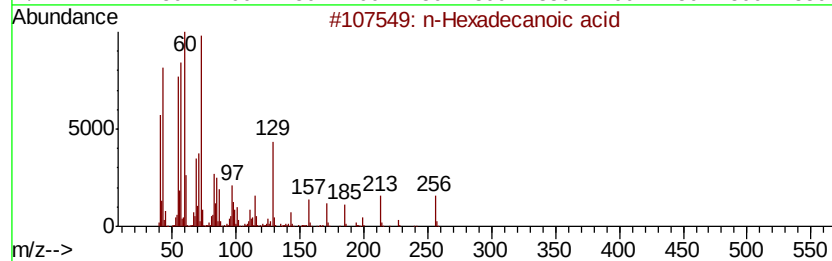
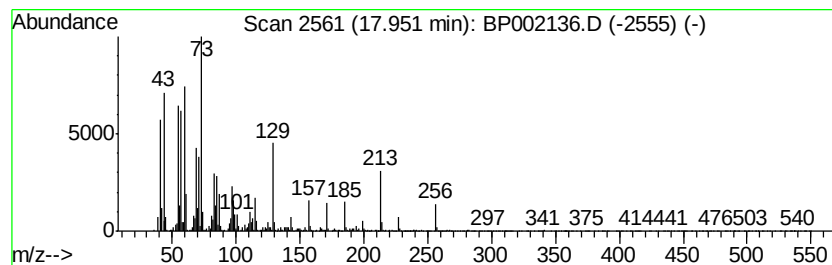
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	2.29 ng/ul	456259	Phenanthrene-d10	17.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Tridecanoic acid	214	C13H26O2	000638-53-9	83



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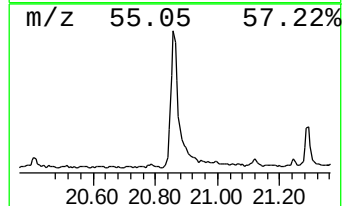
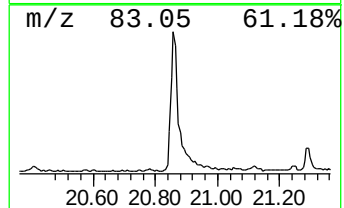
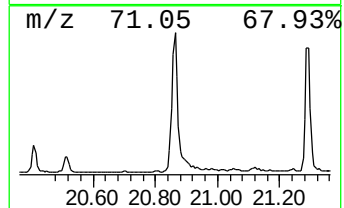
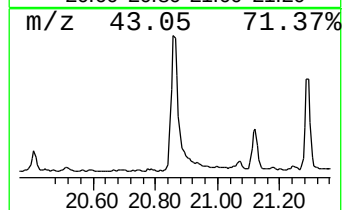
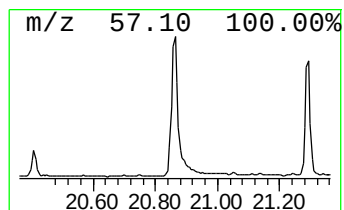
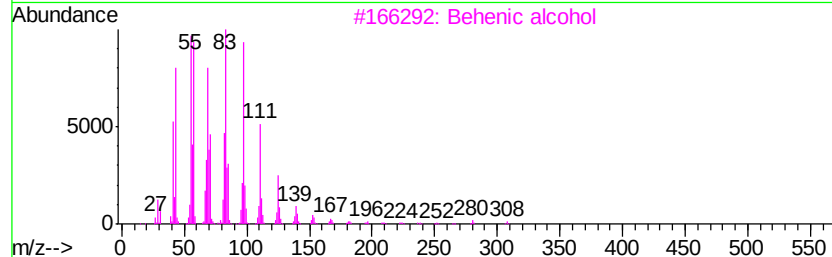
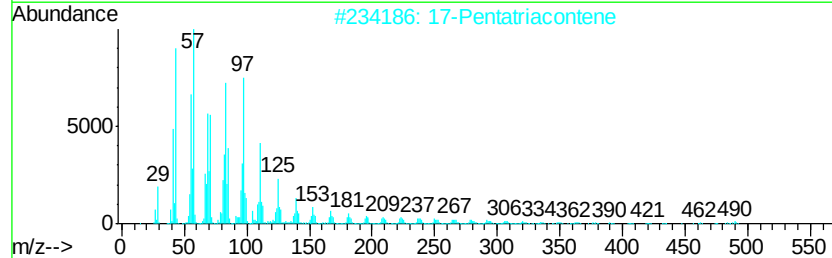
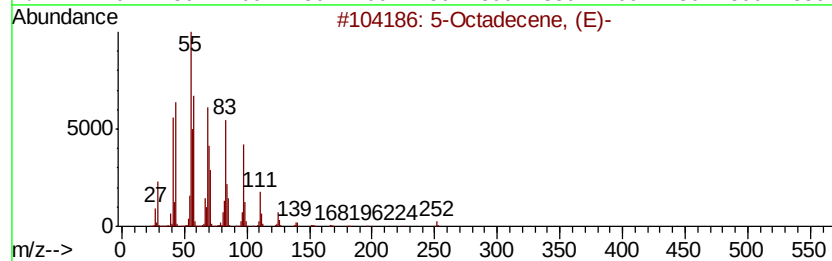
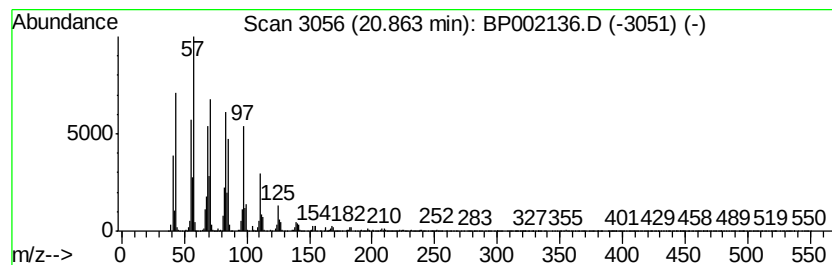
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	3.54 ng/ul	836760	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2		17-Pentatriacontene	491	C35H70	006971-40-0	94
3		Behenic alcohol	326	C22H46O	000661-19-8	93
4		1-Heneicosanol	312	C21H44O	015594-90-8	93
5		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93



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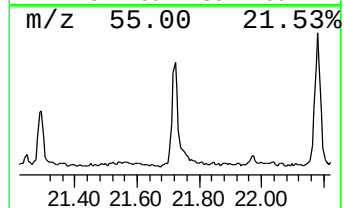
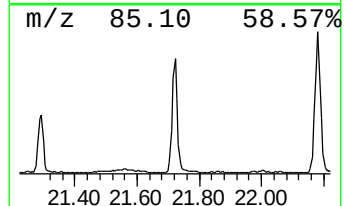
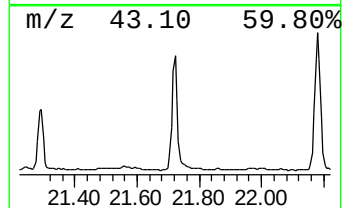
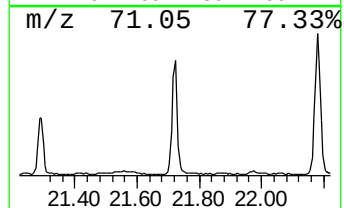
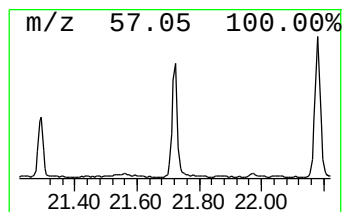
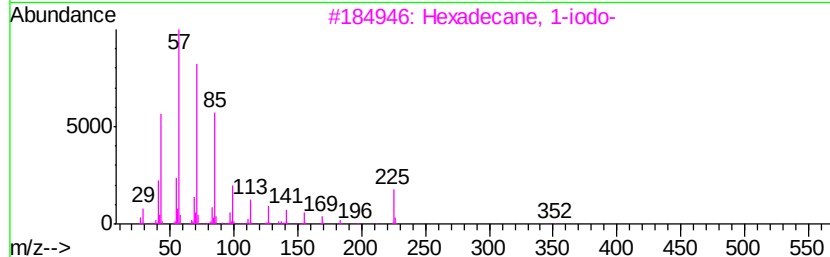
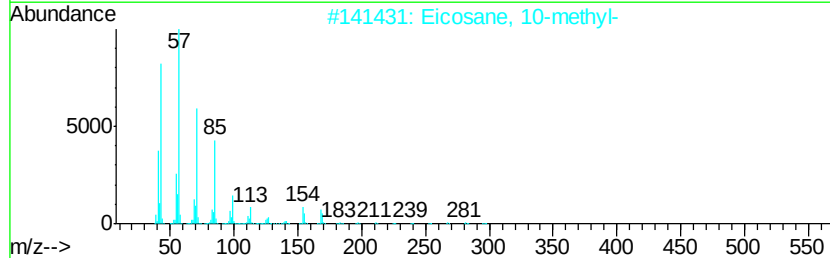
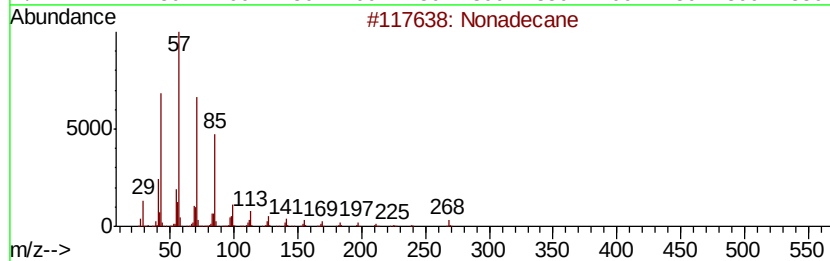
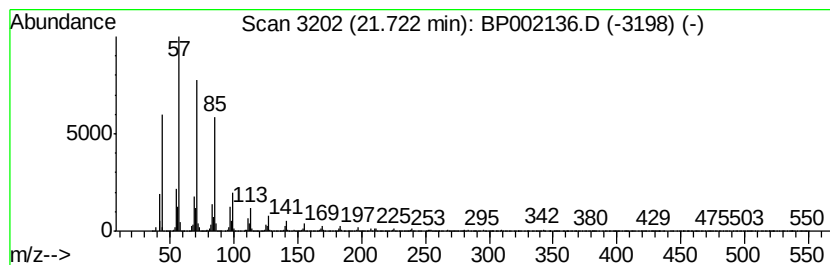
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	2.87 ng/ul	678340	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
5		Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002136.D
Acq On : 10 Mar 2020 00:02
Operator : CG/JU
Sample : L1834-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0DK9

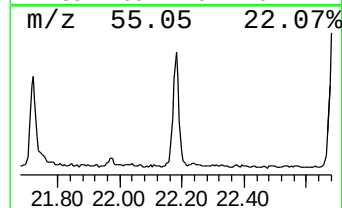
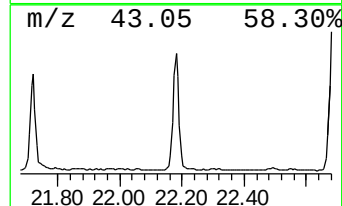
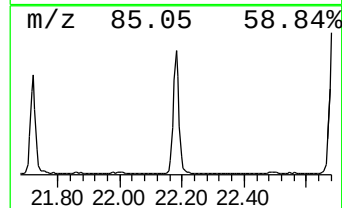
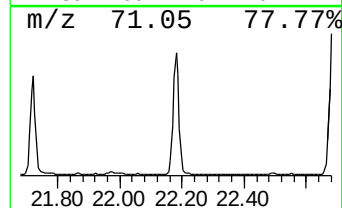
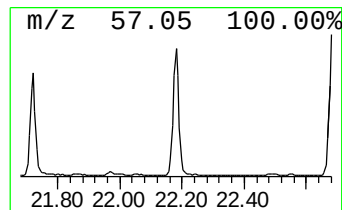
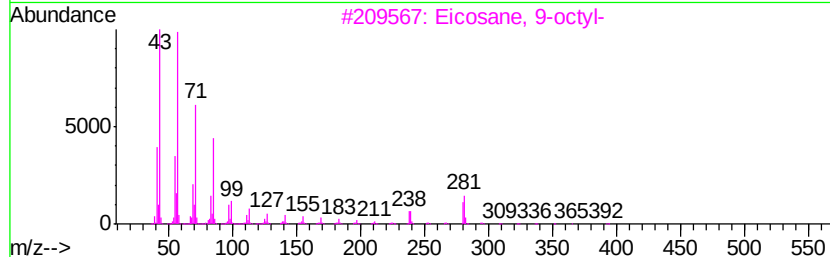
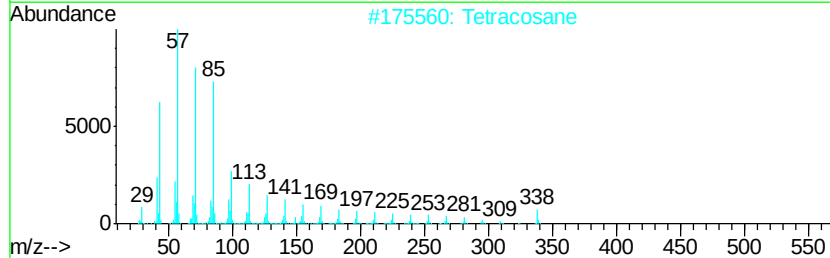
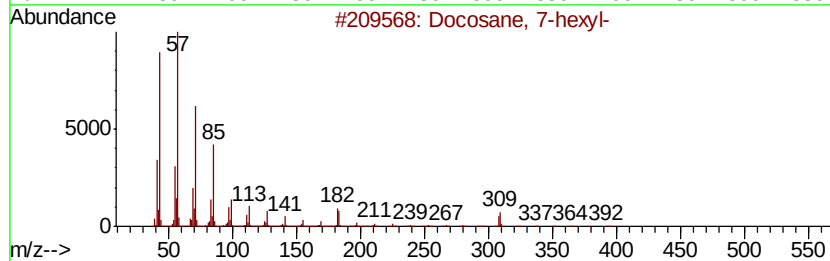
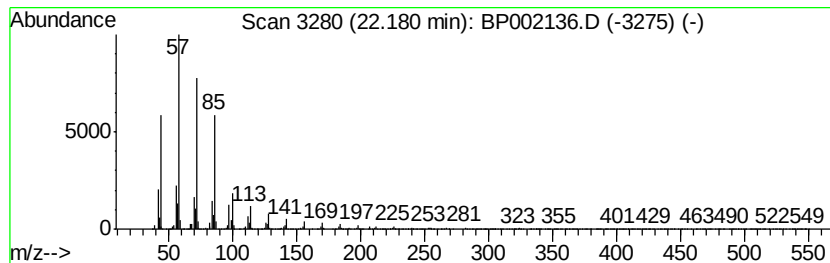
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.18	3.25 ng/ul	769693	Chrysene-d12	21.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
2		Tetracosane	338	C24H50	000646-31-1	94
3		Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
4		Triacontane	422	C30H62	000638-68-6	94
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002136.D
Acq On : 10 Mar 2020 00:02
Operator : CG/JU
Sample : L1834-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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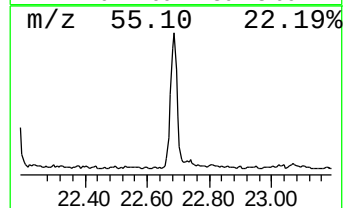
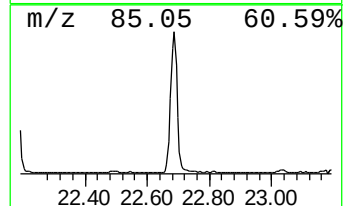
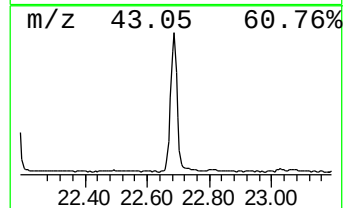
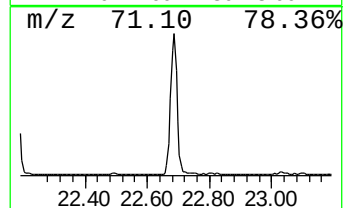
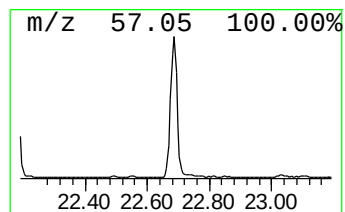
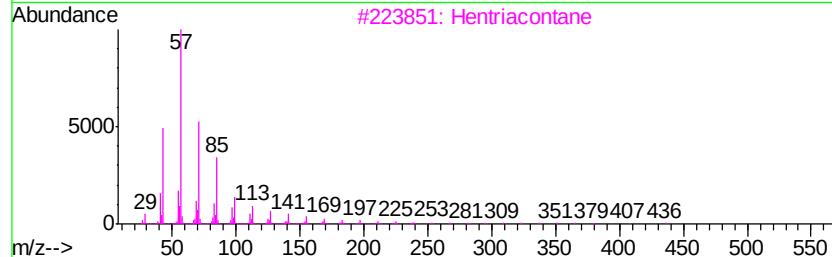
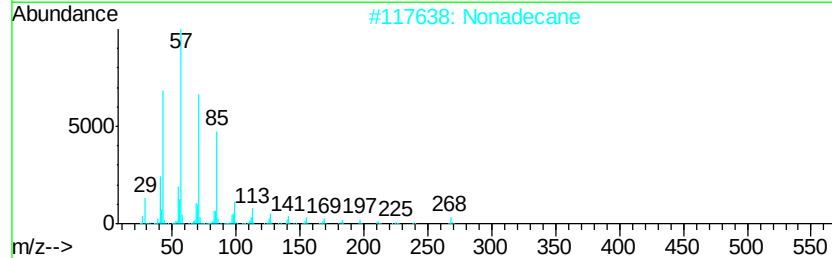
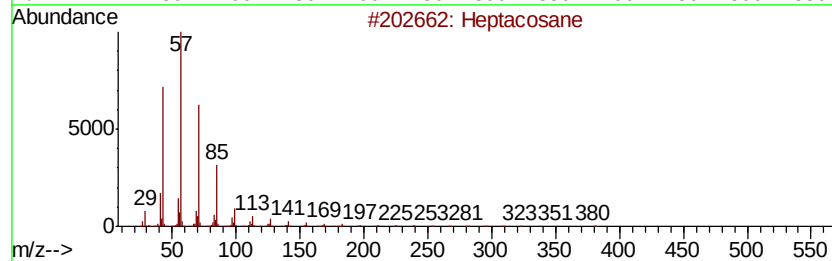
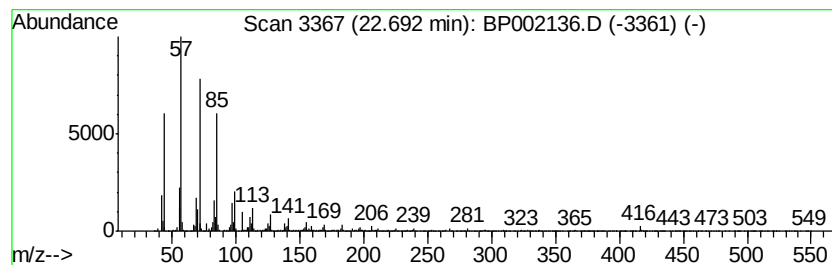
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	5.09 ng/ul	1266260	Perylene-d12	23.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	99
2		Nonadecane	268	C19H40	000629-92-5	93
3		Hentriacontane	437	C31H64	000630-04-6	92
4		Heptacosane	380	C27H56	000593-49-7	92
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002136.D
Acq On : 10 Mar 2020 00:02
Operator : CG/JU
Sample : L1834-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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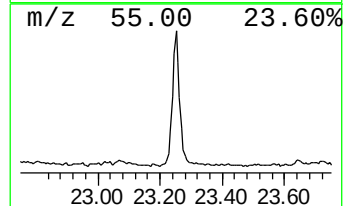
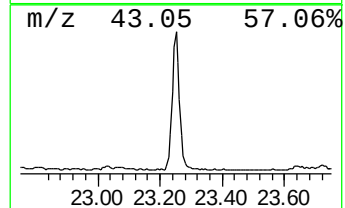
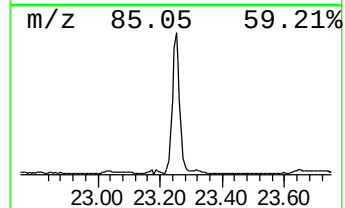
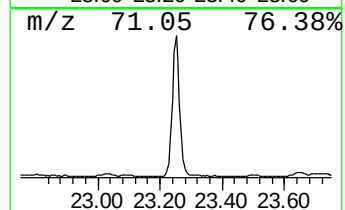
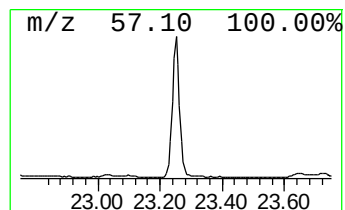
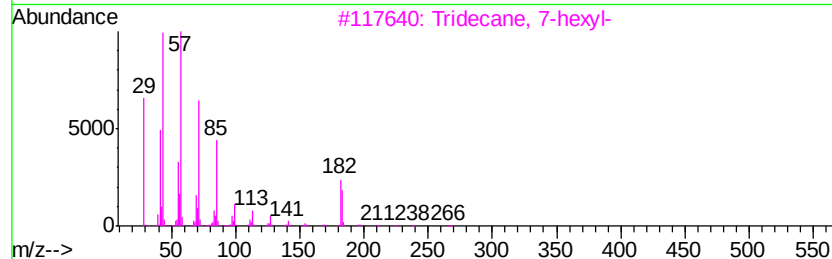
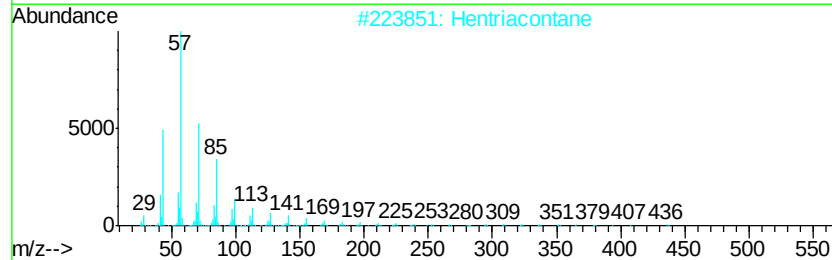
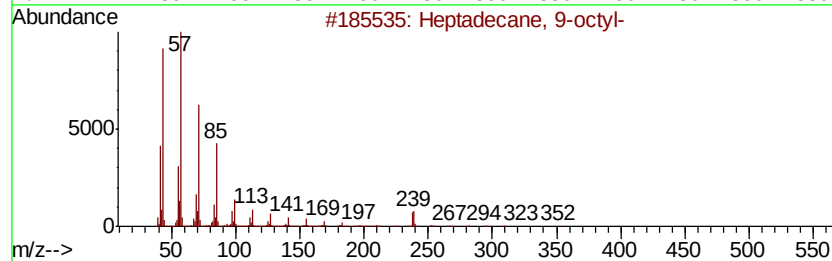
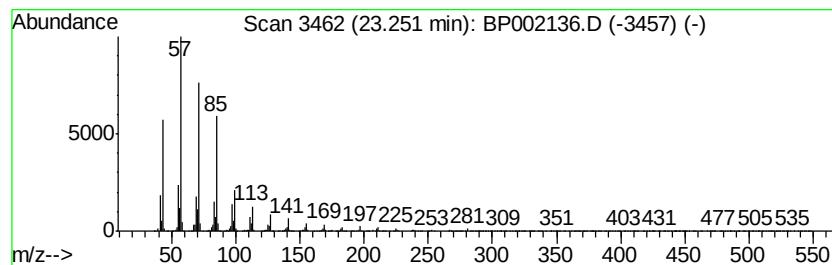
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.25	4.64 ng/ul	1153030	Perylene-d12	23.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2		Hentriacontane	437	C31H64	000630-04-6	93
3		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	93
4		Triacontane	422	C30H62	000638-68-6	91
5		Docosane	310	C22H46	000629-97-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002136.D
Acq On : 10 Mar 2020 00:02
Operator : CG/JU
Sample : L1834-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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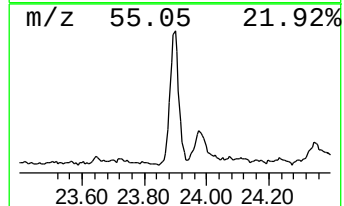
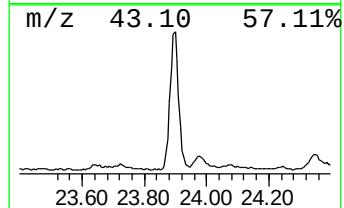
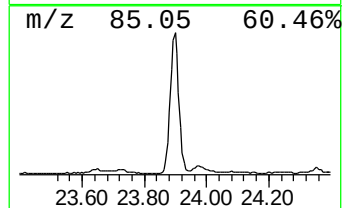
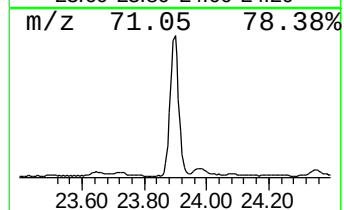
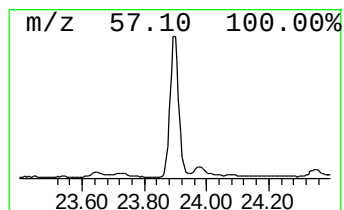
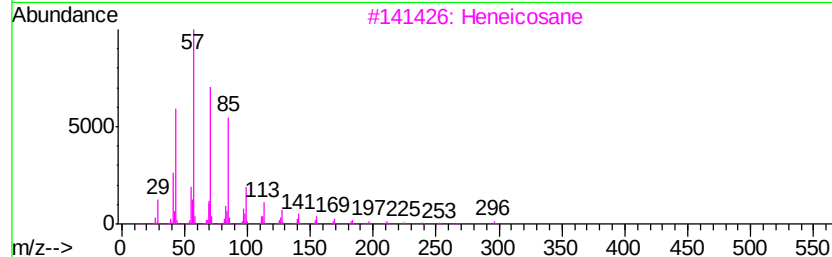
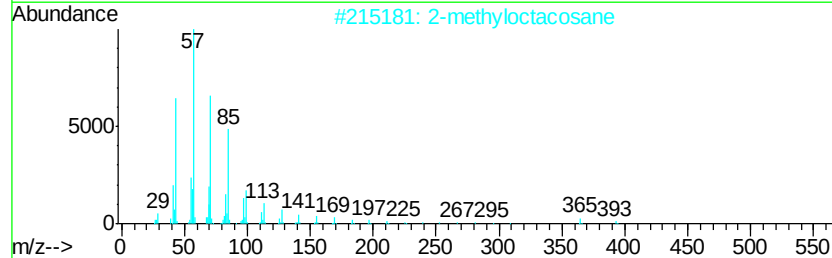
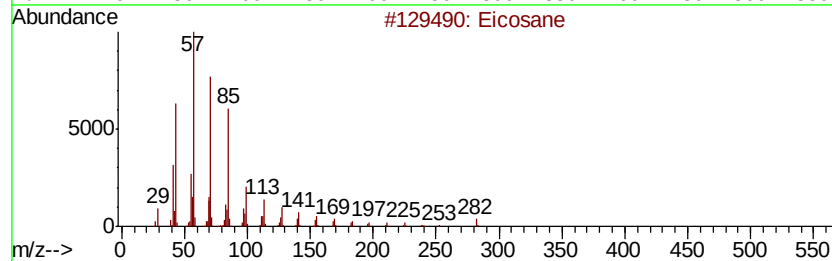
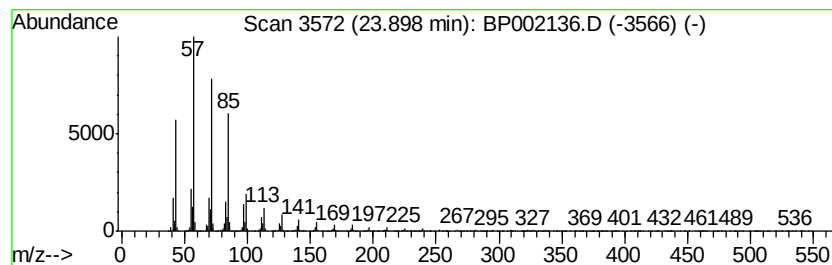
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.90	4.27 ng/ul	1062030	Perylene-d12	23.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP030920\
Data File : BP002136.D
Acq On : 10 Mar 2020 00:02
Operator : CG/JU
Sample : L1834-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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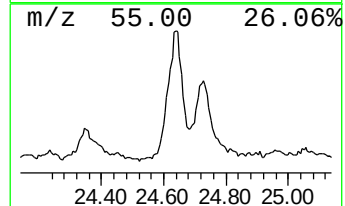
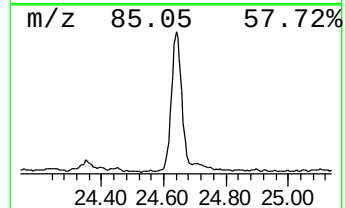
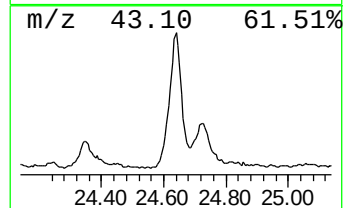
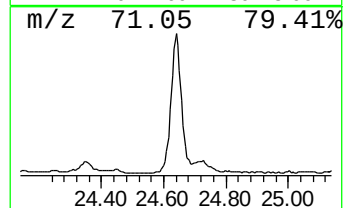
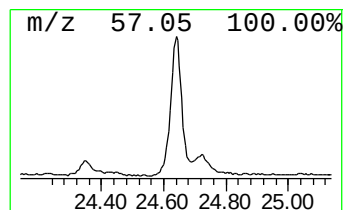
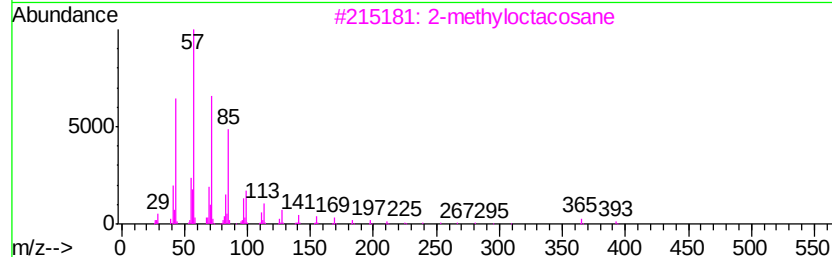
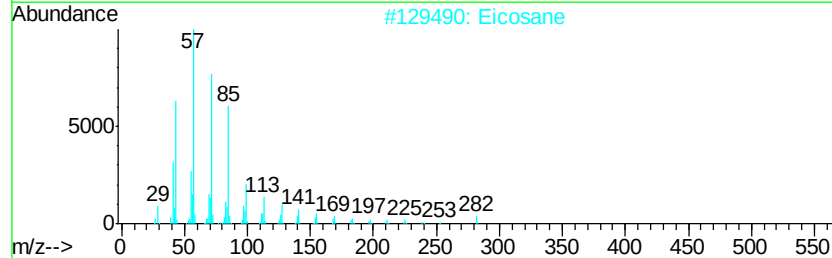
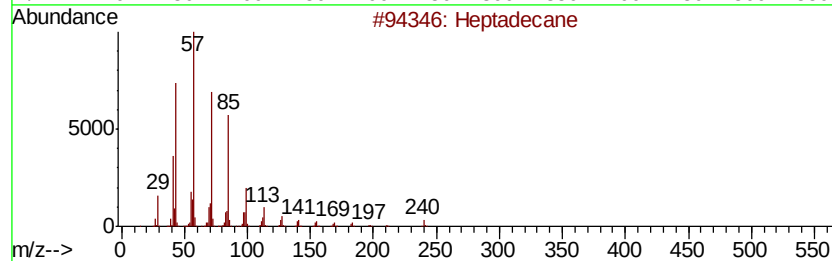
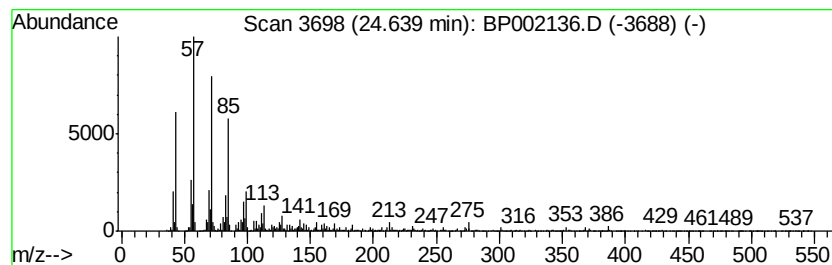
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.64	5.82 ng/ul	1448100	Perylene-d12	23.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	94
2		Eicosane	282	C20H42	000112-95-8	93
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030920\
Data File : BP002136.D
Acq On : 10 Mar 2020 00:02
Operator : CG/JU
Sample : L1834-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA 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
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n-Hexadecanoic acid	17.95	2.3	ng/ul	456259	4	17.02	3986250	20.0
(DEL) Alkane: Str...	20.86	3.5	ng/ul	836760	5	21.12	4732910	20.0
(DEL) Alkane: Str...	21.72	2.9	ng/ul	678340	5	21.12	4732910	20.0
(DEL) Alkane: Str...	22.18	3.3	ng/ul	769693	5	21.12	4732910	20.0
(DEL) Alkane: Str...	22.69	5.1	ng/ul	1266260	6	23.32	4972930	20.0
(DEL) Alkane: Str...	23.25	4.6	ng/ul	1153030	6	23.32	4972930	20.0
(DEL) Alkane: Str...	23.90	4.3	ng/ul	1062030	6	23.32	4972930	20.0
(DEL) Alkane: Str...	24.64	5.8	ng/ul	1448100	6	23.32	4972930	20.0