

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
 Data File : BM023371.D
 Acq On : 24 Oct 2019 10:11
 Operator : JU
 Sample : K5308-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2J25

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_M\METHODS\SOM-EPA-BM102319MA.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	3.157	26	29	34	rVB	71495	93214	0.22%	0.024%
2	4.199	201	206	213	rBV	554573	788918	1.86%	0.203%
3	4.393	236	239	246	rVB	113071	184162	0.43%	0.047%
4	4.787	302	306	316	rVB	534379	775155	1.82%	0.200%
5	6.775	637	644	648	rBV	4528470	7516644	17.70%	1.936%
6	6.810	648	650	663	rVB	1667518	2207055	5.20%	0.568%
7	6.975	672	678	691	rVB	1110345	1813985	4.27%	0.467%
8	7.169	704	711	713	rBV	1640778	2709183	6.38%	0.698%
9	7.198	713	716	724	rVB	3624855	5426025	12.77%	1.397%
10	7.634	783	790	798	rVB	1545785	2503726	5.89%	0.645%
11	8.339	901	910	915	rBV	1763969	2782917	6.55%	0.717%
12	8.398	915	920	924	rVV	2025074	3116457	7.34%	0.803%
13	8.451	924	929	939	rVB	3864493	6114625	14.40%	1.575%
14	8.704	965	972	979	rBV	2652202	4390083	10.34%	1.131%
15	8.781	979	985	994	rVB	1431667	2369512	5.58%	0.610%
16	9.251	1059	1065	1073	rVB	62297	102927	0.24%	0.027%
17	9.498	1100	1107	1114	rBV	1125570	1818906	4.28%	0.468%
18	10.063	1191	1203	1216	rBV2	4913899	10879092	25.61%	2.802%
19	10.410	1254	1262	1266	rBV	2113740	3641129	8.57%	0.938%
20	10.463	1266	1271	1278	rVV	6065676	9913074	23.34%	2.553%
21	10.545	1278	1285	1297	rVV	1680811	2984034	7.03%	0.769%
22	11.339	1414	1420	1441	rBV	253443	778891	1.83%	0.201%
23	11.957	1517	1525	1531	rBV4	82223	164574	0.39%	0.042%
24	12.010	1531	1534	1538	rVB	35173	51006	0.12%	0.013%
25	12.086	1538	1547	1554	rBV	7146065	11671229	27.48%	3.006%
26	12.198	1560	1566	1576	rVB7	26971	66040	0.16%	0.017%
27	12.304	1576	1584	1590	rBV	85212	144198	0.34%	0.037%
28	12.751	1655	1660	1666	rBV	62342	106468	0.25%	0.027%
29	13.098	1714	1719	1721	rBV4	60434	91157	0.21%	0.023%
30	13.145	1721	1727	1734	rVB	7549942	11720195	27.59%	3.018%
31	13.698	1814	1821	1825	rVV	3332706	4602447	10.84%	1.185%
32	13.745	1825	1829	1837	rVV	874301	1223370	2.88%	0.315%
33	13.969	1860	1867	1875	rBV	3874908	5684434	13.38%	1.464%
34	14.280	1912	1920	1925	rBV	3194431	4693080	11.05%	1.209%

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35	14.345	1925	1931	1938	rVB	5466257	7876329	18.54%	2.028%
36	14.480	1948	1954	1970	rBV	1468903	2152683	5.07%	0.554%
37	14.716	1990	1994	2000	rVB8	51705	85829	0.20%	0.022%
38	15.168	2066	2071	2075	rVV2	40363	53441	0.13%	0.014%
39	15.274	2083	2089	2094	rBV	5003628	7128342	16.78%	1.836%
40	15.333	2094	2099	2105	rVV	6277190	8655092	20.38%	2.229%
41	15.410	2105	2112	2124	rVV3	3701837	6126721	14.42%	1.578%
42	15.515	2124	2130	2134	rVV2	76132	140094	0.33%	0.036%
43	16.445	2283	2288	2292	rBV4	83512	123484	0.29%	0.032%
44	16.510	2292	2299	2303	rBV	7142869	9489909	22.34%	2.444%
45	16.545	2303	2305	2309	rVB2	106625	110220	0.26%	0.028%
46	16.845	2348	2356	2361	rVB	139553	240646	0.57%	0.062%
47	17.027	2380	2387	2390	rBV2	3812683	5603432	13.19%	1.443%
48	17.068	2390	2394	2399	rVV	8953808	13215491	31.11%	3.404%
49	17.121	2399	2403	2412	rVB	5363029	7455877	17.55%	1.920%
50	17.280	2426	2430	2435	rVB5	42302	74250	0.17%	0.019%
51	17.427	2448	2455	2466	rBV	12152698	16730321	39.39%	4.309%
52	17.951	2538	2544	2551	rBV	652890	1054643	2.48%	0.272%
53	18.009	2552	2554	2558	rVB	69480	96033	0.23%	0.025%
54	18.251	2590	2595	2604	rBV2	205310	330559	0.78%	0.085%
55	18.892	2701	2704	2708	rVB3	59227	66162	0.16%	0.017%
56	19.056	2728	2732	2735	rBV	92843	119466	0.28%	0.031%
57	19.233	2758	2762	2771	rBV2	236046	402975	0.95%	0.104%
58	19.333	2772	2779	2782	rVB2	123778	259619	0.61%	0.067%
59	19.421	2788	2794	2796	rBV2	6567688	9176000	21.60%	2.363%
60	19.451	2796	2799	2806	rVB	13748888	18217960	42.89%	4.692%
61	20.374	2951	2956	2961	rBV	10365027	11904828	28.03%	3.066%
62	20.503	2976	2978	2982	rBV	127902	108005	0.25%	0.028%
63	20.909	3043	3047	3051	rBV	333023	578371	1.36%	0.149%
64	20.962	3052	3056	3061	rVV3	1632855	2429546	5.72%	0.626%
65	21.009	3061	3064	3070	rVB	589449	752450	1.77%	0.194%
66	21.162	3085	3090	3094	rBV	11859561	13223620	31.13%	3.406%
67	21.215	3094	3099	3105	rVB	5119527	6463103	15.22%	1.665%
68	21.403	3127	3131	3137	rBV	1836260	2022762	4.76%	0.521%
69	21.833	3200	3204	3211	rBV	3232113	4082747	9.61%	1.051%
70	22.292	3277	3282	3290	rBV	4957796	6761523	15.92%	1.741%
71	22.797	3362	3368	3382	rVB	6011345	9271573	21.83%	2.388%

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72	23.274	3443	3449	3456	rVB	5802702	10005529	23.56%	2.577%
73	23.356	3457	3463	3468	rVV	6332438	10671737	25.12%	2.748%
74	23.409	3468	3472	3485	rVB2	3654064	6972207	16.41%	1.796%
75	23.997	3565	3572	3581	rBV	4807874	9410411	22.15%	2.424%
76	24.733	3690	3697	3709	rVB	2948473	6621768	15.59%	1.705%
77	25.627	3837	3849	3861	rVB2	13488074	42475982	100.00%	10.939%
78	26.291	3952	3962	3976	rVB	5666076	16417373	38.65%	4.228%

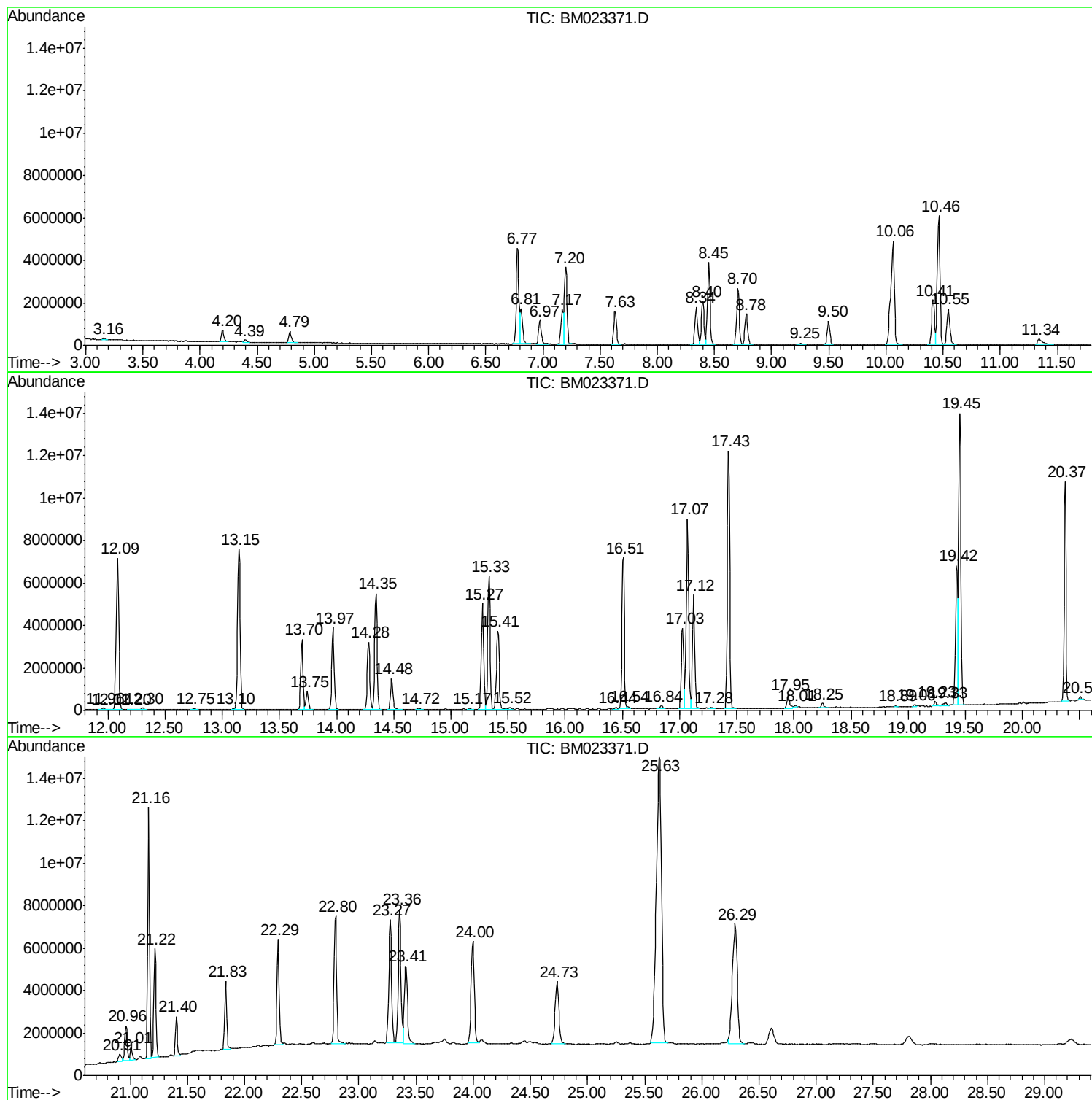
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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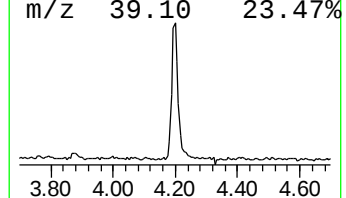
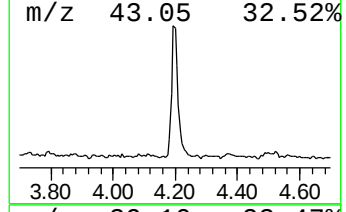
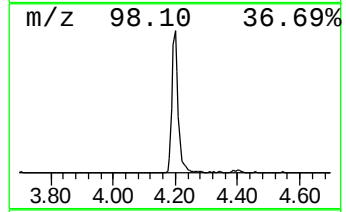
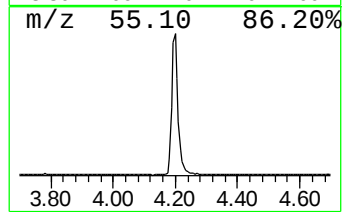
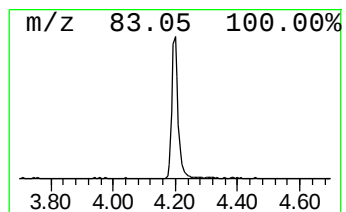
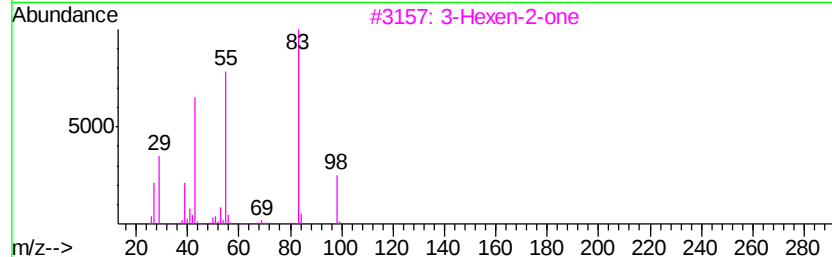
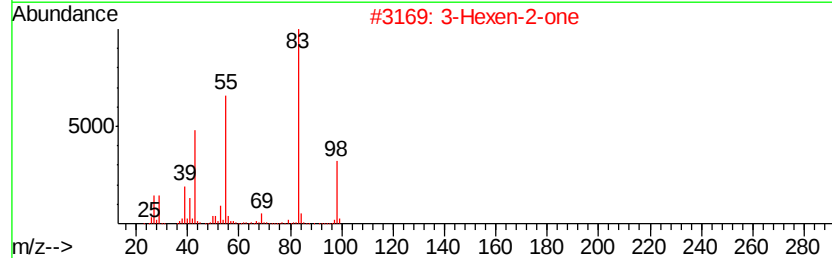
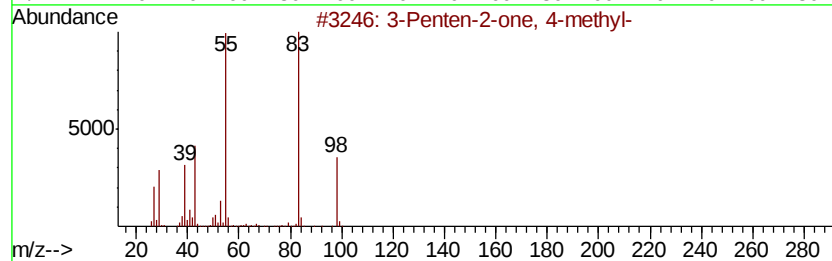
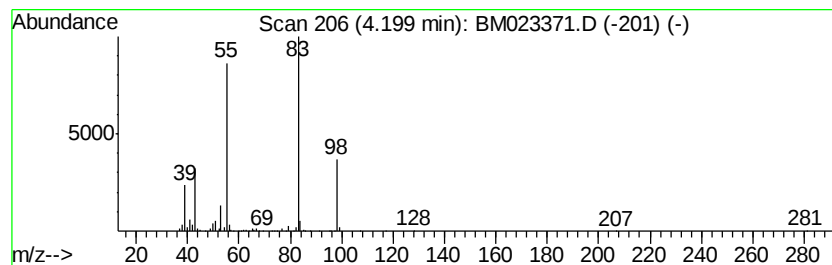
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	6.30 ng/ul	788918	1,4-Dichlorobenzene-d4	7.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2			3-Hexen-2-one	98	C6H10O	000763-93-9	91
3			3-Hexen-2-one	98	C6H10O	000763-93-9	90
4			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
5			3-Hexen-2-one	98	C6H10O	000763-93-9	86



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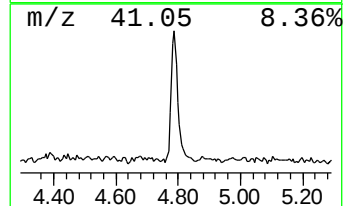
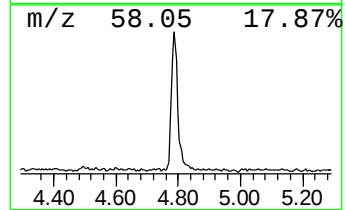
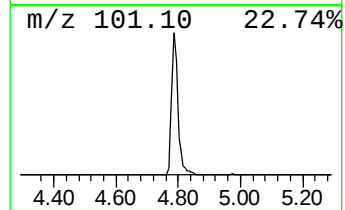
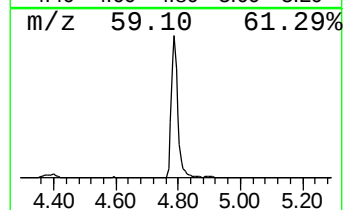
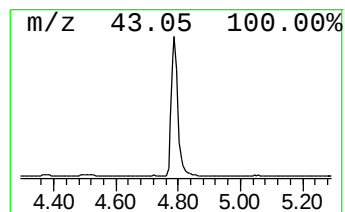
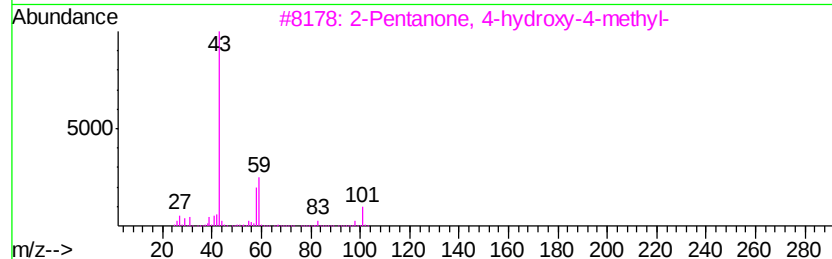
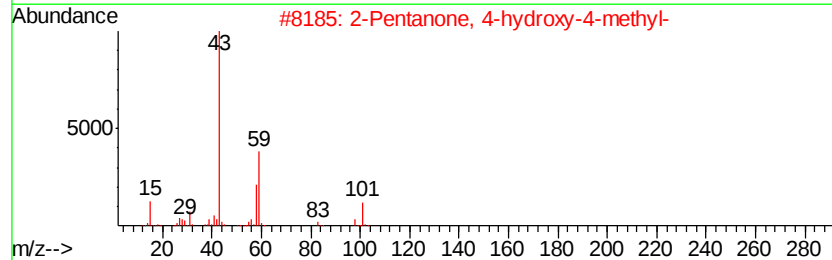
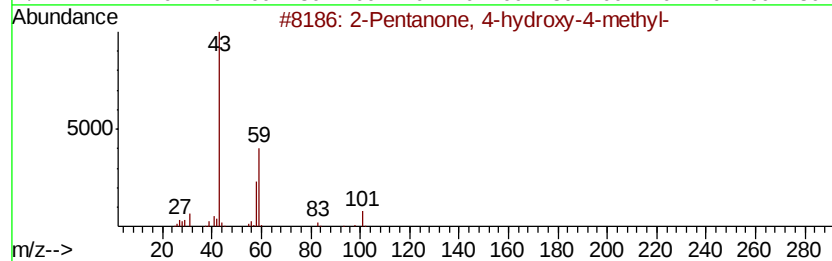
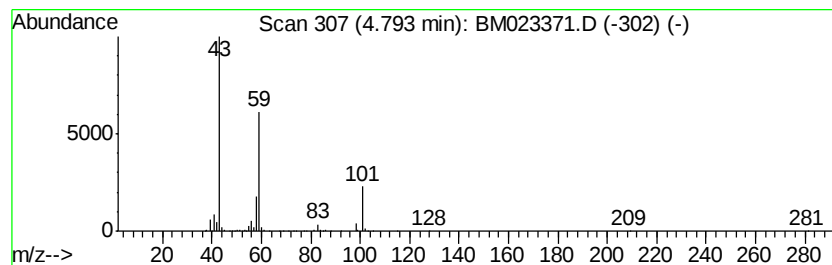
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	6.19 ng/ul	775155	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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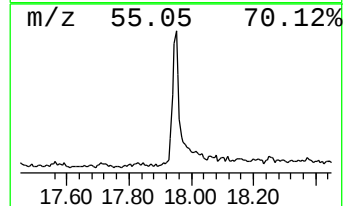
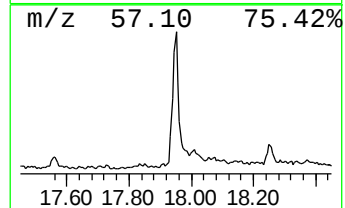
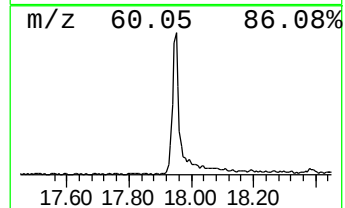
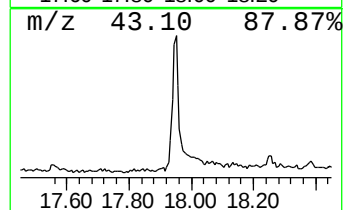
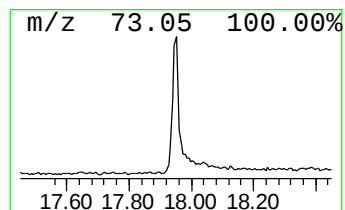
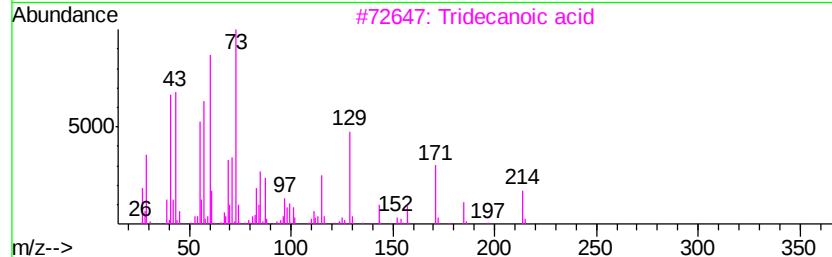
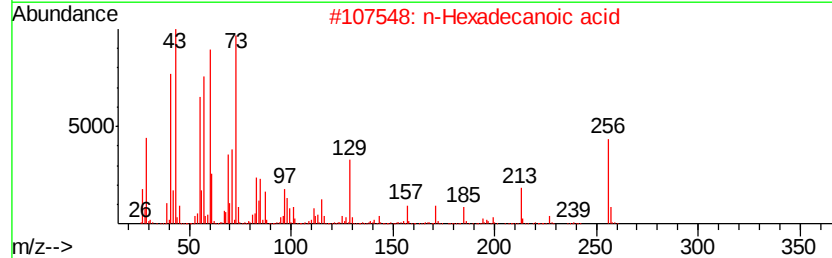
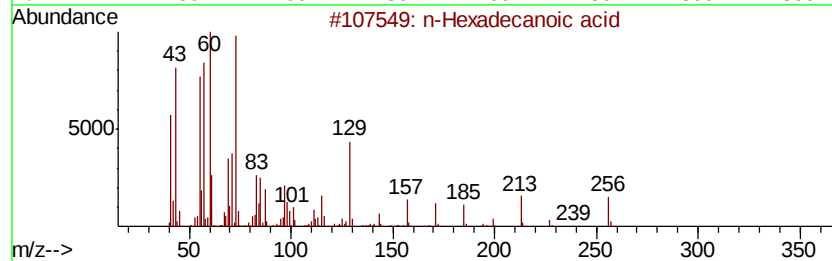
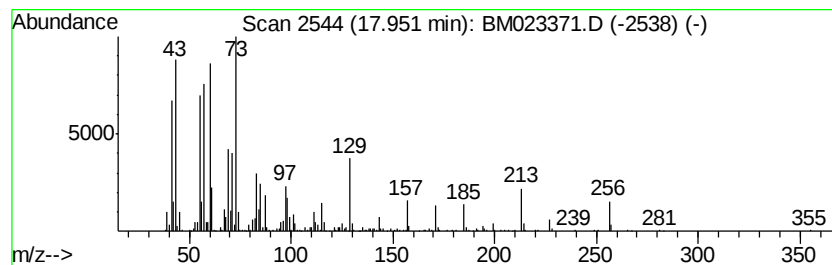
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.95	3.76 ng/ul	1054640	Phenanthrene-d10	17.03		
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	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	94
4		Octadecanoic acid	284	C18H36O2	000057-11-4	91
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	87



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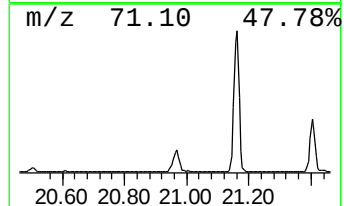
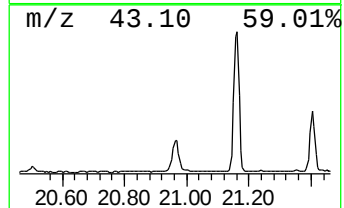
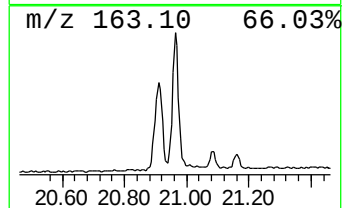
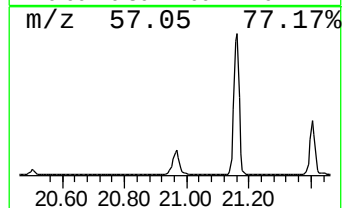
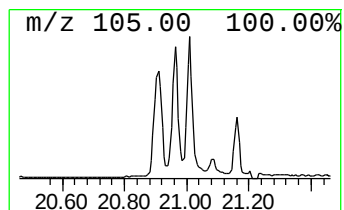
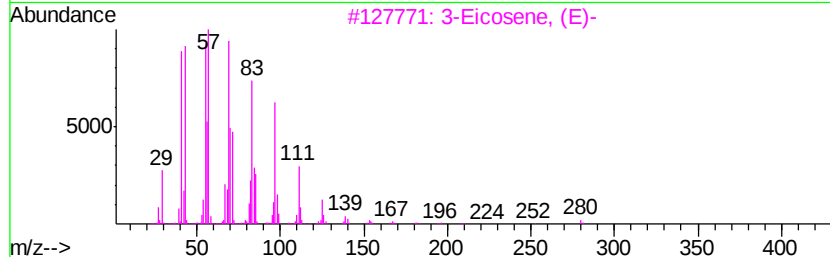
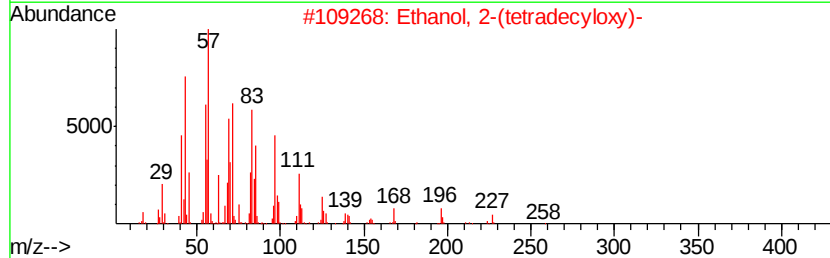
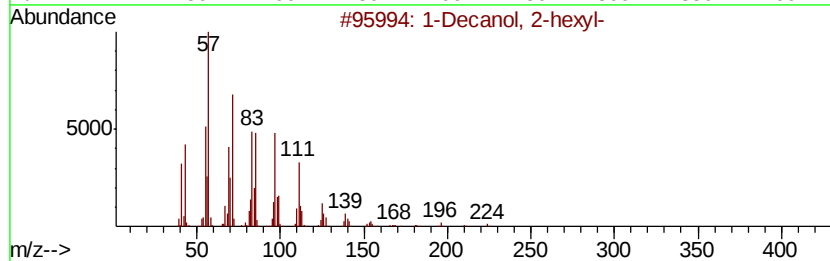
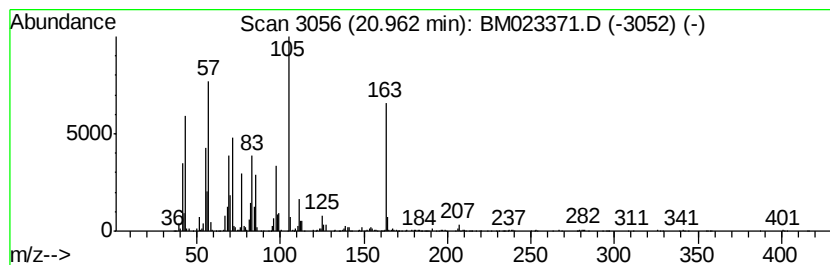
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Decanol, 2-hexyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	7.52 ng/ul	2429550	Chrysene-d12	21.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	64	
2	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	64	
3	3-Eicosene, (E)-	280	C20H40	074685-33-9	60	
4	Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	58	
5	Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	58	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
Data File : BM023371.D
Acq On : 24 Oct 2019 10:11
Operator : JU
Sample : K5308-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
X2J25

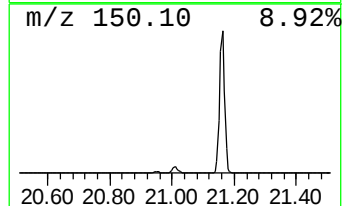
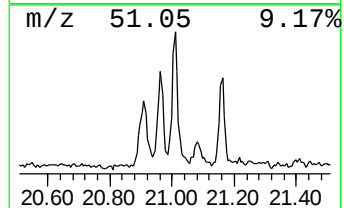
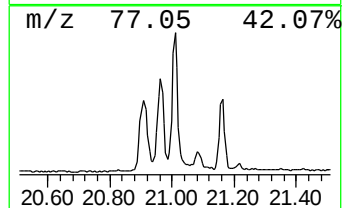
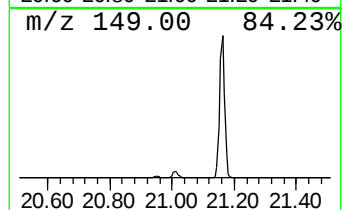
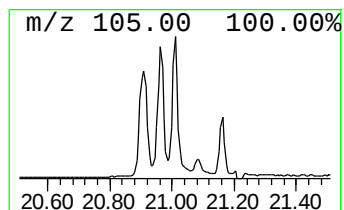
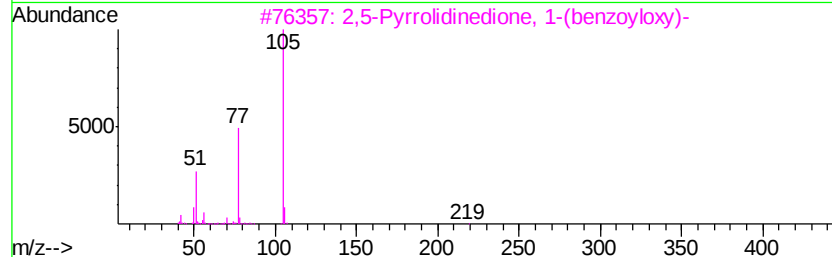
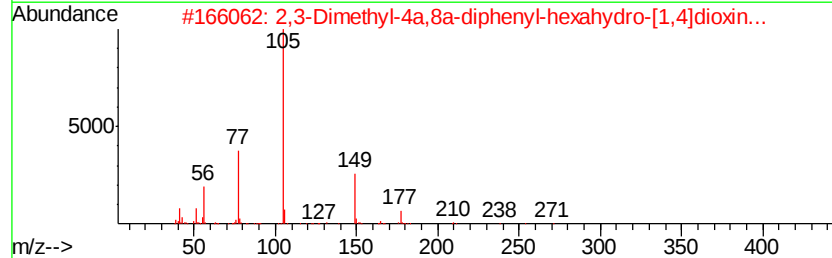
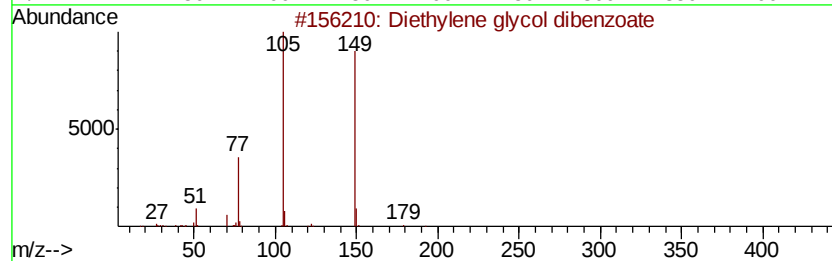
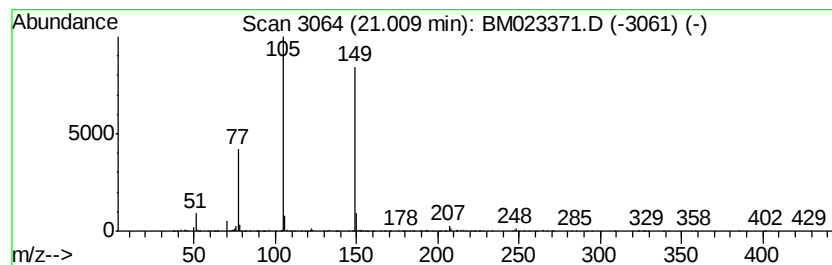
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Diethylene glycol dibenzoate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.01	2.33 ng/ul	752450	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2		2,3-Dimethyl-4a,8a-diphenyl-hexa...	326	C20H22O4	1000297-07-0	43
3		2,5-Pyrrolidinedione, 1-(benzoyl...	219	C11H9NO4	023405-15-4	35
4		1,2-Bis(benzoyloxy)benzene	318	C20H14O4	000643-94-7	35
5		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
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Sample : K5308-07
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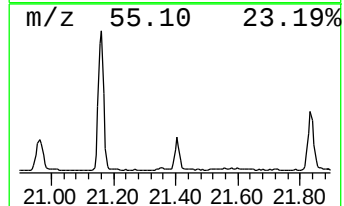
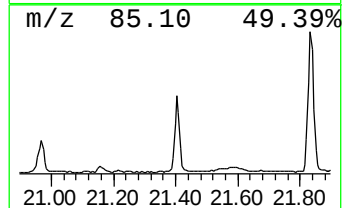
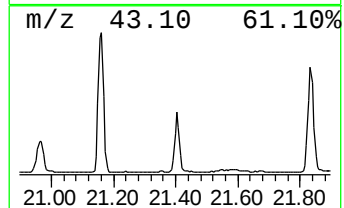
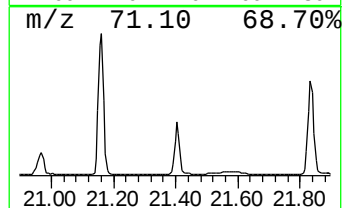
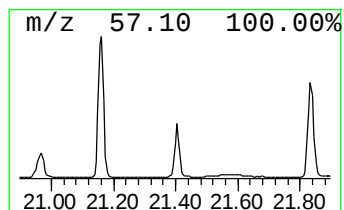
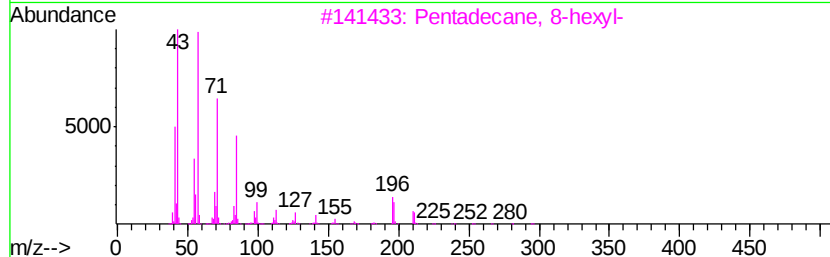
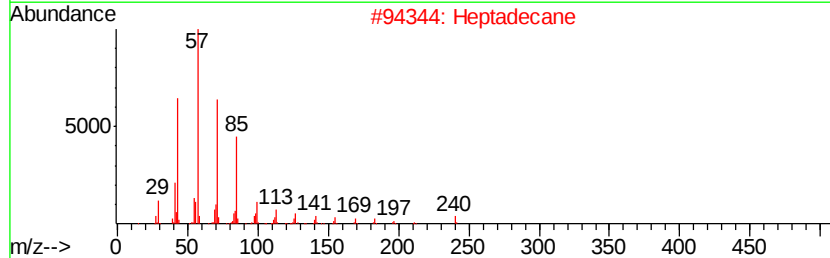
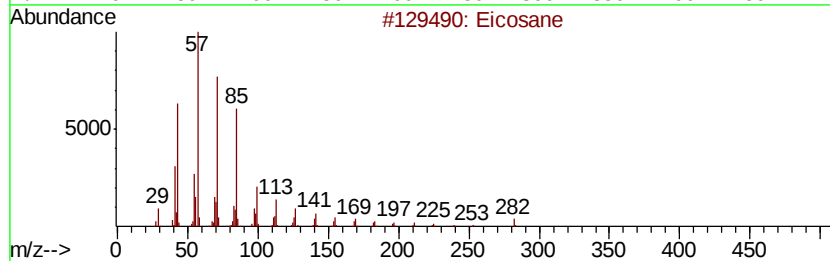
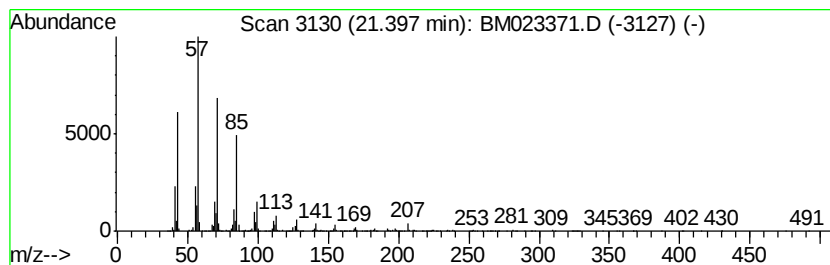
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.40	6.26 ng/ul	2022760	Chrysene-d12	21.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Heptadecane	240	C17H36	000629-78-7	94
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5	Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
Data File : BM023371.D
Acq On : 24 Oct 2019 10:11
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Sample : K5308-07
Misc :
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Instrument :
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ClientSampleId :
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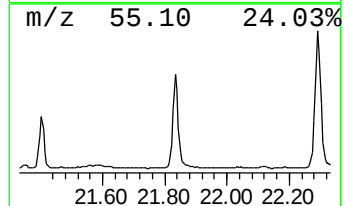
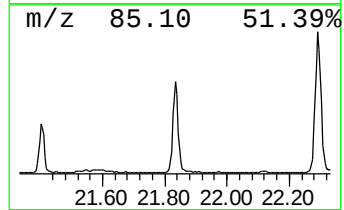
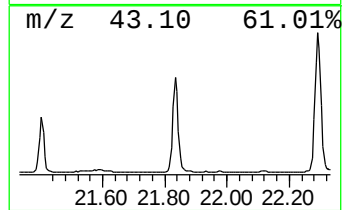
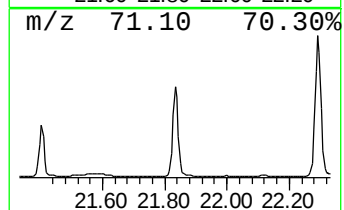
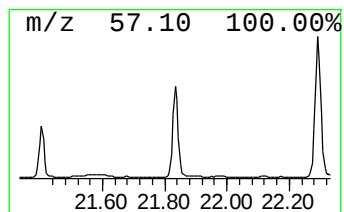
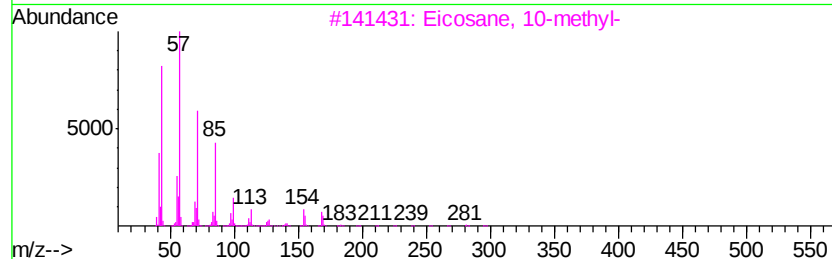
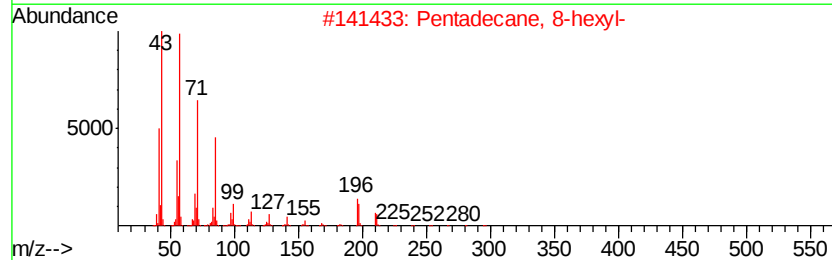
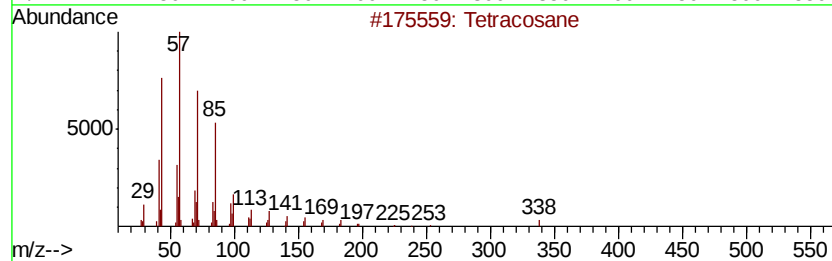
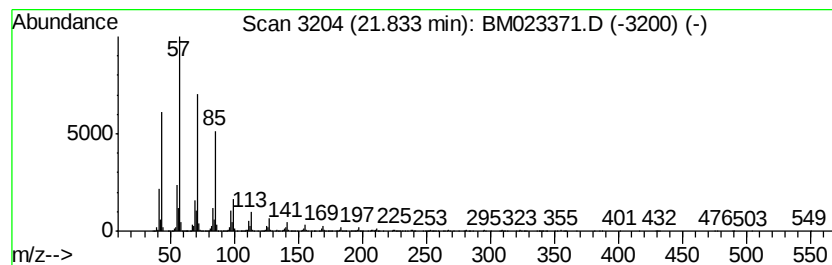
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	12.63 ng/ul	4082750	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
4		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
Data File : BM023371.D
Acq On : 24 Oct 2019 10:11
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Misc :
ALS Vial : 23 Sample Multiplier: 1

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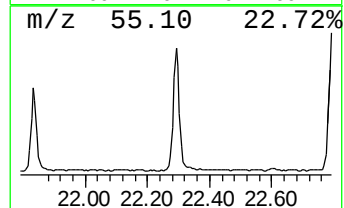
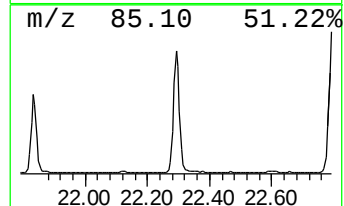
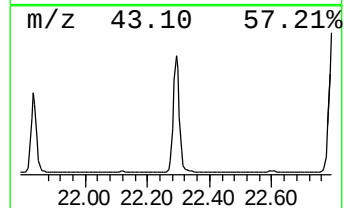
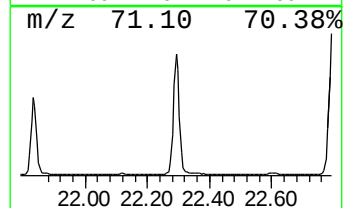
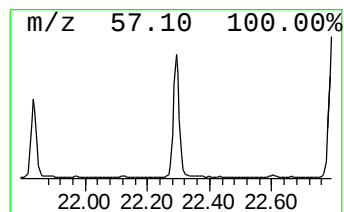
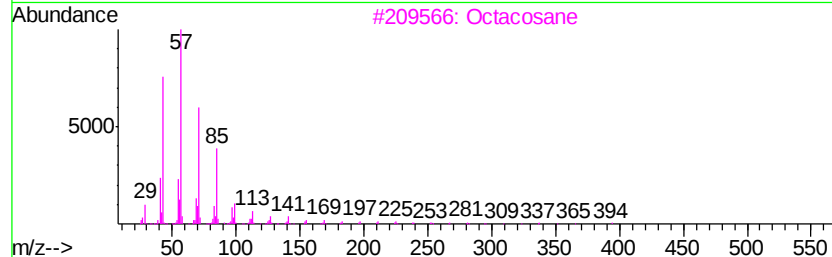
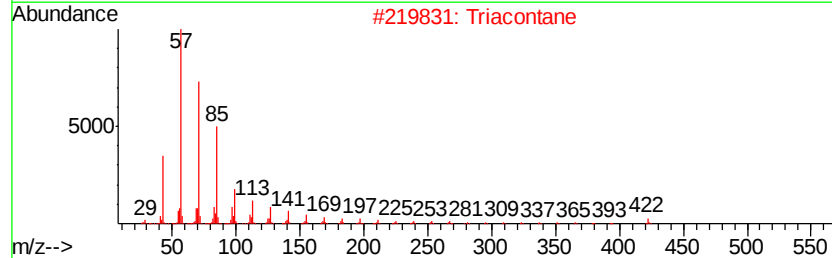
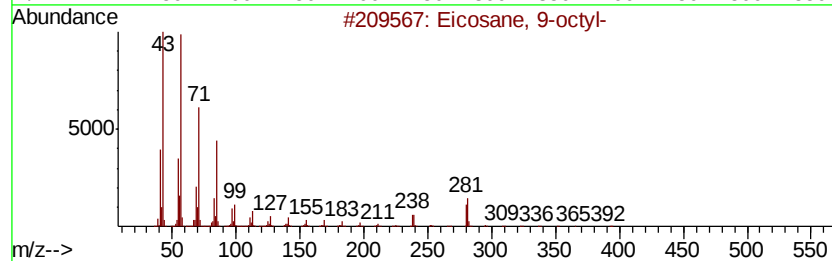
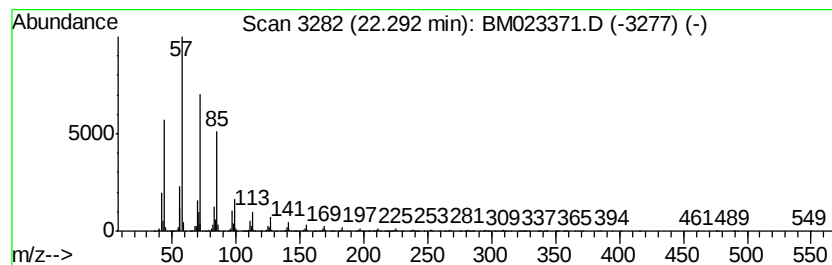
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.29	20.92 ng/ul	6761520	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
2		Triacontane	422	C30H62	000638-68-6	91
3		Octacosane	394	C28H58	000630-02-4	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Docosane	310	C22H46	000629-97-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
 Data File : BM023371.D
 Acq On : 24 Oct 2019 10:11
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Instrument :
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 ClientSampleId :
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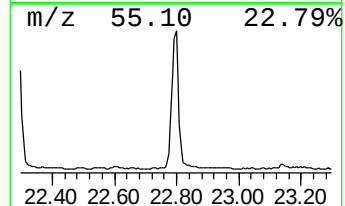
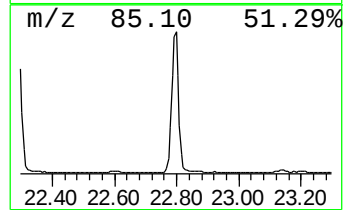
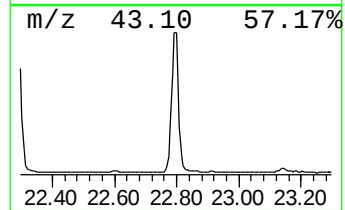
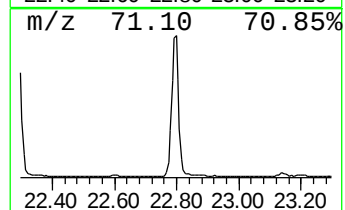
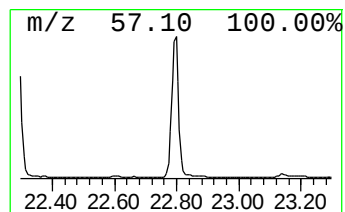
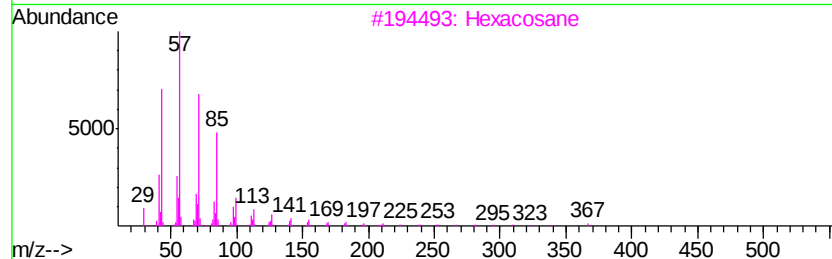
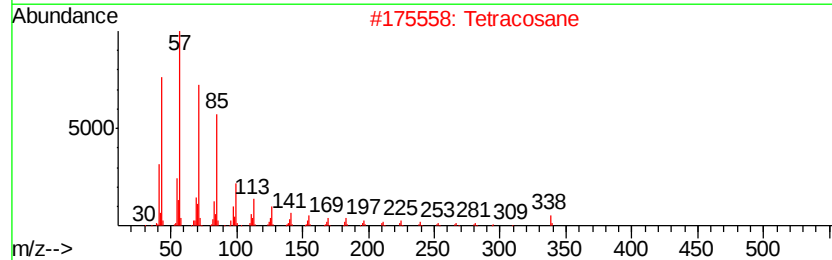
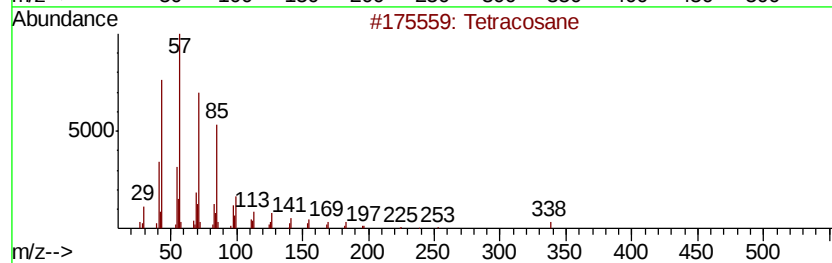
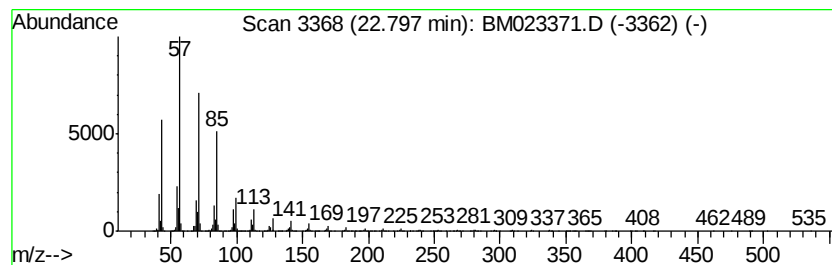
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.80	26.60 ng/ul	9271570	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	97
2		Tetracosane	338	C24H50	000646-31-1	97
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



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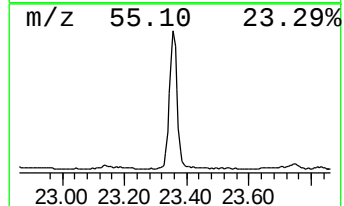
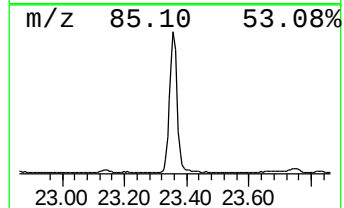
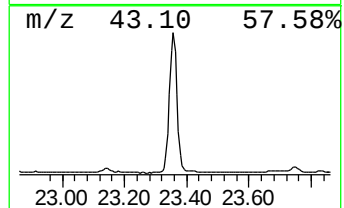
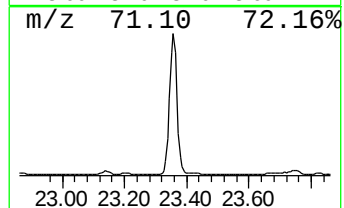
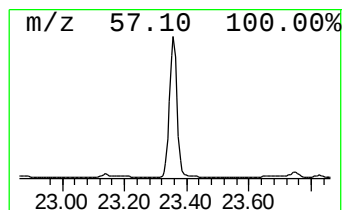
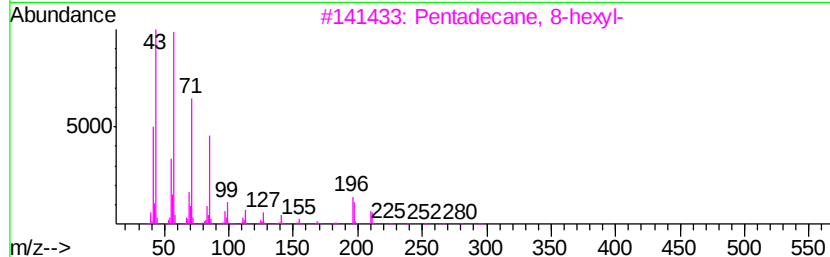
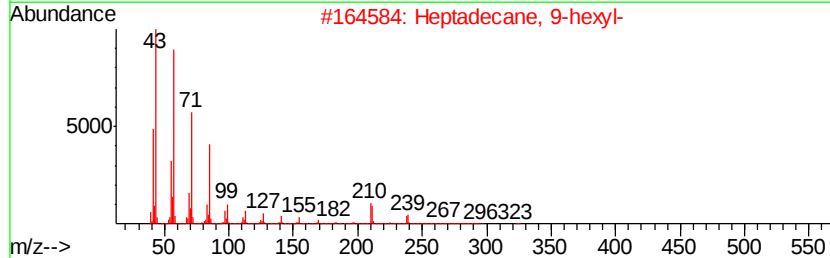
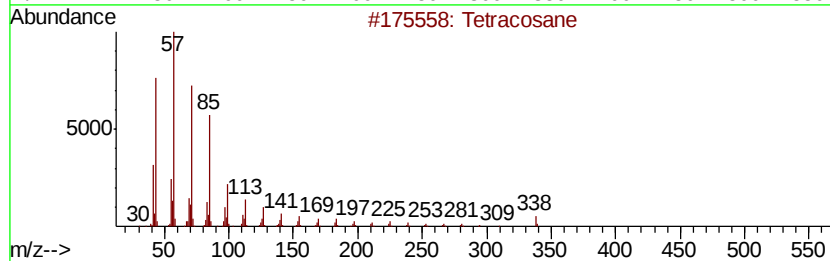
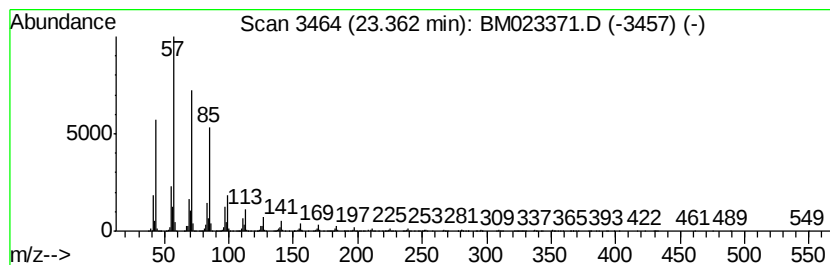
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.36	30.61 ng/ul	10671700	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



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ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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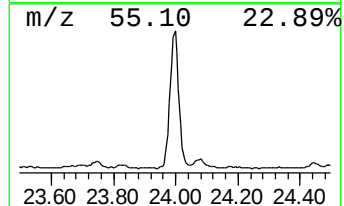
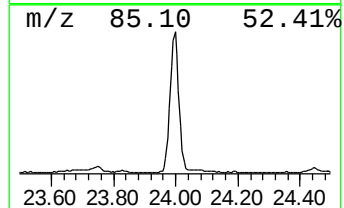
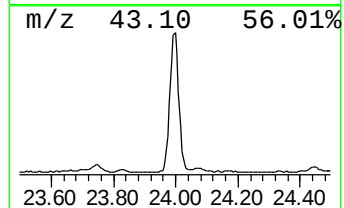
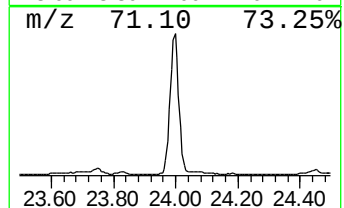
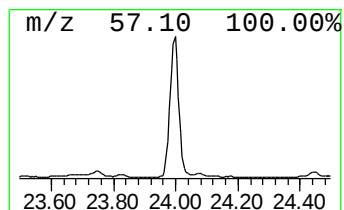
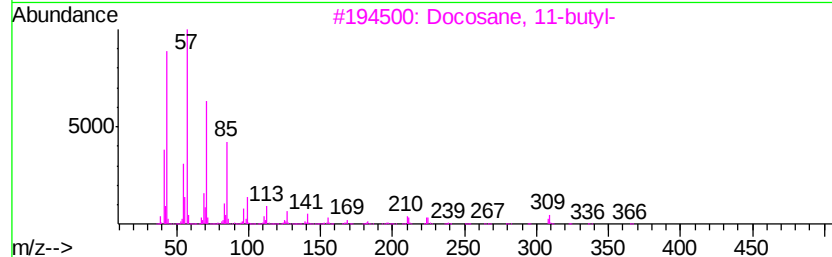
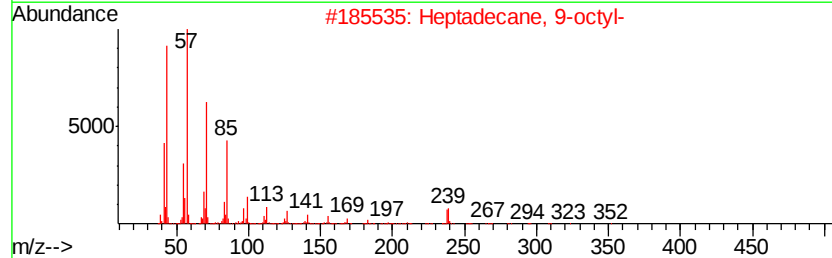
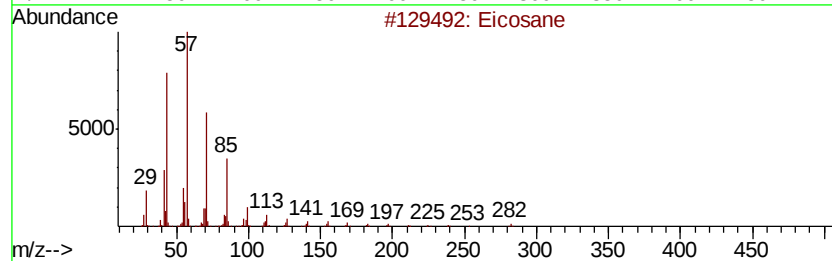
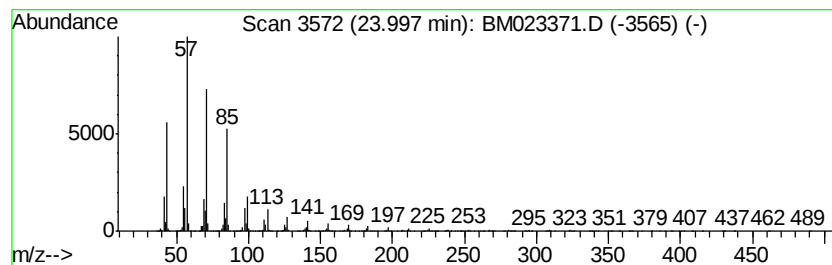
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.00	26.99 ng/ul	9410410	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	95
4		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
Data File : BM023371.D
Acq On : 24 Oct 2019 10:11
Operator : JU
Sample : K5308-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
X2J25

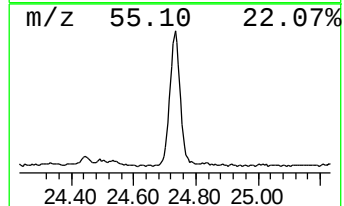
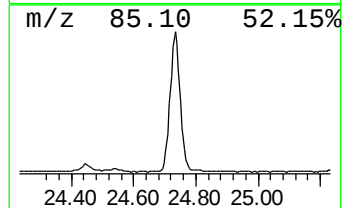
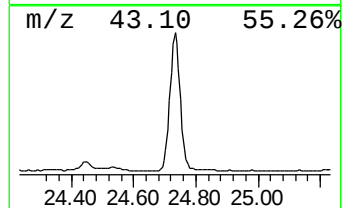
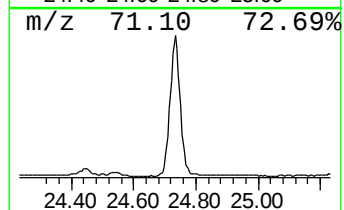
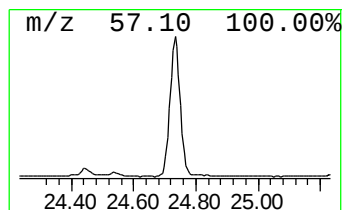
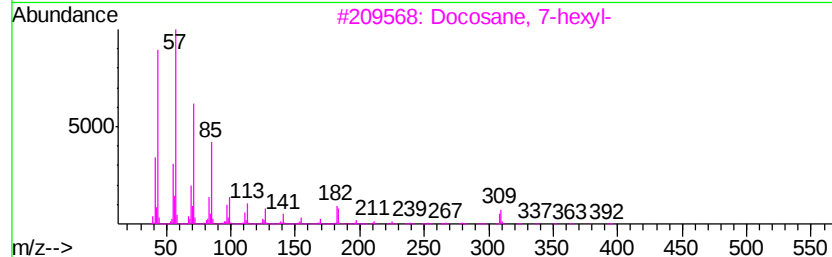
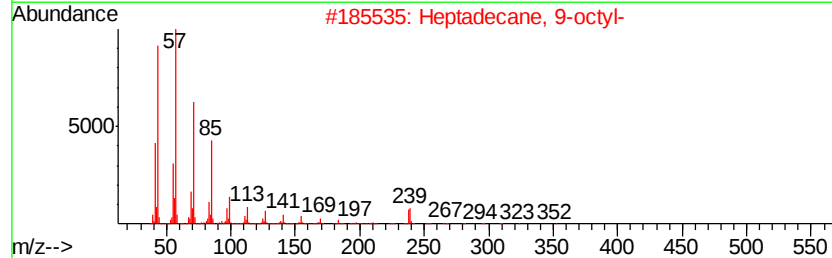
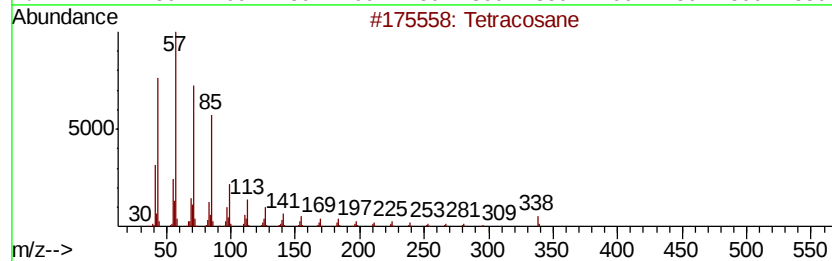
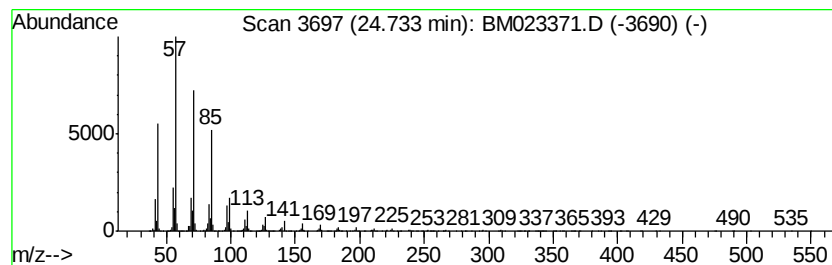
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Quant Title : SVOA CALIBRATION

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.
24.73	18.99 ng/ul	6621770	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	97
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3		Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
4		Heptacosane	380	C27H56	000593-49-7	91
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102319\
 Data File : BM023371.D
 Acq On : 24 Oct 2019 10:11
 Operator : JU
 Sample : K5308-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2J25

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.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
3-Penten-2-one, 4...	4.20	6.3	ng/ul	788918	1	7.63	2503730	20.0
2-Pentanone, 4-hy...	4.79	6.2	ng/ul	775155	1	7.63	2503730	20.0
n-Hexadecanoic acid	17.95	3.8	ng/ul	1054640	4	17.03	5603430	20.0
1-Decanol, 2-hexyl-	20.96	7.5	ng/ul	2429550	5	21.22	6463100	20.0
Diethylene glycol...	21.01	2.3	ng/ul	752450	5	21.22	6463100	20.0
(DEL) Alkane: Str...	21.40	6.3	ng/ul	2022760	5	21.22	6463100	20.0
(DEL) Alkane: Str...	21.83	12.6	ng/ul	4082750	5	21.22	6463100	20.0
(DEL) Alkane: Str...	22.29	20.9	ng/ul	6761520	5	21.22	6463100	20.0
(DEL) Alkane: Str...	22.80	26.6	ng/ul	9271570	6	23.41	6972210	20.0
(DEL) Alkane: Str...	23.36	30.6	ng/ul	10671700	6	23.41	6972210	20.0
(DEL) Alkane: Str...	24.00	27.0	ng/ul	9410410	6	23.41	6972210	20.0
(DEL) Alkane: Str...	24.73	19.0	ng/ul	6621770	6	23.41	6972210	20.0