

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
 Data File : BP006319.D
 Acq On : 24 Jul 2021 23:03
 Operator : CG/JU
 Sample : M2973-14
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGE19

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006319.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.305	33	37	44	rVB	73339	106681	1.41%	0.126%
2	3.587	82	85	93	rVB	12820	15890	0.21%	0.019%
3	3.711	101	106	126	rBV	902807	1468566	19.38%	1.732%
4	4.028	155	160	170	rVB	41107	66345	0.88%	0.078%
5	5.811	459	463	469	rBV9	3276	8497	0.11%	0.010%
6	6.781	623	628	635	rBV8	9641	17729	0.23%	0.021%
7	6.964	651	659	669	rBV	1247468	1958722	25.85%	2.309%
8	7.134	681	688	702	rVB	1032271	1576818	20.81%	1.859%
9	7.322	713	720	730	rBV	1293154	1988848	26.25%	2.345%
10	7.428	734	738	744	rVB5	8148	14531	0.19%	0.017%
11	7.793	789	800	806	rBV	1135698	1833359	24.20%	2.162%
12	7.858	806	811	818	rVB	29834	50504	0.67%	0.060%
13	8.034	834	841	846	rBV	6813	13457	0.18%	0.016%
14	8.493	912	919	932	rBV	1547094	2453721	32.38%	2.893%
15	8.946	988	996	1004	rBV	1114333	1821110	24.04%	2.147%
16	9.058	1009	1015	1020	rVV3	13173	21969	0.29%	0.026%
17	9.116	1021	1025	1030	rVB7	7085	9209	0.12%	0.011%
18	9.381	1063	1070	1076	rVB5	8488	14301	0.19%	0.017%
19	9.663	1108	1118	1134	rBV	858313	1415848	18.69%	1.669%
20	9.799	1138	1141	1145	rBV6	4164	7644	0.10%	0.009%
21	9.887	1151	1156	1160	rBV7	6650	10515	0.14%	0.012%
22	10.193	1200	1208	1219	rBV	1395228	2288951	30.21%	2.699%
23	10.363	1231	1237	1246	rBV	395496	610192	8.05%	0.719%
24	10.575	1264	1273	1281	rBV	1570796	2629180	34.70%	3.100%
25	10.716	1288	1297	1308	rBV	1345962	2272354	29.99%	2.679%
26	11.046	1349	1353	1357	rVB6	6395	9036	0.12%	0.011%
27	11.222	1378	1383	1391	rBV6	3808	9902	0.13%	0.012%
28	11.363	1403	1407	1412	rBV8	7506	12300	0.16%	0.015%
29	12.110	1524	1534	1538	rBV2	19407	43280	0.57%	0.051%
30	12.163	1539	1543	1548	rVB	26795	44452	0.59%	0.052%
31	12.746	1638	1642	1646	rBV6	5249	11132	0.15%	0.013%
32	12.781	1646	1648	1653	rVB4	8853	10079	0.13%	0.012%
33	13.034	1686	1691	1697	rBV5	10731	17944	0.24%	0.021%
34	13.263	1725	1730	1736	rVV2	21941	32565	0.43%	0.038%
35	13.334	1736	1742	1747	rVV5	12297	22992	0.30%	0.027%
36	13.387	1747	1751	1756	rVB5	9145	14325	0.19%	0.017%

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Stop Thrs : 0

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Title : SVOA CALIBRATION

37	13.728	1802	1809	1811	rBV7	6050	11162	0.15%	0.013%
38	13.834	1821	1827	1836	rBV	2249610	3220204	42.50%	3.797%
39	14.110	1864	1874	1882	rBV	2643202	3925144	51.80%	4.628%
40	14.334	1908	1912	1918	rVB3	10187	13027	0.17%	0.015%
41	14.416	1918	1926	1932	rVV	2278796	3318183	43.79%	3.912%
42	14.475	1932	1936	1945	rVB	56299	91260	1.20%	0.108%
43	14.610	1950	1959	1977	rBV	947039	1345384	17.76%	1.586%
44	14.840	1992	1998	2005	rBV5	20178	29293	0.39%	0.035%
45	14.963	2015	2019	2025	rVB4	10586	19244	0.25%	0.023%
46	15.051	2032	2034	2044	rVB10	6950	11835	0.16%	0.014%
47	15.287	2070	2074	2078	rVB	15525	19854	0.26%	0.023%
48	15.334	2078	2082	2086	rBV2	11313	17401	0.23%	0.021%
49	15.410	2088	2095	2108	rVB	3303087	4796693	63.31%	5.656%
50	15.522	2108	2114	2123	rBV	765171	1097789	14.49%	1.294%
51	15.651	2130	2136	2140	rBV	38490	59300	0.78%	0.070%
52	15.840	2163	2168	2169	rBV5	5780	7590	0.10%	0.009%
53	15.863	2169	2172	2176	rVB5	8453	12449	0.16%	0.015%
54	16.098	2207	2212	2216	rBV8	12081	23413	0.31%	0.028%
55	16.204	2226	2230	2239	rVB2	41906	67092	0.89%	0.079%
56	16.316	2245	2249	2254	rVB6	16300	22991	0.30%	0.027%
57	16.410	2261	2265	2272	rBV2	21950	35212	0.46%	0.042%
58	16.569	2287	2292	2298	rVB2	54533	85560	1.13%	0.101%
59	16.634	2300	2303	2309	rVB2	12647	15531	0.20%	0.018%
60	16.687	2309	2312	2316	rBV6	7421	10178	0.13%	0.012%
61	16.745	2316	2322	2327	rBV2	25063	47637	0.63%	0.056%
62	16.845	2336	2339	2345	rVB8	7678	11728	0.15%	0.014%
63	16.916	2345	2351	2354	rBV	48989	78232	1.03%	0.092%
64	16.945	2354	2356	2360	rVB3	18371	21064	0.28%	0.025%
65	17.163	2387	2393	2403	rBV	2578239	3726327	49.18%	4.393%
66	17.263	2403	2410	2417	rVV2	3540526	5031555	66.41%	5.932%
67	17.316	2417	2419	2423	rVV	30579	42690	0.56%	0.050%
68	17.357	2423	2426	2432	rVB7	12239	26008	0.34%	0.031%
69	17.857	2507	2511	2522	rVB	67946	96272	1.27%	0.114%
70	17.945	2522	2526	2528	rBV5	18392	26445	0.35%	0.031%
71	18.075	2542	2548	2557	rBV	578732	787472	10.39%	0.928%
72	18.369	2595	2598	2602	rVB5	14708	17200	0.23%	0.020%
73	18.481	2612	2617	2622	rBV	238601	296084	3.91%	0.349%
74	18.604	2635	2638	2643	rBV7	20684	34332	0.45%	0.040%
75	18.892	2684	2687	2690	rBV3	17780	25560	0.34%	0.030%
76	18.969	2698	2700	2703	rBV4	15619	15012	0.20%	0.018%

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 Stop Thrs : 0 Peak Location: TOP

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77	19.081	2715	2719	2723	rBV3	43772	60702	0.80%	0.072%
78	19.151	2727	2731	2735	rBV	44429	80282	1.06%	0.095%
79	19.310	2753	2758	2769	rBV	432928	588217	7.76%	0.694%
80	19.504	2786	2791	2800	rBV	4176525	5418451	71.51%	6.389%
81	20.057	2882	2885	2889	rVB	64168	62114	0.82%	0.073%
82	20.498	2956	2960	2965	rBV	7095557	7576843	100.00%	8.933%
83	20.880	3022	3025	3029	rBV4	65381	81885	1.08%	0.097%
84	20.998	3040	3045	3053	rBV2	950840	1580478	20.86%	1.863%
85	21.263	3085	3090	3099	rVB	3116760	3838383	50.66%	4.526%
86	21.886	3192	3196	3206	rBV	1458568	2014906	26.59%	2.376%
87	22.927	3368	3373	3377	rBV	715508	1089273	14.38%	1.284%
88	22.969	3377	3380	3389	rVB2	410648	634681	8.38%	0.748%
89	23.457	3456	3463	3473	rVB2	2768469	5322642	70.25%	6.276%
90	23.610	3482	3489	3496	rBV2	1969952	3808466	50.26%	4.490%
91	24.257	3595	3599	3606	rVB	283211	549771	7.26%	0.648%
92	24.345	3609	3614	3623	rVB3	313380	665148	8.78%	0.784%

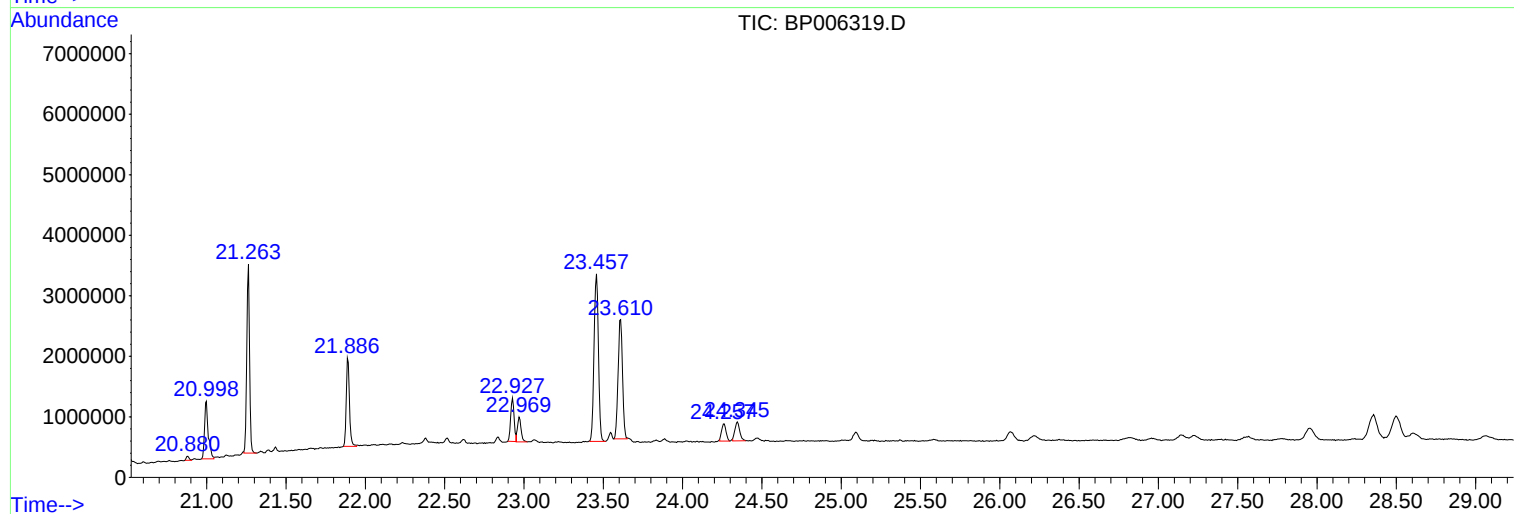
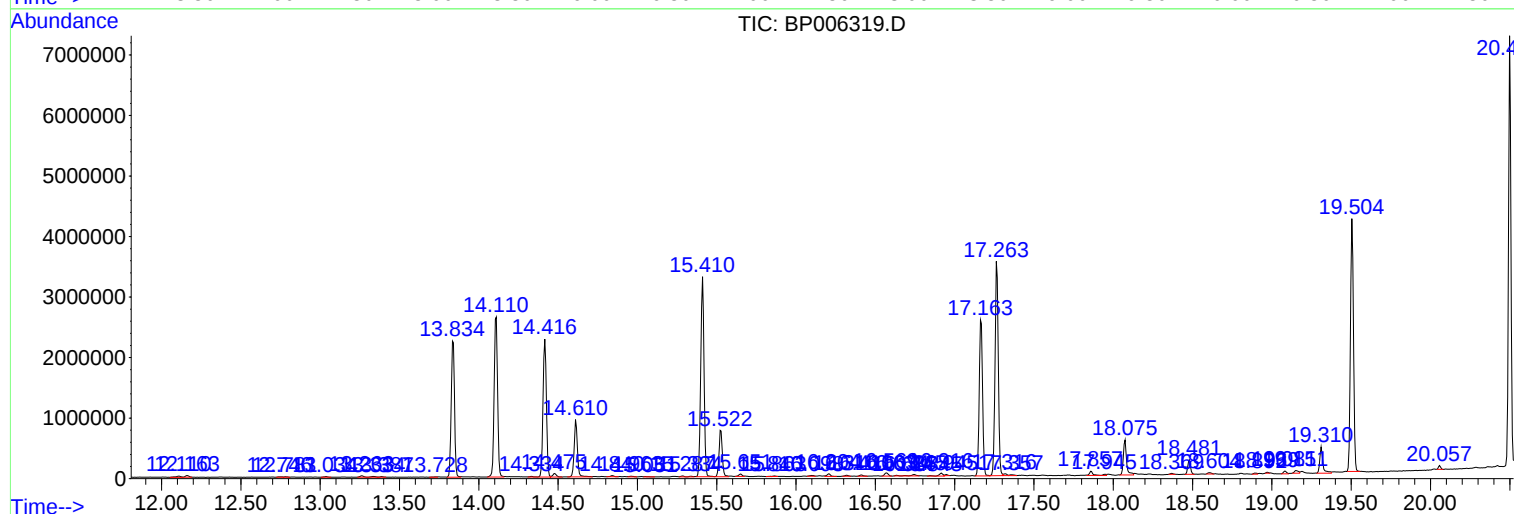
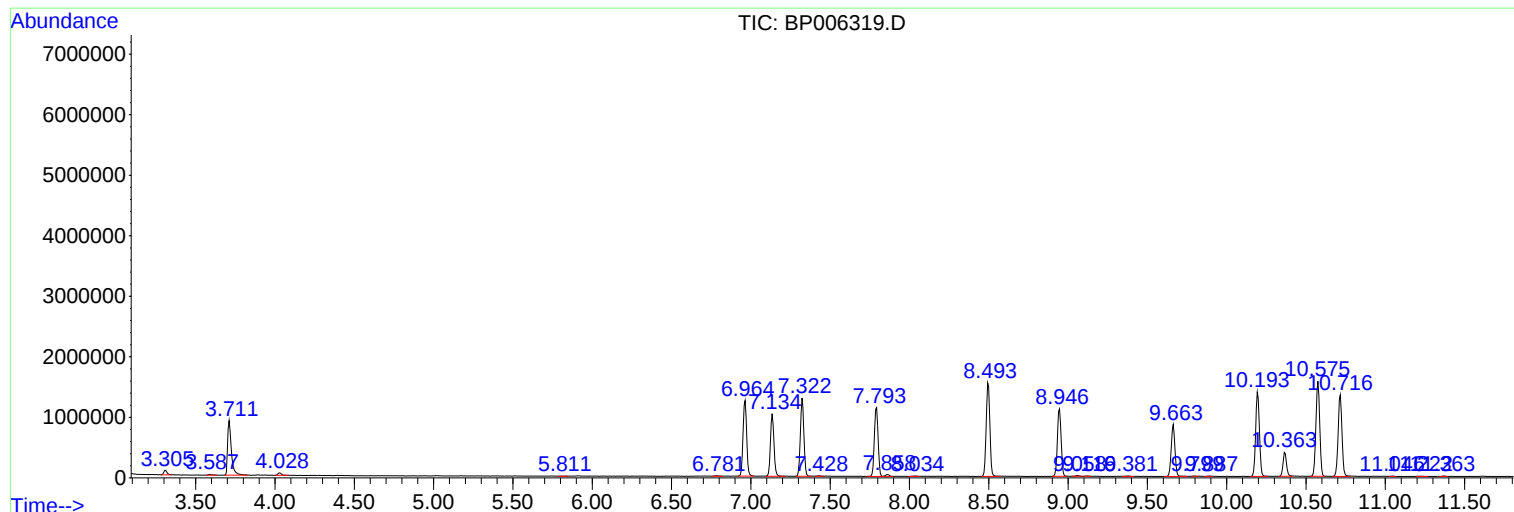
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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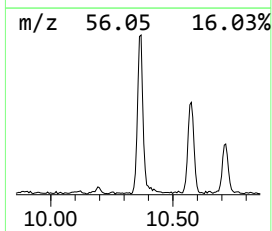
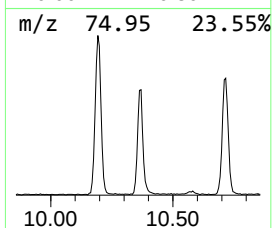
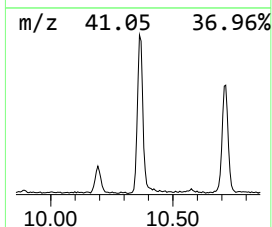
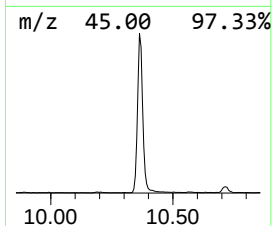
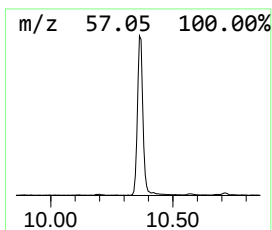
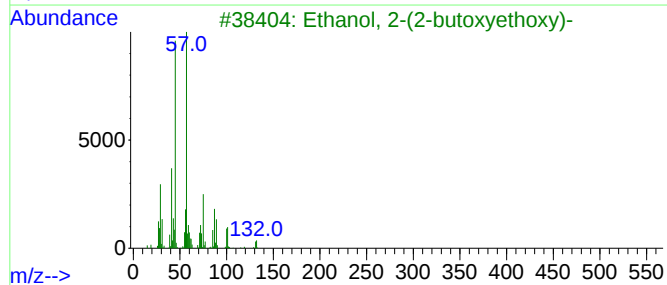
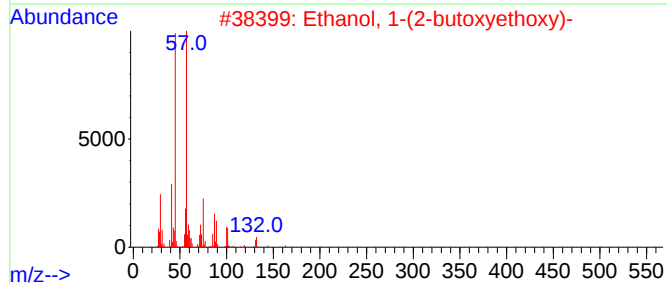
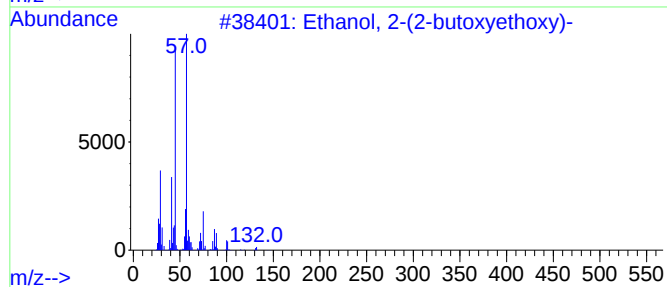
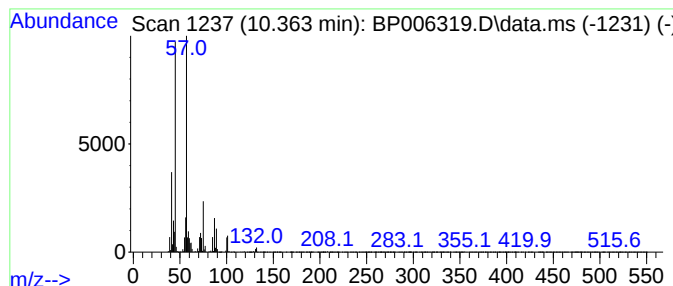
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	4.64 ng/ul	610192	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83



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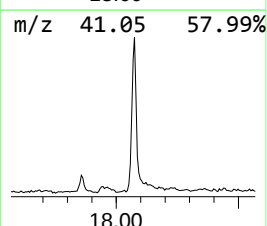
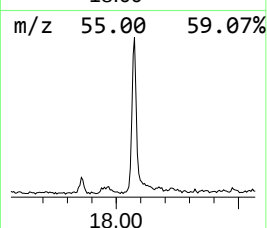
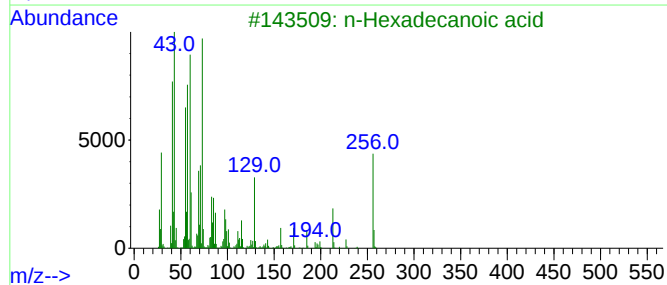
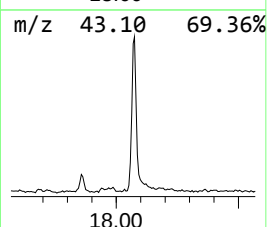
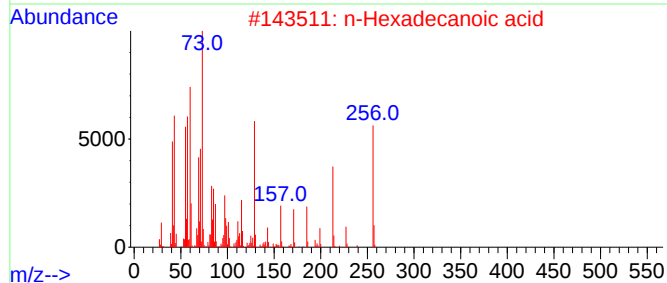
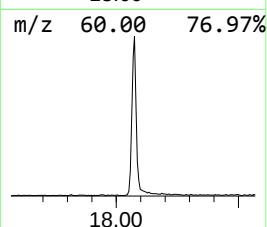
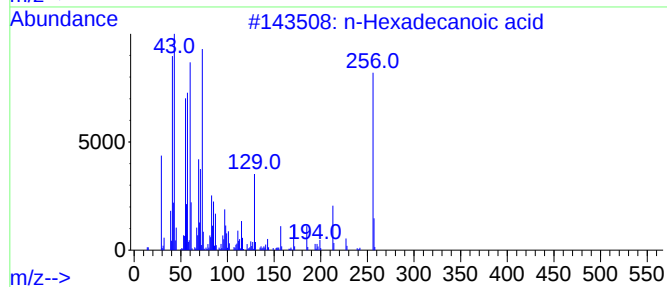
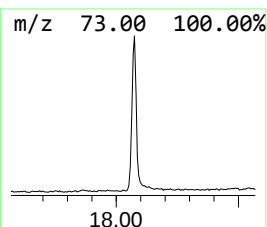
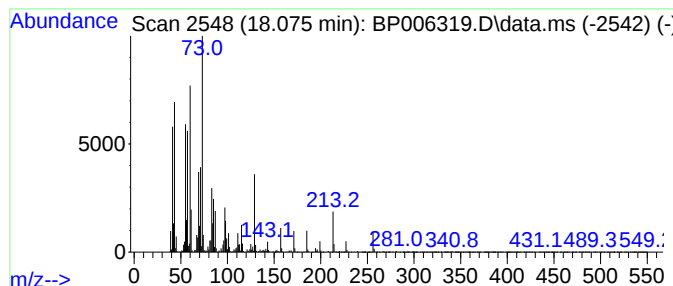
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	4.23 ng/ul	787472	Phenanthrene-d10	17.163

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96



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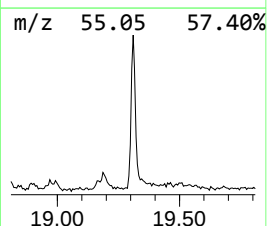
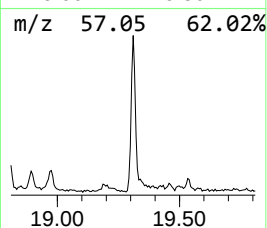
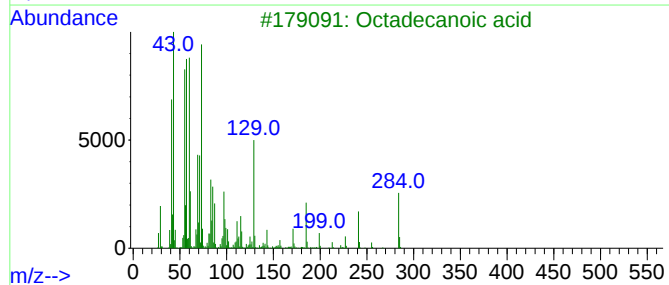
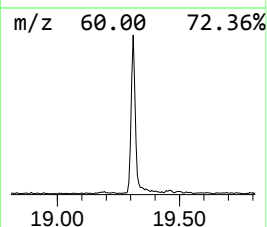
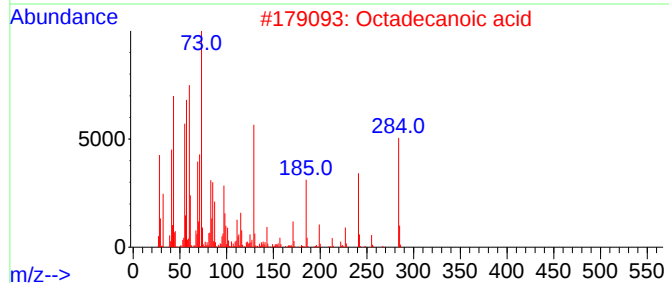
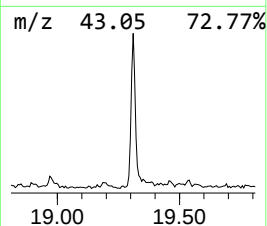
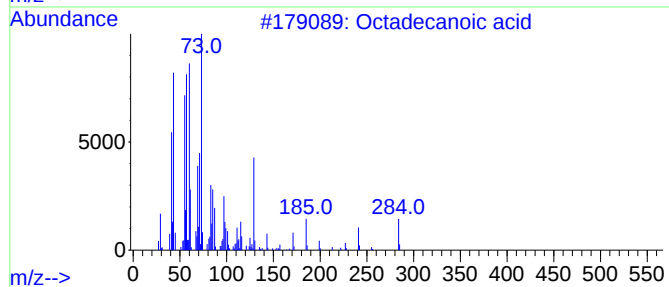
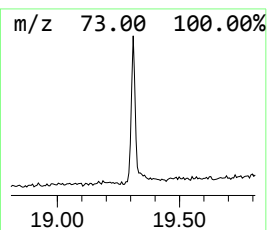
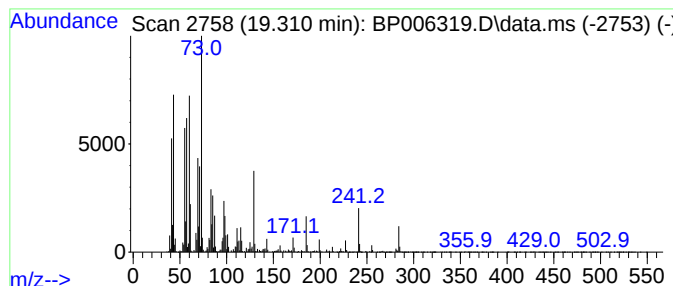
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.310	3.06 ng/ul	588217	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
Operator : CG/JU
Sample : M2973-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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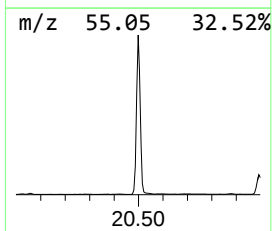
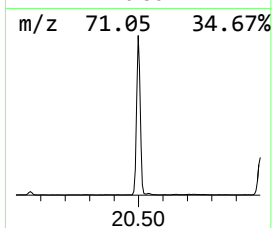
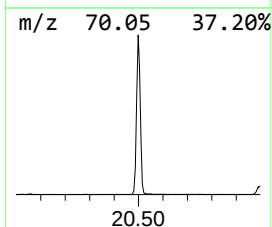
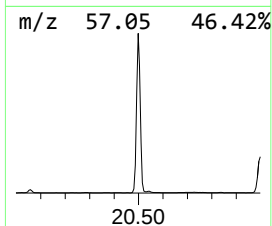
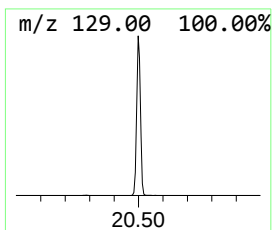
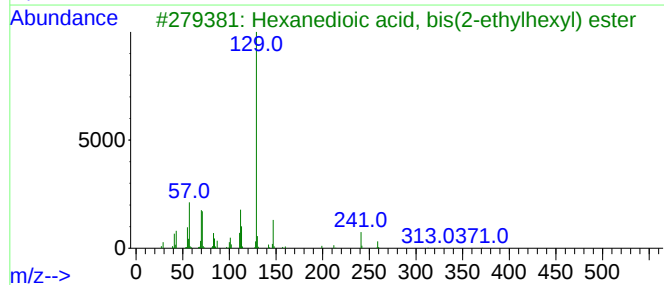
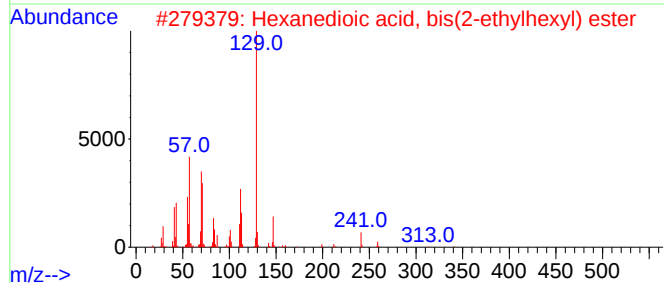
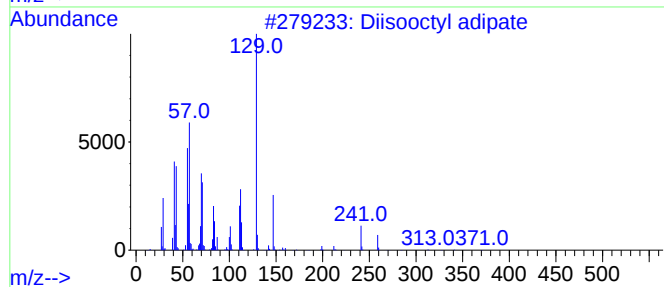
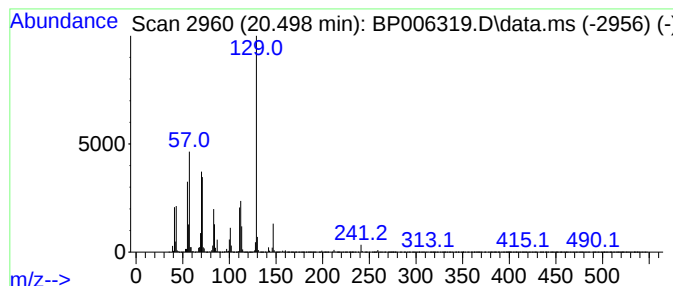
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Diisooctyl adipate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.500	39.48 ng/ul	7576840	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisooctyl adipate	370	C22H42O4	001330-86-5	90
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	83
4			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	64
5			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
Operator : CG/JU
Sample : M2973-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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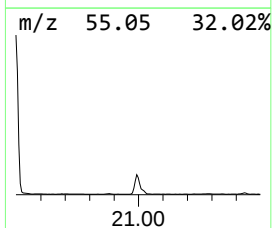
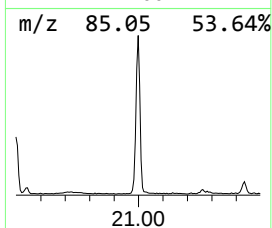
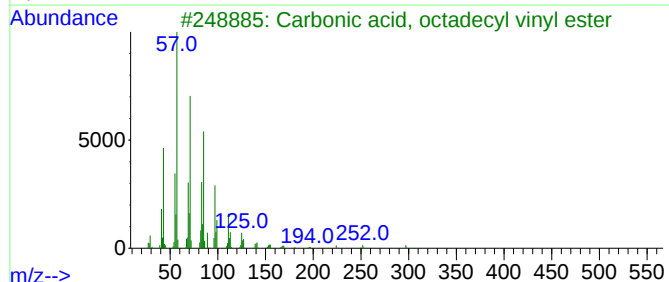
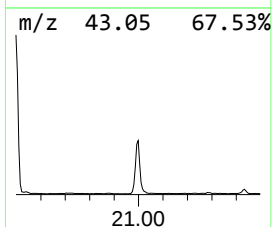
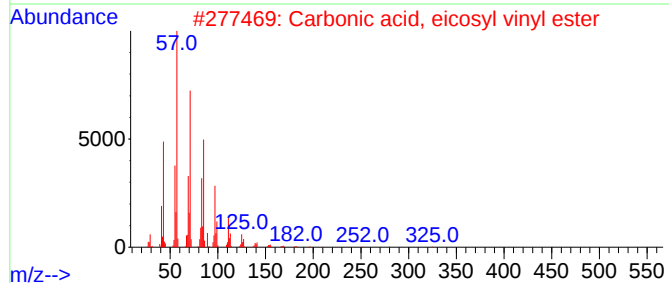
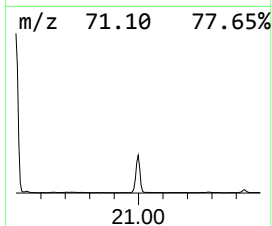
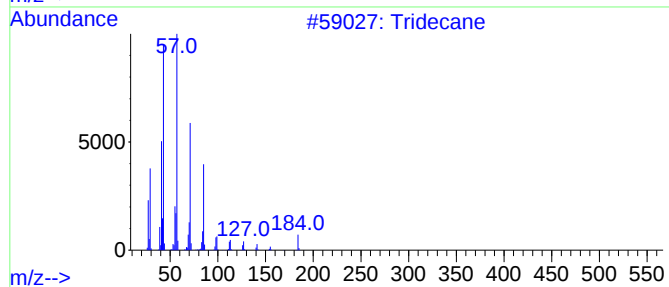
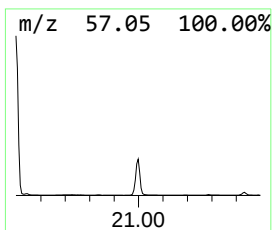
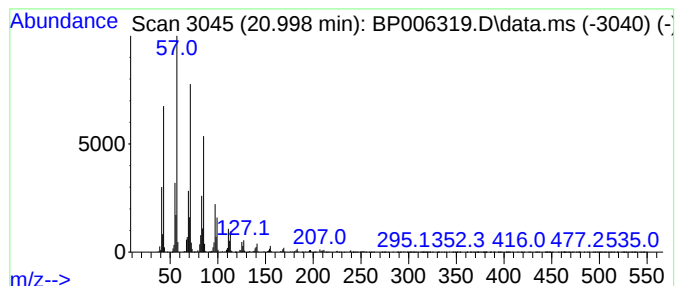
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.000	8.24 ng/ul	1580480	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	91
3			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	91
4			Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	91
5			Tritetracontane	605	C43H88	007098-21-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
Operator : CG/JU
Sample : M2973-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

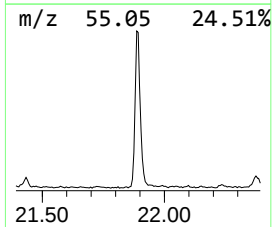
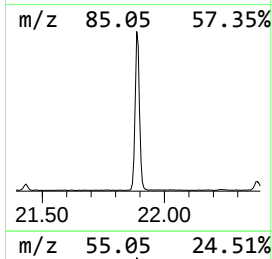
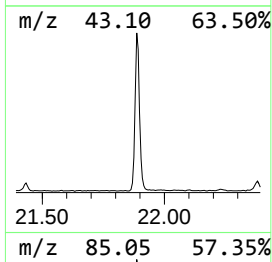
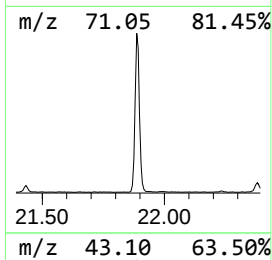
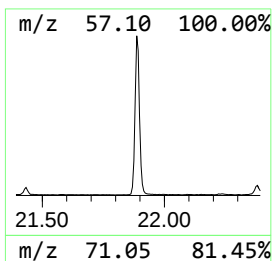
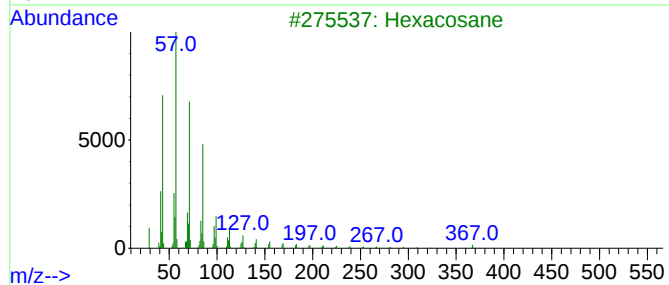
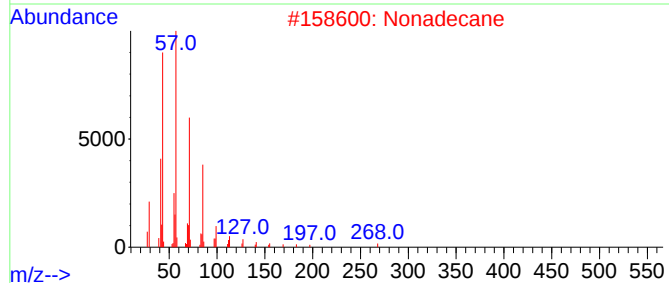
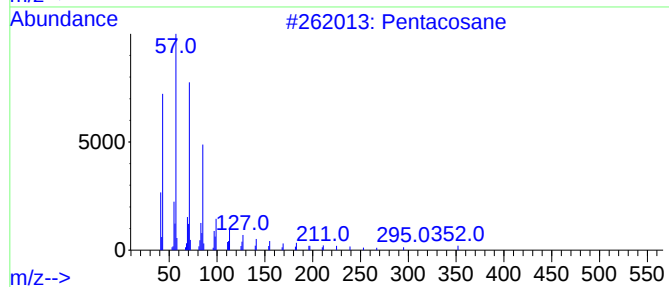
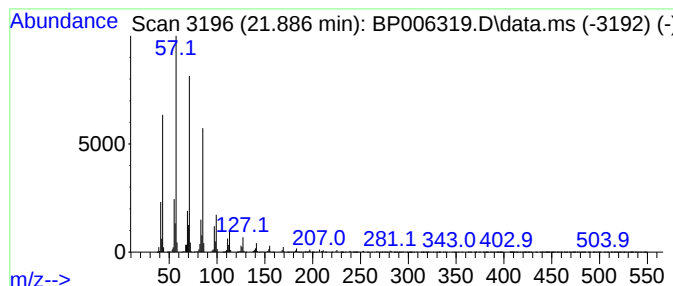
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.890	10.50 ng/ul	2014910	Chrysene-d12	21.263	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	94
2	Nonadecane	268	C19H40	000629-92-5	93
3	Hexacosane	366	C26H54	000630-01-3	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Nonadecane	268	C19H40	000629-92-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
Operator : CG/JU
Sample : M2973-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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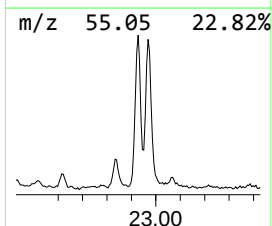
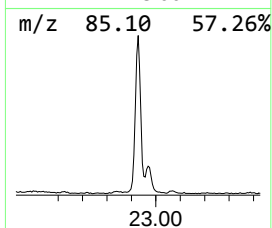
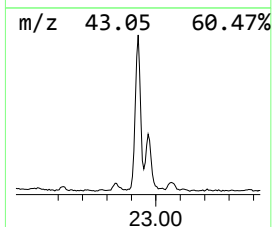
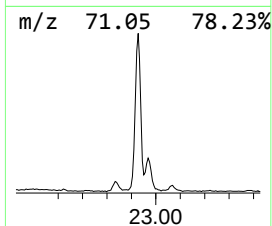
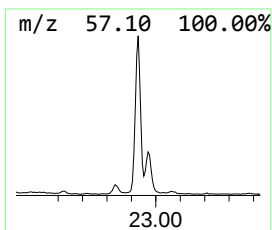
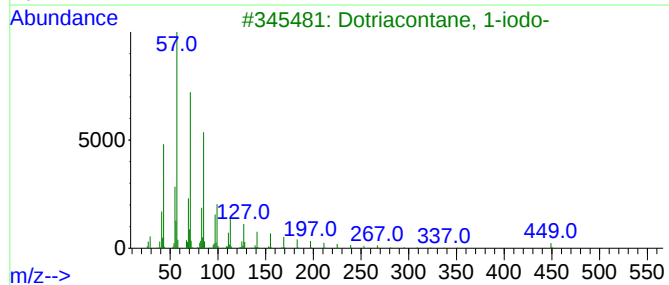
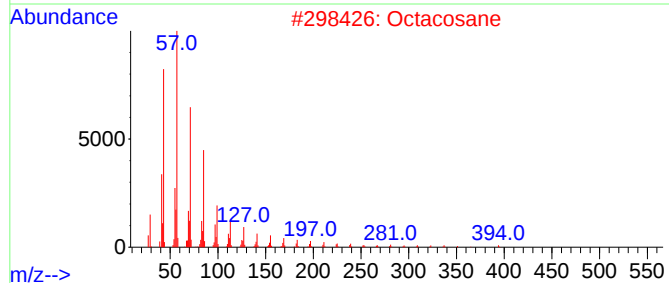
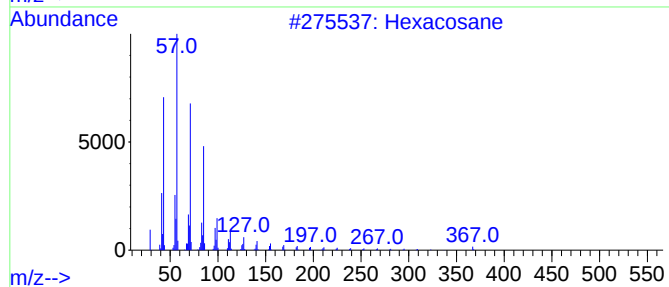
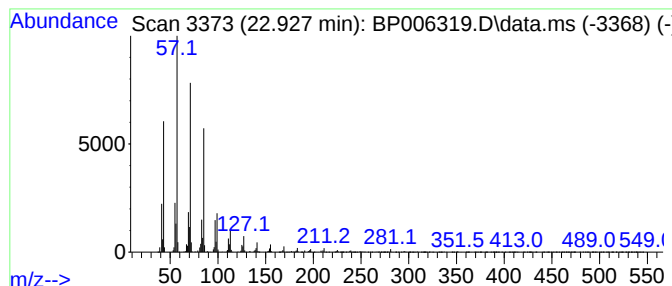
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.930	5.72 ng/ul	1089270	Perylene-d12	23.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Octacosane	394	C28H58	000630-02-4	91
3			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
Operator : CG/JU
Sample : M2973-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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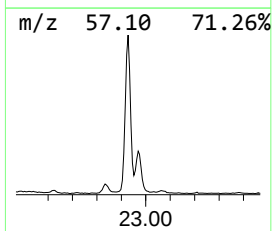
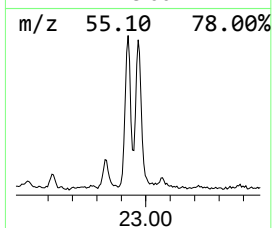
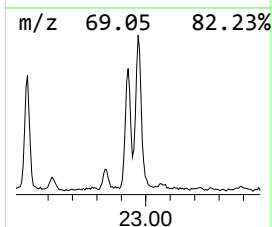
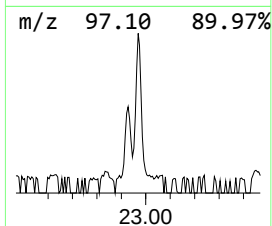
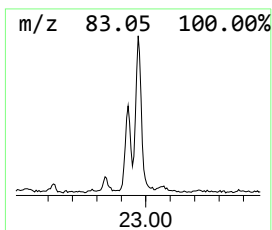
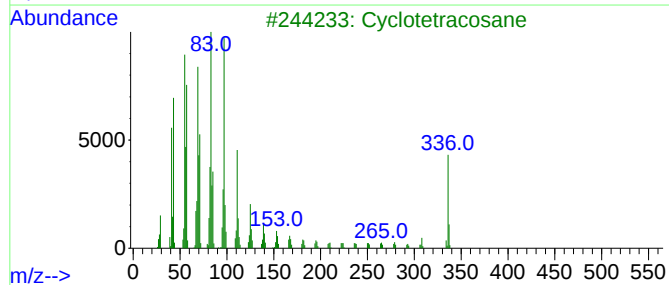
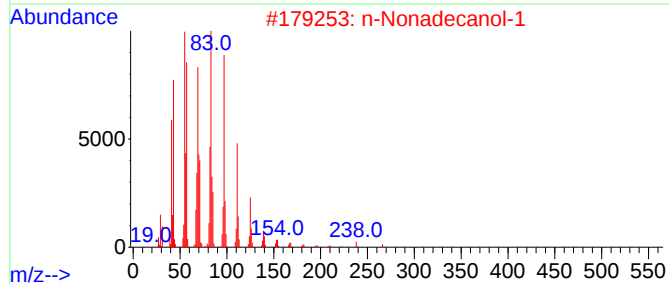
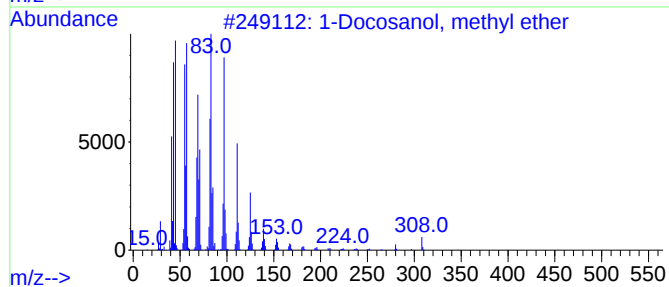
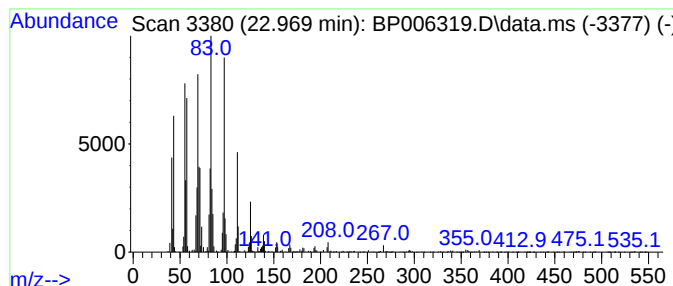
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1-Docosanol, methyl ether Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.970	3.33 ng/ul	634681	Perylene-d12	23.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	95
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3		Cyclotetracosane	336	C24H48	000297-03-0	94
4		Pentacos-1-ene	350	C25H50	016980-85-1	94
5		Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
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Sample : M2973-14
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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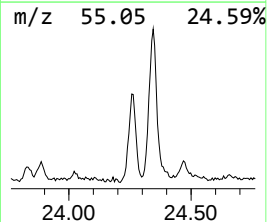
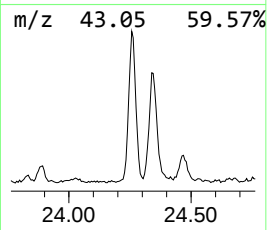
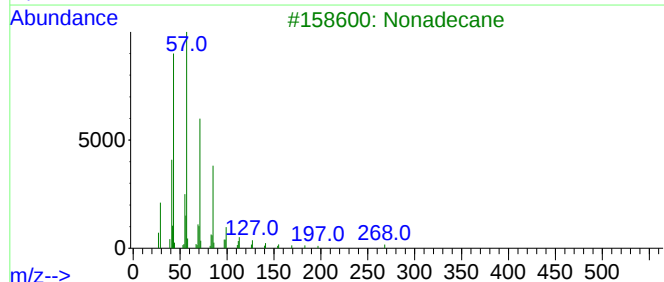
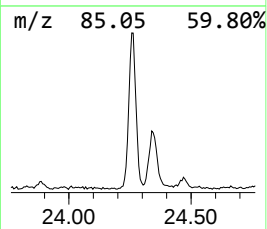
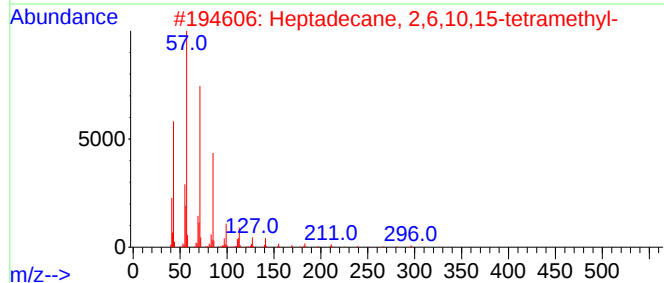
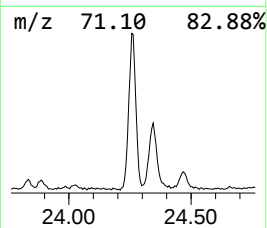
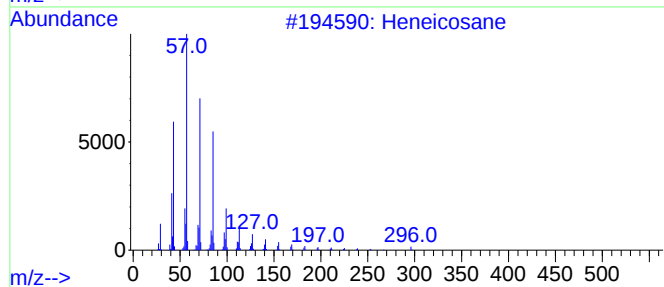
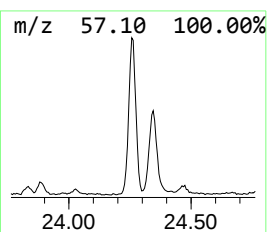
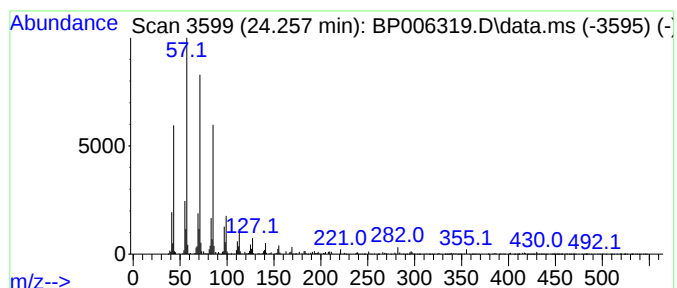
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.260	2.89 ng/ul	549771	Perylene-d12	23.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	99
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
3		Nonadecane	268	C19H40	000629-92-5	93
4		Tetracosane	338	C24H50	000646-31-1	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006319.D
Acq On : 24 Jul 2021 23:03
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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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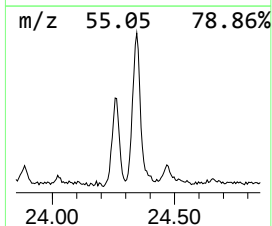
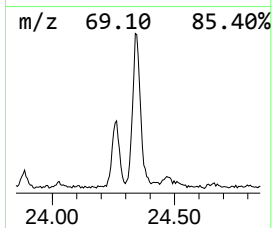
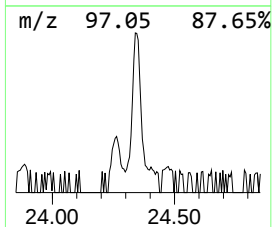
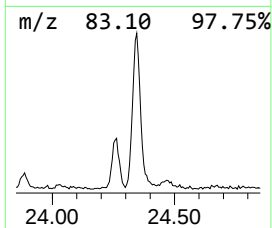
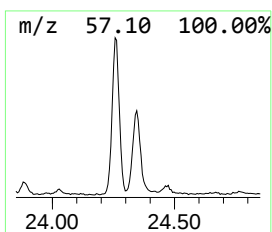
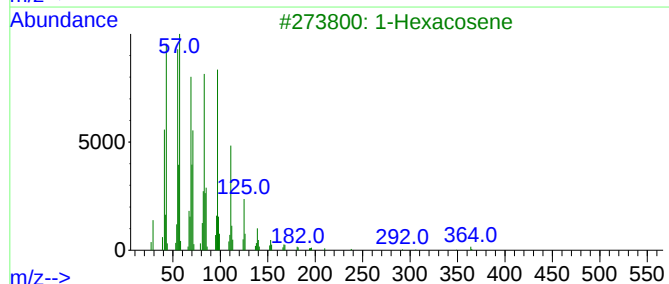
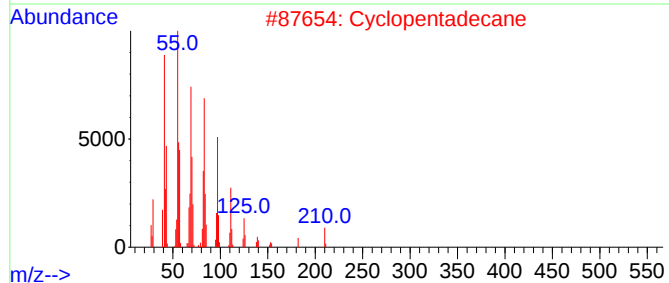
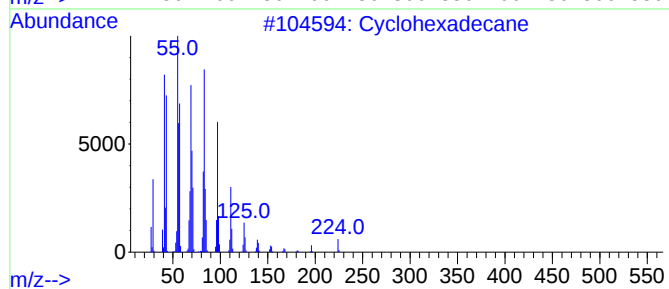
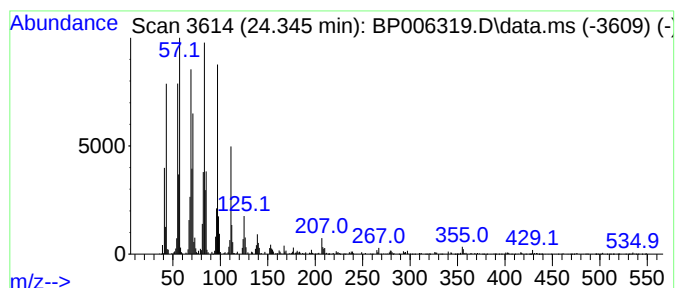
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Cyclic24.35 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.350	3.49 ng/ul	665148	Perylene-d12	23.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	99
2			Cyclopentadecane	210	C15H30	000295-48-7	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
 Data File : BP006319.D
 Acq On : 24 Jul 2021 23:03
 Operator : CG/JU
 Sample : M2973-14
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGE19

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.360	4.6	ng/ul	610192	2	10.575	2629180	20.0
n-Hexadecanoic ...	18.070	4.2	ng/ul	787472	4	17.163	3726330	20.0
Octadecanoic acid	19.310	3.1	ng/ul	588217	5	21.263	3838380	20.0
Diisooctyl adipate	20.500	39.5	ng/ul	7576840	5	21.263	3838380	20.0
(DEL) Alkane: S...	21.000	8.2	ng/ul	1580480	5	21.263	3838380	20.0
(DEL) Alkane: S...	21.890	10.5	ng/ul	2014910	5	21.263	3838380	20.0
(DEL) Alkane: S...	22.930	5.7	ng/ul	1089270	6	23.610	3808470	20.0
1-Docosanol, me...	22.970	3.3	ng/ul	634681	6	23.610	3808470	20.0
(DEL) Alkane: S...	24.260	2.9	ng/ul	549771	6	23.610	3808470	20.0
(DEL) Alkane: C...	24.350	3.5	ng/ul	665148	6	23.610	3808470	20.0