

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
 Data File : BP016616.D
 Acq On : 16 Aug 2023 22:39
 Operator : MA/JU
 Sample : 03967-15
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A48M6

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-BP081423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016616.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	15	18	26	rVB	43393	62935	0.99%	0.063%
2	3.840	91	94	104	rBV	45796	79623	1.25%	0.080%
3	3.946	108	112	120	rVV	42979	67726	1.07%	0.068%
4	4.034	122	127	135	rVB	44823	79327	1.25%	0.080%
5	4.152	143	147	152	rVB	16094	25722	0.41%	0.026%
6	4.734	243	246	254	rVB4	22524	34114	0.54%	0.034%
7	4.987	287	289	295	rVB7	8228	11880	0.19%	0.012%
8	5.093	301	307	311	rBV	35467	54820	0.86%	0.055%
9	5.205	323	326	333	rVB3	12020	15549	0.24%	0.016%
10	6.004	458	462	466	rVB3	10633	13797	0.22%	0.014%
11	6.493	540	545	548	rBV7	6446	11999	0.19%	0.012%
12	6.987	624	629	635	rBV7	7527	14726	0.23%	0.015%
13	7.193	656	664	677	rBV	703678	1227451	19.34%	1.234%
14	7.387	691	697	706	rVB	567687	900785	14.19%	0.905%
15	7.569	722	728	737	rBV	651494	1050921	16.56%	1.056%
16	7.646	737	741	753	rVB2	25247	49152	0.77%	0.049%
17	8.051	802	810	820	rBV	920409	1498231	23.61%	1.506%
18	8.310	847	854	857	rBV8	6544	13480	0.21%	0.014%
19	8.751	922	929	941	rBV	814634	1336052	21.05%	1.343%
20	8.898	948	954	961	rVB	304697	490829	7.73%	0.493%
21	9.034	971	977	983	rVB2	13946	29408	0.46%	0.030%
22	9.251	1003	1014	1024	rBV	669601	1135553	17.89%	1.141%
23	9.481	1048	1053	1057	rVB5	10593	13937	0.22%	0.014%
24	9.963	1127	1135	1142	rBV	519460	859417	13.54%	0.864%
25	10.075	1147	1154	1161	rBV	46554	91929	1.45%	0.092%
26	10.328	1192	1197	1204	rBV	25211	41479	0.65%	0.042%
27	10.487	1213	1224	1238	rBV	844514	1466916	23.11%	1.474%
28	10.887	1284	1292	1299	rBV	1434876	2445634	38.53%	2.458%
29	10.939	1299	1301	1309	rVB6	13576	20618	0.32%	0.021%
30	11.039	1311	1318	1337	rBV	319117	649262	10.23%	0.652%
31	11.422	1376	1383	1393	rBV	17125	33292	0.52%	0.033%
32	11.628	1414	1418	1421	rBV6	8225	11990	0.19%	0.012%
33	11.692	1425	1429	1438	rVB	14784	30276	0.48%	0.030%
34	11.898	1456	1464	1466	rBV9	9162	18398	0.29%	0.018%
35	12.263	1517	1526	1533	rBV6	11314	29235	0.46%	0.029%
36	12.457	1555	1559	1567	rVV4	16080	28111	0.44%	0.028%

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Integrator: RTE

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
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37	12.534	1567	1572	1576	rVB7	8274	13783	0.22%	0.014%
38	12.757	1605	1610	1616	rVB8	17185	28604	0.45%	0.029%
39	13.039	1653	1658	1663	rBV4	12874	25184	0.40%	0.025%
40	13.169	1677	1680	1688	rVB4	20353	28863	0.45%	0.029%
41	13.469	1726	1731	1735	rBV3	13481	22933	0.36%	0.023%
42	13.551	1739	1745	1752	rVB	55671	96809	1.53%	0.097%
43	13.639	1754	1760	1762	rBV7	9039	15608	0.25%	0.016%
44	13.881	1798	1801	1805	rVB6	11072	15303	0.24%	0.015%
45	14.016	1821	1824	1829	rVB7	6819	11795	0.19%	0.012%
46	14.110	1834	1840	1846	rBV	1739549	2373640	37.40%	2.385%
47	14.404	1883	1890	1904	rBV	1950982	2945968	46.42%	2.961%
48	14.628	1924	1928	1932	rVB4	20592	27683	0.44%	0.028%
49	14.704	1932	1941	1949	rBV	2371694	3506649	55.25%	3.524%
50	14.763	1949	1951	1957	rVB6	13411	21278	0.34%	0.021%
51	14.898	1968	1974	1989	rBV	678089	1180633	18.60%	1.186%
52	15.080	2000	2005	2011	rBV5	52734	100886	1.59%	0.101%
53	15.151	2015	2017	2024	rVB8	14887	22669	0.36%	0.023%
54	15.298	2039	2042	2047	rBV6	10180	18339	0.29%	0.018%
55	15.457	2062	2069	2070	rBV7	10954	18003	0.28%	0.018%
56	15.492	2072	2075	2084	rVB7	17590	34275	0.54%	0.034%
57	15.698	2099	2110	2117	rBV	2623468	3898307	61.42%	3.918%
58	15.816	2124	2130	2137	rBV	569373	832652	13.12%	0.837%
59	15.939	2147	2151	2152	rBV3	12581	16061	0.25%	0.016%
60	16.033	2164	2167	2177	rVB3	14989	30130	0.47%	0.030%
61	16.363	2219	2223	2228	rVV2	25196	41843	0.66%	0.042%
62	16.498	2243	2246	2249	rBV5	12326	17792	0.28%	0.018%
63	16.680	2274	2277	2279	rBV3	13722	16259	0.26%	0.016%
64	17.163	2355	2359	2362	rBV	30899	41045	0.65%	0.041%
65	17.198	2363	2365	2370	rVB5	23630	28171	0.44%	0.028%
66	17.322	2383	2386	2391	rBV7	30026	51691	0.81%	0.052%
67	17.469	2403	2411	2416	rBV	3294122	4754093	74.91%	4.778%
68	17.510	2416	2418	2422	rVV	394077	466023	7.34%	0.468%
69	17.569	2422	2428	2437	rVV	3065428	4319381	68.06%	4.341%
70	17.904	2478	2485	2488	rBV2	91150	146256	2.30%	0.147%
71	17.980	2495	2498	2503	rVB5	26621	32728	0.52%	0.033%
72	18.051	2505	2510	2514	rVV2	110882	185777	2.93%	0.187%
73	18.251	2541	2544	2549	rBV2	114726	147417	2.32%	0.148%
74	18.310	2549	2554	2562	rVV2	1457596	2244275	35.36%	2.255%
75	18.369	2562	2564	2573	rVB2	148059	272491	4.29%	0.274%
76	18.504	2583	2587	2592	rBV3	173142	289072	4.55%	0.291%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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77	18.551	2592	2595	2600	rVV3	90641	150107	2.37%	0.151%
78	18.810	2636	2639	2644	rVB	71838	107483	1.69%	0.108%
79	18.863	2645	2648	2652	rVV3	63210	97049	1.53%	0.098%
80	18.898	2652	2654	2659	rVV3	67430	93960	1.48%	0.094%
81	18.945	2660	2662	2667	rVB5	40403	48891	0.77%	0.049%
82	19.121	2688	2692	2696	rBV	91794	144458	2.28%	0.145%
83	19.186	2700	2703	2710	rBV2	94781	215280	3.39%	0.216%
84	19.468	2742	2751	2756	rVB	979352	1555444	24.51%	1.563%
85	19.539	2759	2763	2774	rVB	449412	673757	10.62%	0.677%
86	19.798	2801	2807	2811	rBV	4484768	6346562	100.00%	6.378%
87	20.263	2882	2886	2890	rBV	314408	435940	6.87%	0.438%
88	21.215	3043	3048	3055	rBV3	1826148	3104263	48.91%	3.120%
89	21.562	3101	3107	3111	rBV	4255708	6078242	95.77%	6.108%
90	21.604	3111	3114	3118	rVB	529905	646994	10.19%	0.650%
91	22.145	3202	3206	3209	rBV	1250650	1715157	27.02%	1.724%
92	22.180	3209	3212	3223	rVB	1654325	2727208	42.97%	2.741%
93	22.957	3340	3344	3351	rVB	983456	1465723	23.09%	1.473%
94	23.274	3392	3398	3404	rBV	2839376	4786949	75.43%	4.811%
95	23.345	3404	3410	3423	rVB2	2035726	4721305	74.39%	4.745%
96	23.974	3510	3517	3523	rVB2	2614866	5738639	90.42%	5.767%
97	24.145	3539	3546	3554	rBV	2648171	5594818	88.16%	5.623%
98	24.356	3576	3582	3591	rVB	1466029	2973120	46.85%	2.988%
99	24.751	3642	3649	3661	rVB	1700675	4035496	63.59%	4.055%
100	24.880	3664	3671	3686	rVV	1320376	3765847	59.34%	3.784%

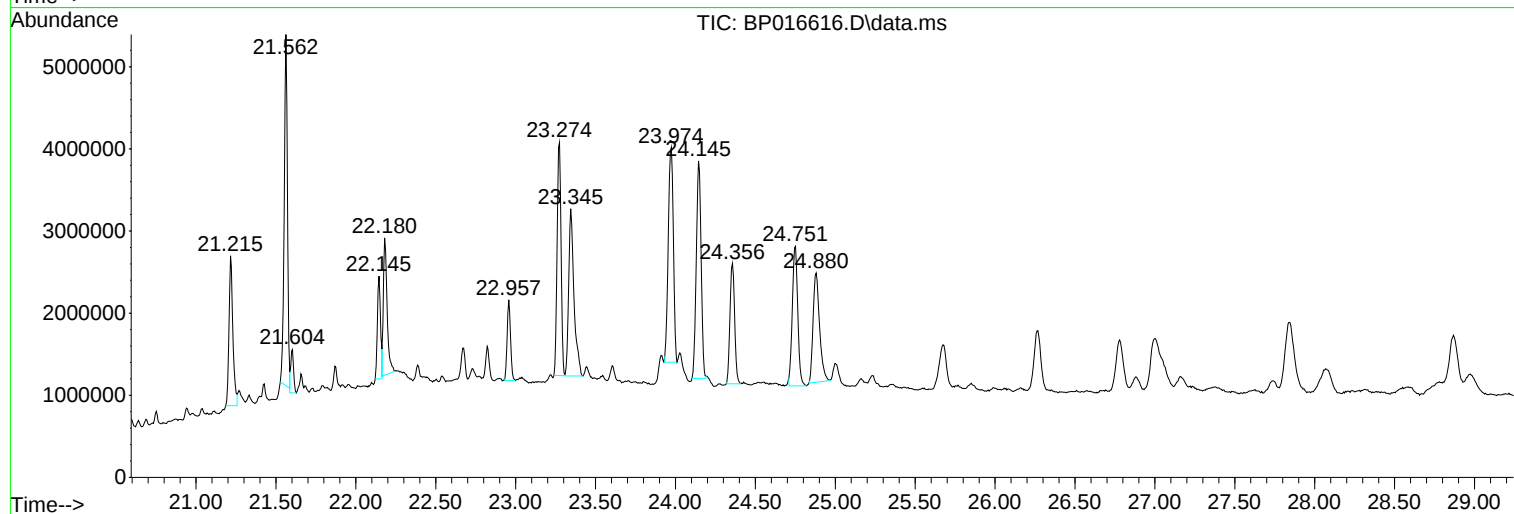
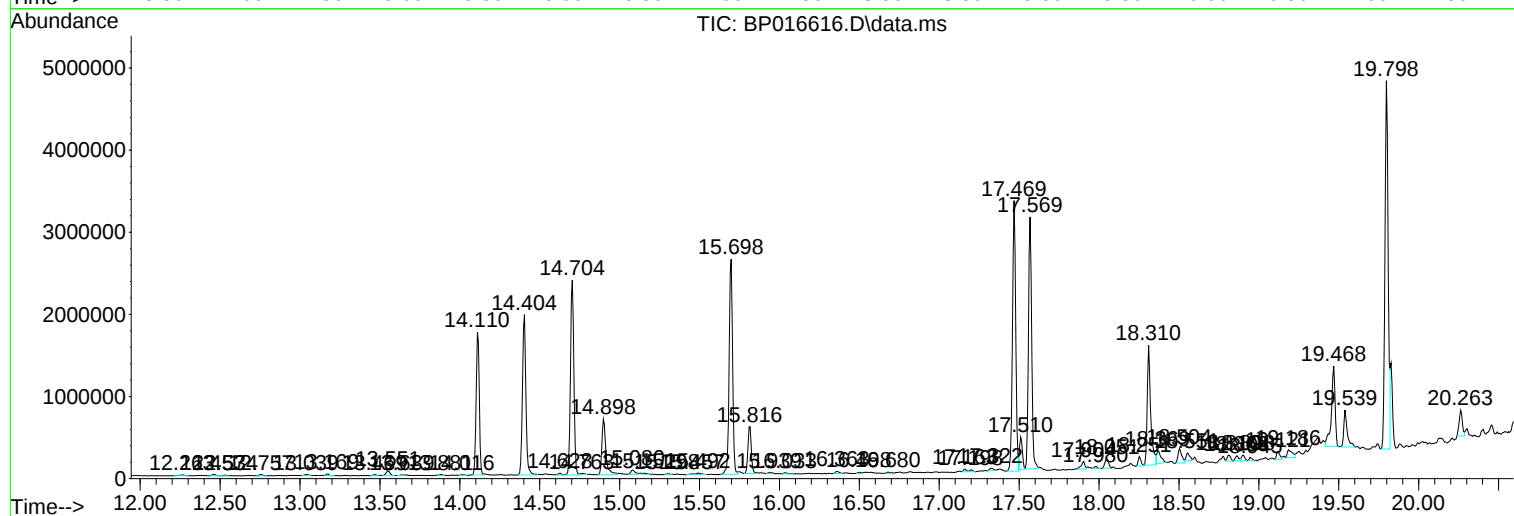
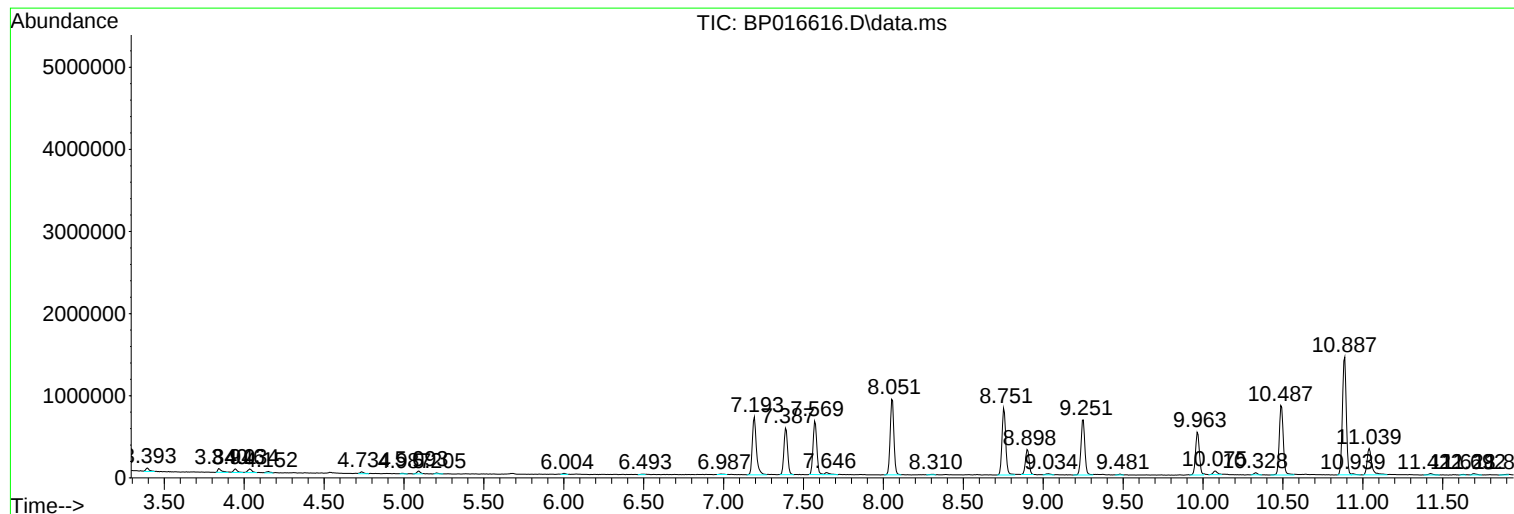
Sum of corrected areas: 99507635

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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



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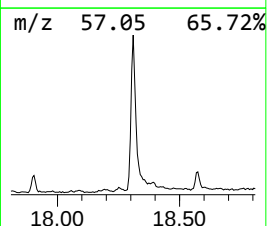
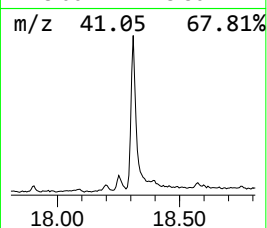
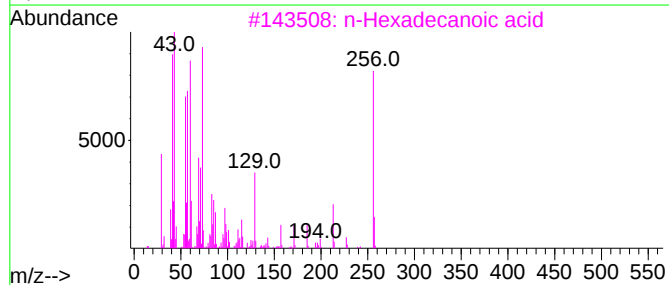
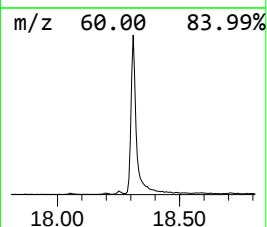
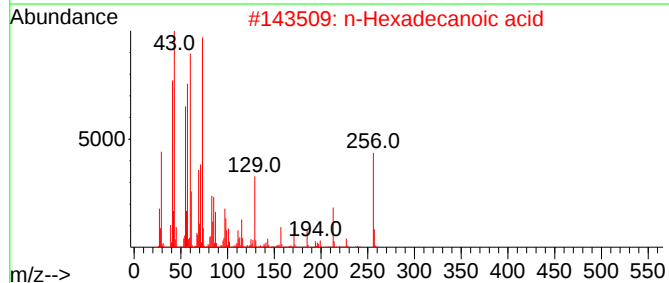
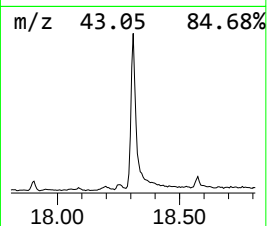
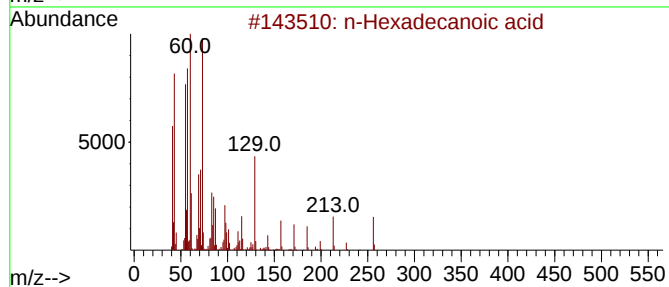
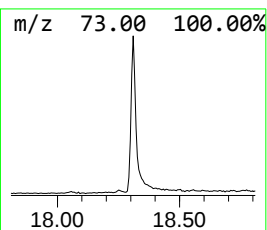
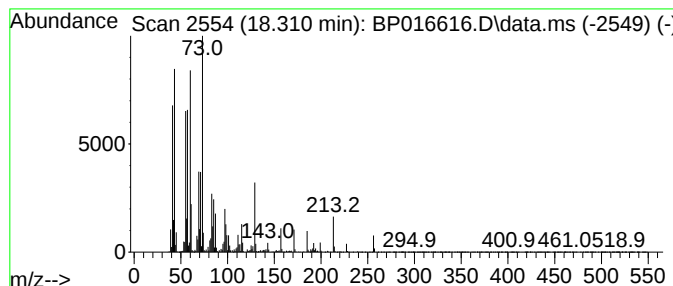
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.310	9.44 ng/ul	2244280	Phenanthrene-d10	17.469

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	97
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93



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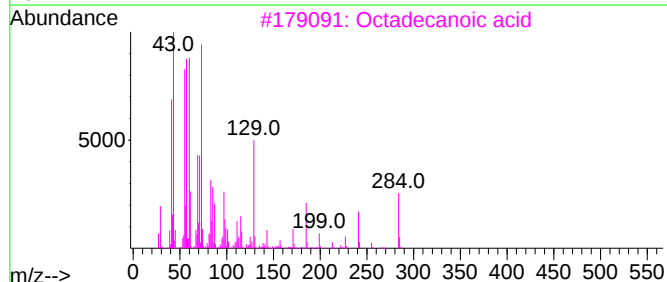
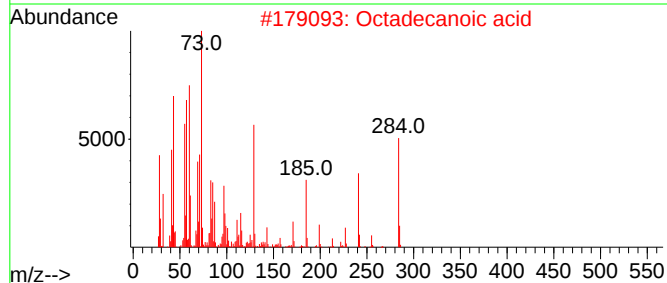
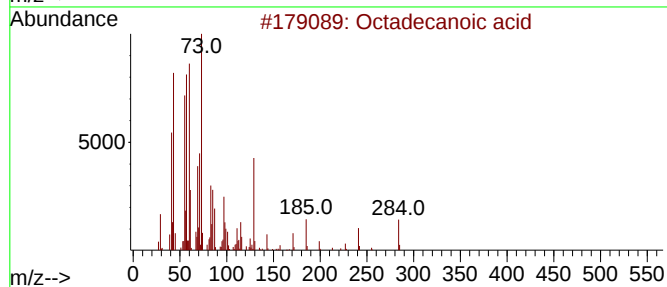
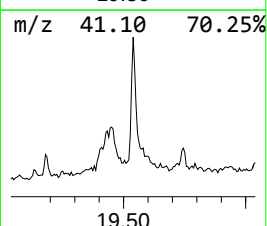
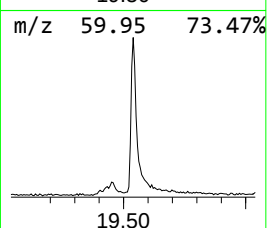
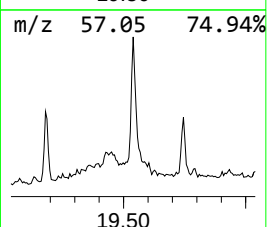
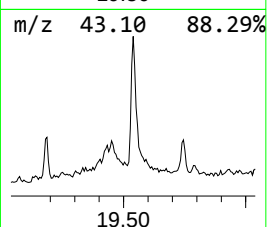
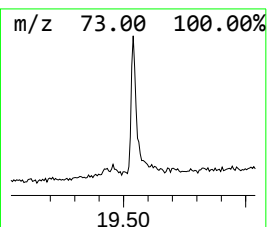
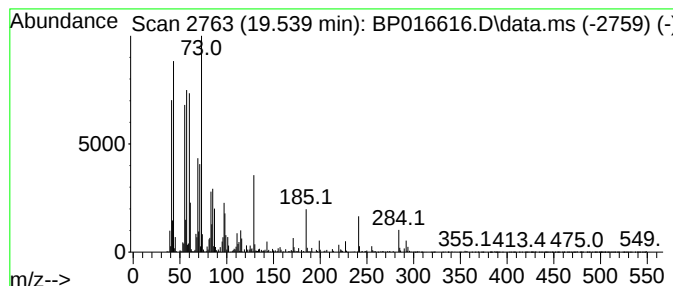
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	2.22 ng/ul	673757	Chrysene-d12	21.562

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	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



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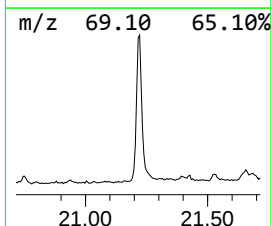
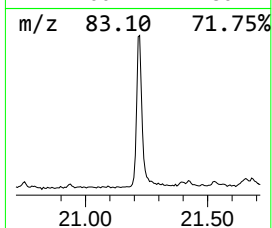
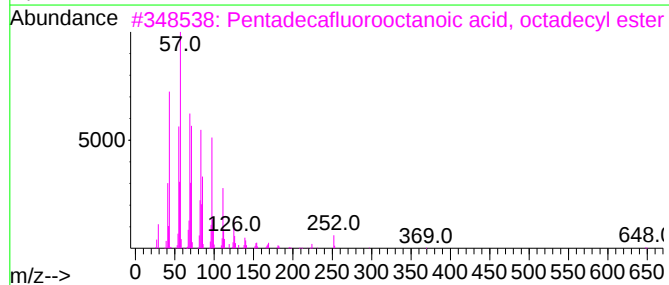
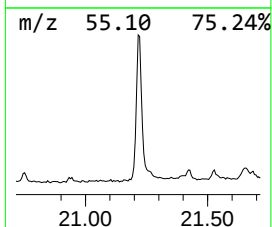
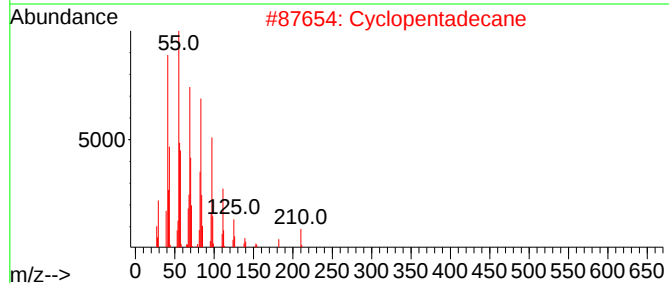
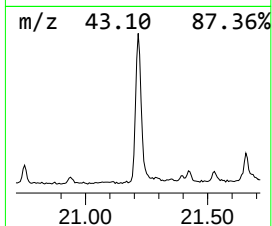
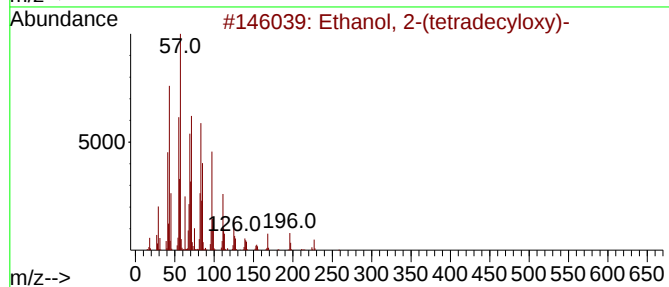
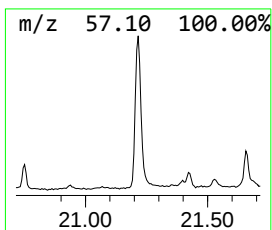
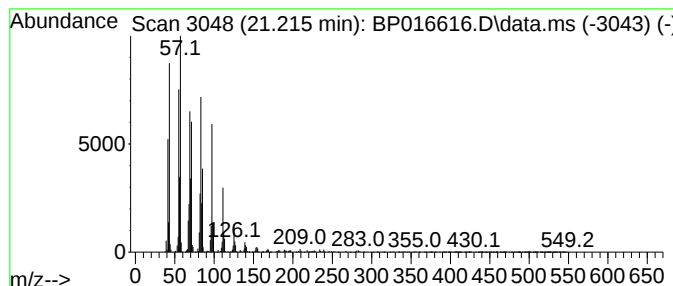
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(tetradecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.215	10.21 ng/ul	3104260	Chrysene-d12	21.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	96
2			Cyclopentadecane	210	C15H30	000295-48-7	96
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
4			9-Tricosene, (Z)-	322	C23H46	027519-02-4	93
5			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016616.D
Acq On : 16 Aug 2023 22:39
Operator : MA/JU
Sample : 03967-15
Misc :
ALS Vial : 19 Sample Multiplier: 1

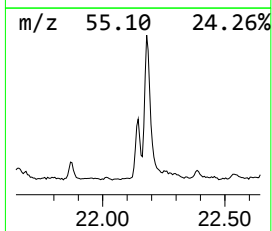
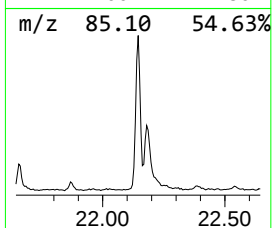
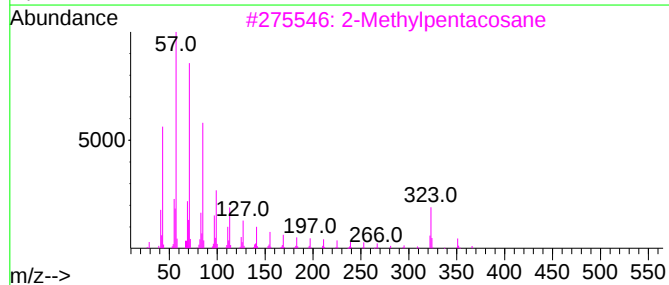
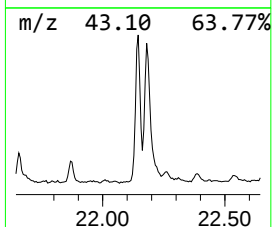
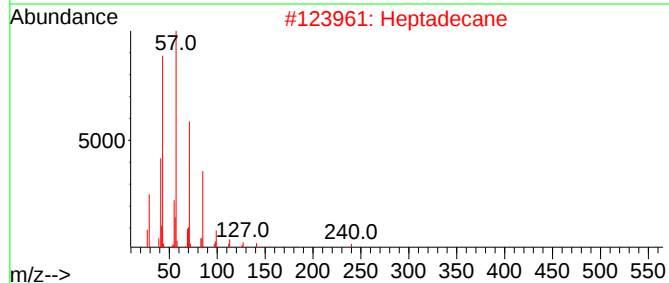
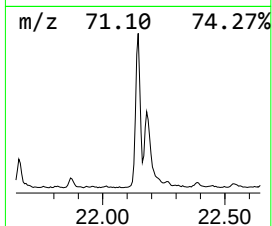
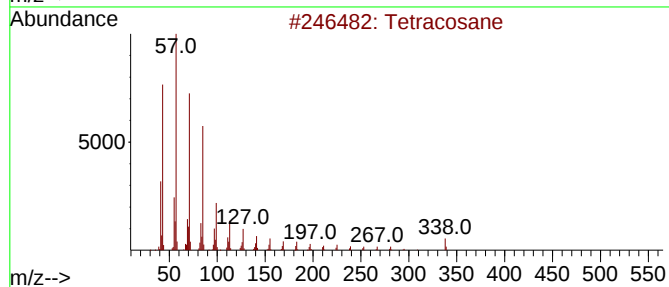
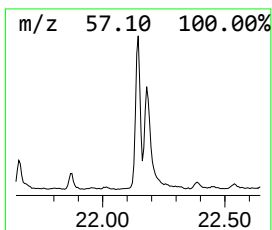
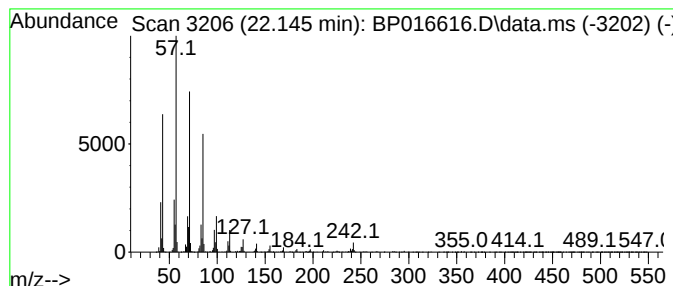
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.145	5.64 ng/ul	1715160	Chrysene-d12	21.562	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	96
2	Heptadecane	240	C17H36	000629-78-7	94
3	2-Methylpentacosane	366	C26H54	000629-87-8	93
4	Heneicosane	296	C21H44	000629-94-7	91
5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016616.D
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ClientSampleId :
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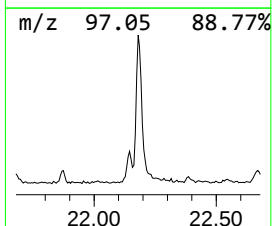
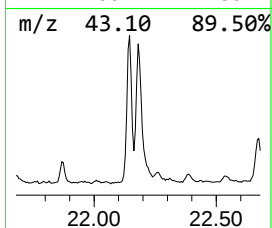
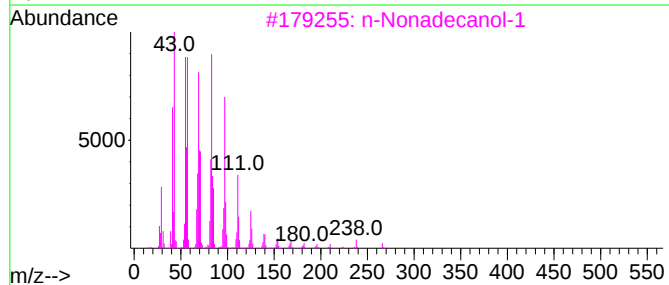
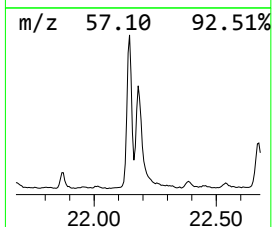
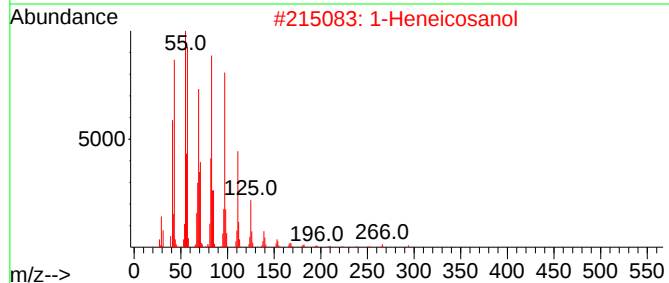
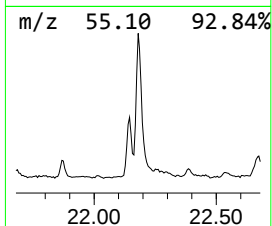
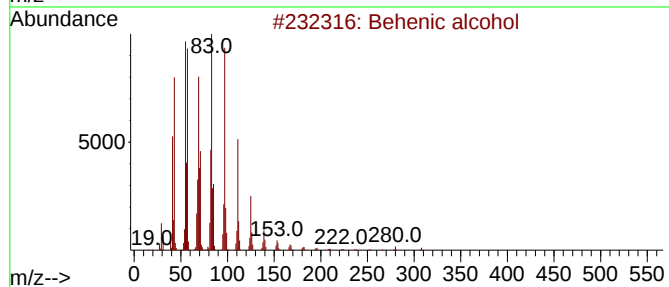
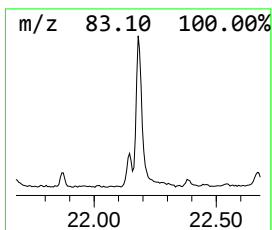
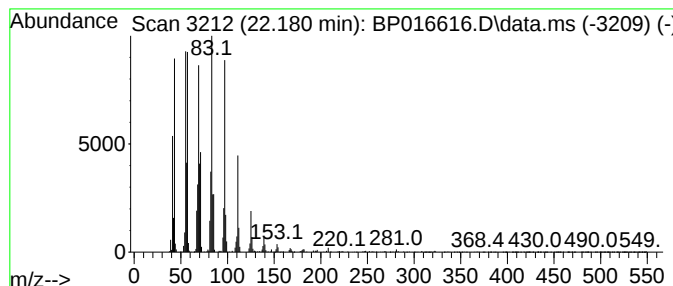
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Behenic alcohol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.180	8.97 ng/ul	2727210	Chrysene-d12	21.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	93
4	9-Tricosene, (Z)-	322	C23H46	027519-02-4	91
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016616.D
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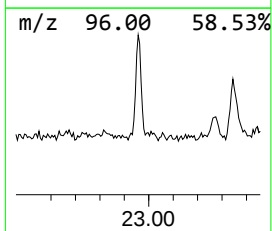
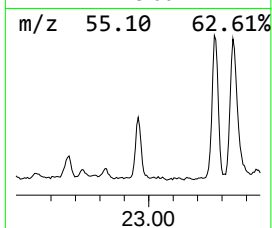
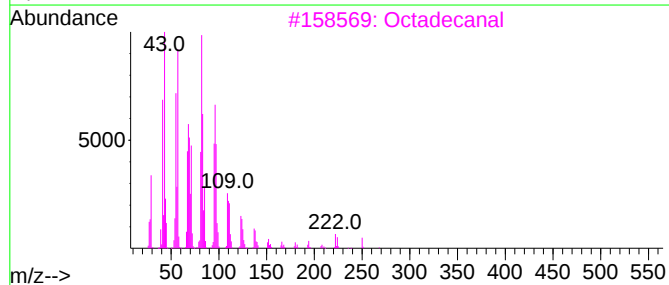
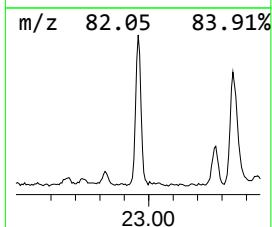
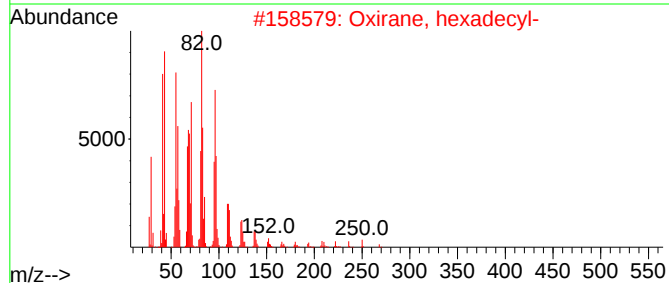
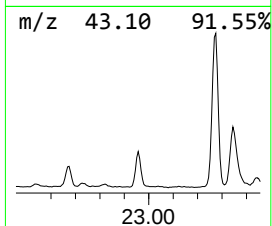
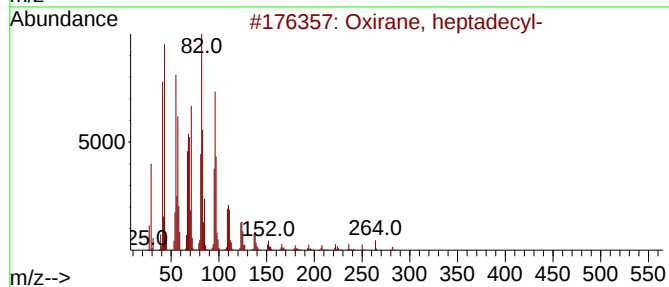
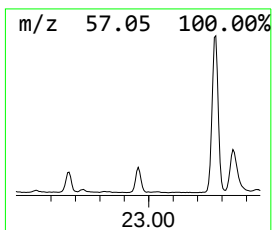
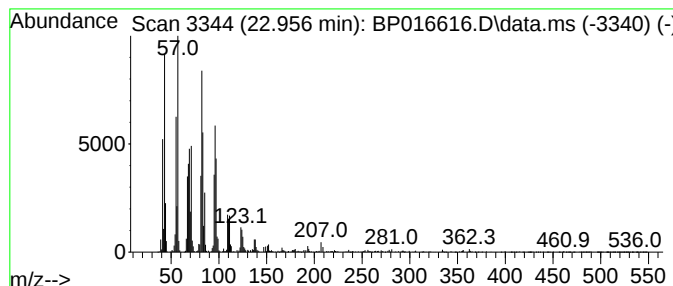
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Oxirane, heptadecyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.956	5.24 ng/ul	1465720	Perylene-d12		24.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Oxirane, heptadecyl-		282	C19H38O	067860-04-2	94
2	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	94
3	Octadecanal		268	C18H36O	000638-66-4	91
4	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	91
5	Octadecanal		268	C18H36O	000638-66-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
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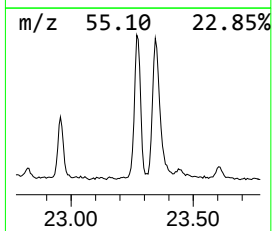
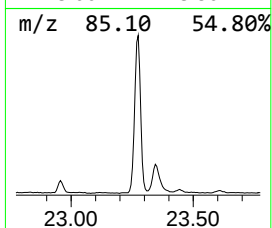
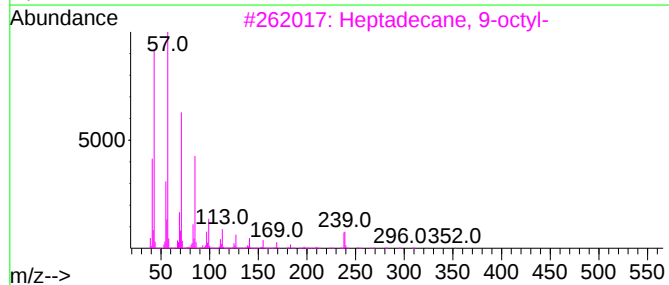
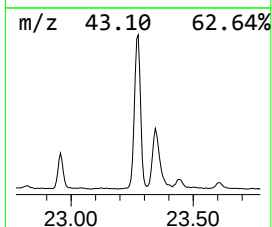
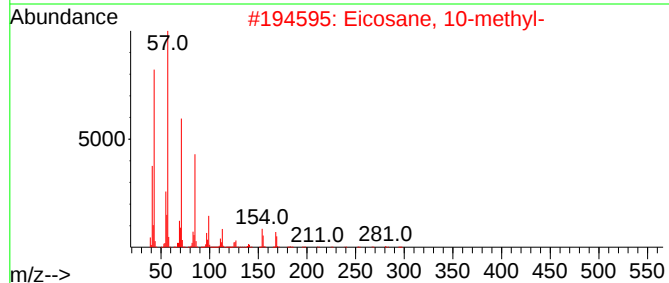
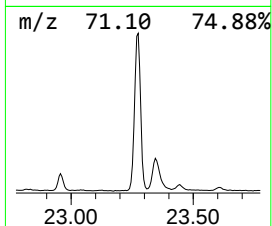
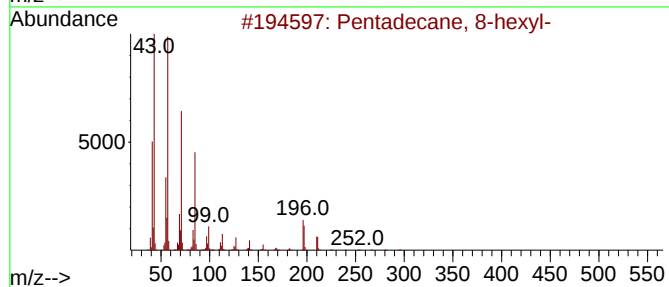
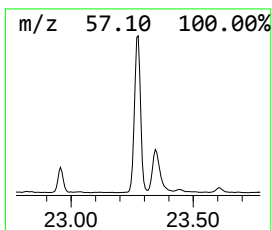
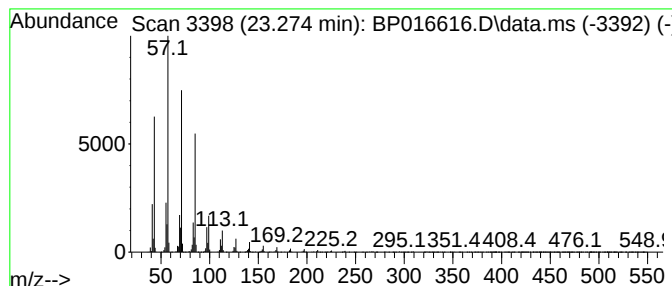
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.274	17.11 ng/ul	4786950	Perylene-d12	24.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	92
4	Heneicosane	296	C21H44	000629-94-7	91
5	Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
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Acq On : 16 Aug 2023 22:39
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ALS Vial : 19 Sample Multiplier: 1

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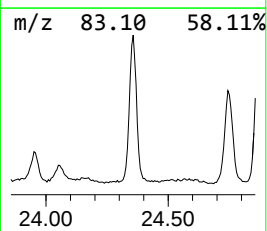
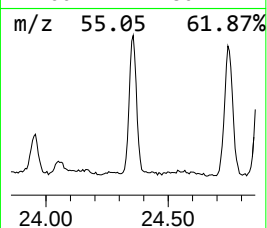
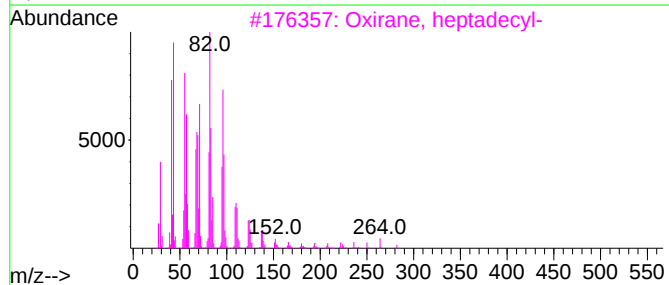
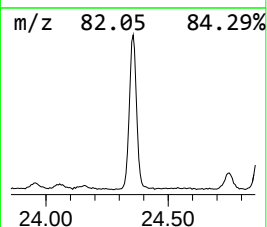
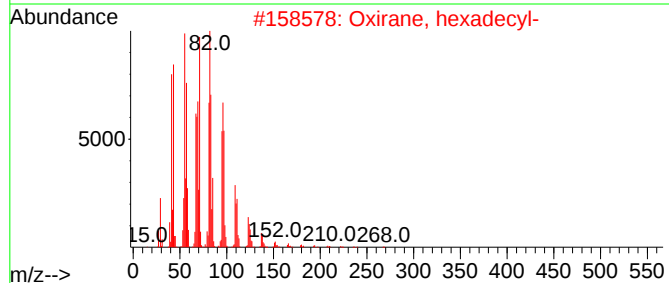
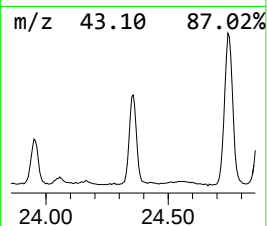
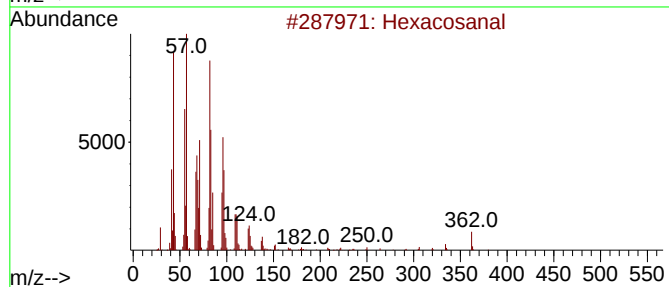
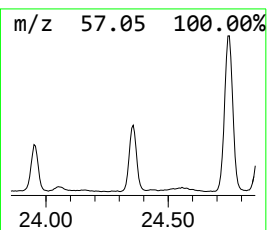
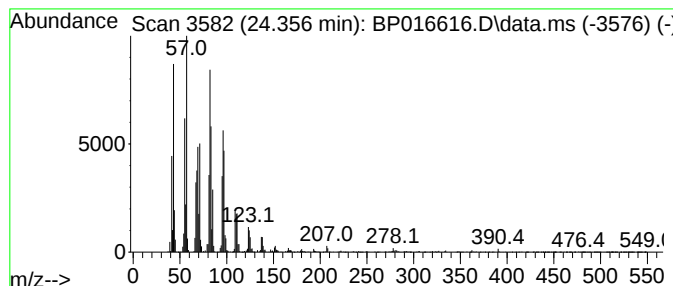
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexacosanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.356	10.63 ng/ul	2973120	Perylene-d12	24.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosanal	380	C26H52O	026627-85-0	94
2			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
5			Octacosanal	408	C28H56O	022725-64-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
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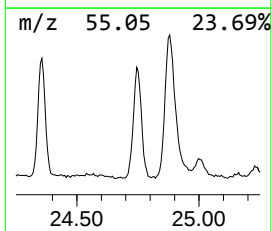
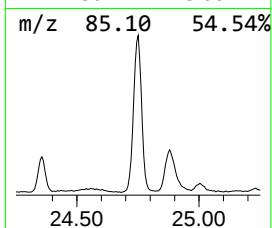
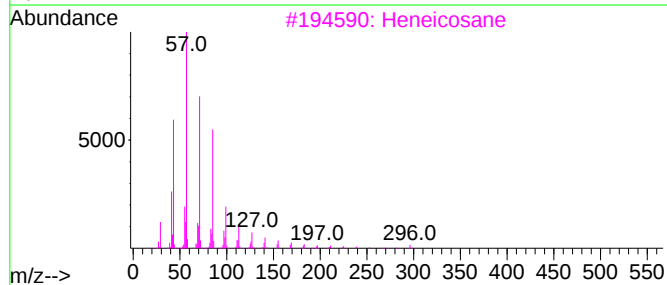
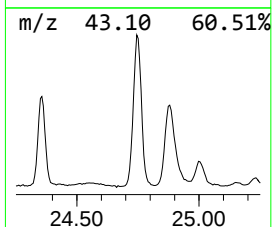
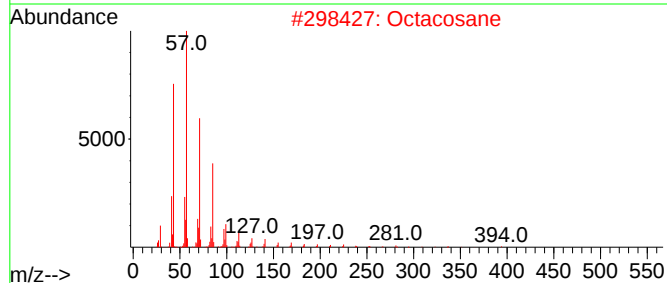
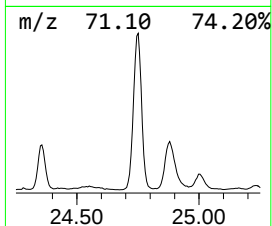
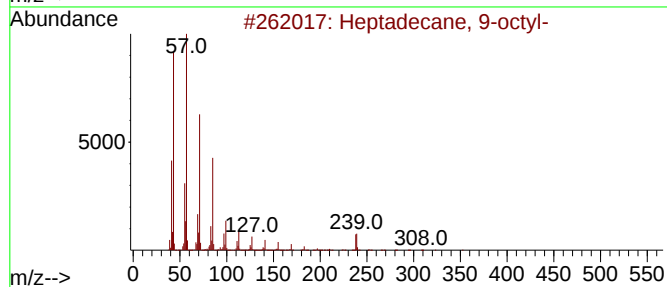
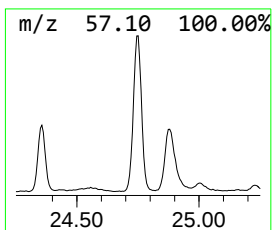
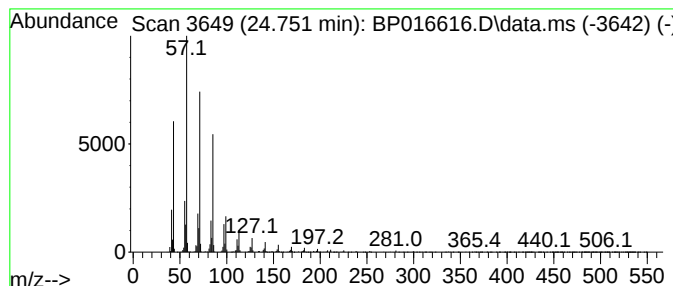
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.750	14.43 ng/ul	4035500	Perylene-d12	24.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2	Octacosane	394	C28H58	000630-02-4	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016616.D
Acq On : 16 Aug 2023 22:39
Operator : MA/JU
Sample : 03967-15
Misc :
ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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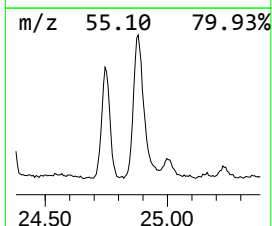
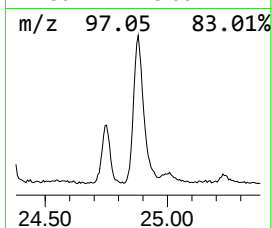
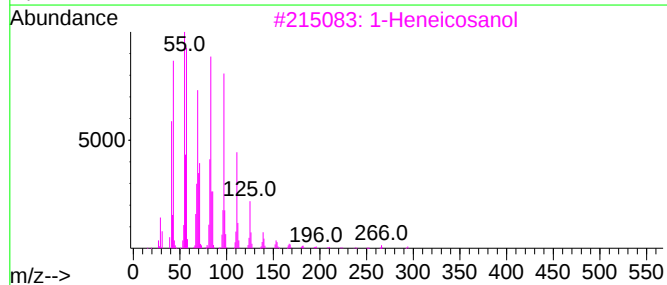
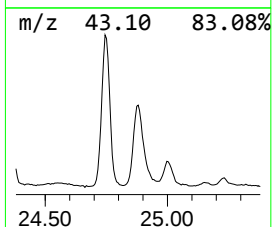
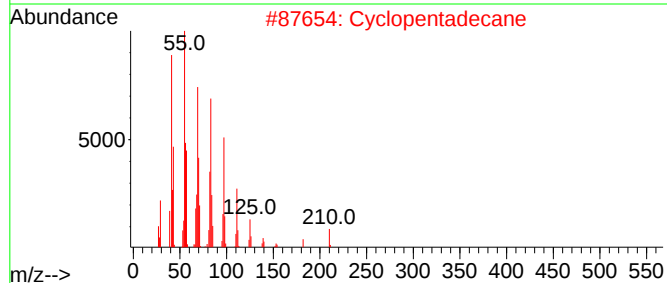
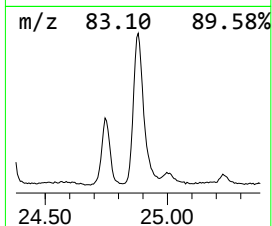
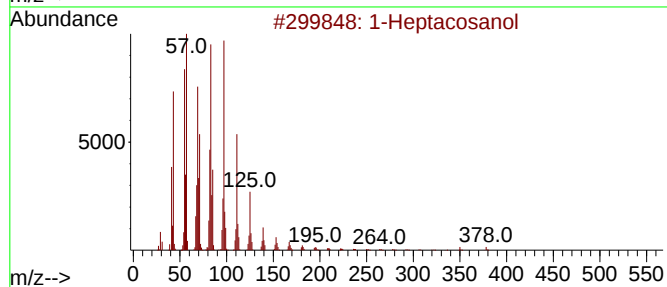
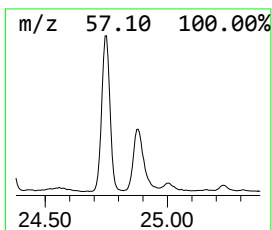
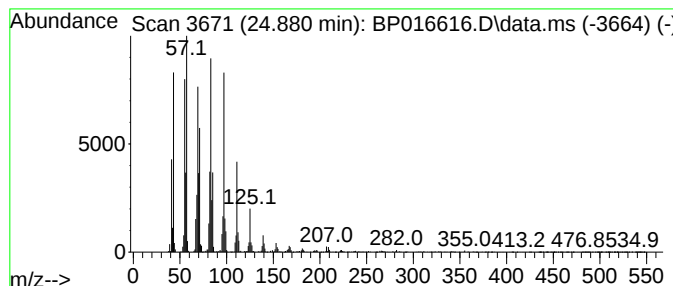
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1-Heptacosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.880	13.46 ng/ul	3765850	Perylene-d12	24.145

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heptacosanol	396	C27H56O	002004-39-9	95
2		Cyclopentadecane	210	C15H30	000295-48-7	93
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		Behenic alcohol	326	C22H46O	000661-19-8	93
5		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016616.D
Acq On : 16 Aug 2023 22:39
Operator : MA/JU
Sample : 03967-15
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A48M6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexadecanoic ...	18.310	9.4	ng/ul	2244280	4	17.469	4754090	20.0
Octadecanoic acid	19.539	2.2	ng/ul	673757	5	21.562	6078240	20.0
Ethanol, 2-(tet...	21.215	10.2	ng/ul	3104260	5	21.562	6078240	20.0
(DEL) Alkane: S...	22.145	5.6	ng/ul	1715160	5	21.562	6078240	20.0
Behenic alcohol	22.180	9.0	ng/ul	2727210	5	21.562	6078240	20.0
Oxirane, heptad...	22.956	5.2	ng/ul	1465720	6	24.145	5594820	20.0
(DEL) Alkane: S...	23.274	17.1	ng/ul	4786950	6	24.145	5594820	20.0
Hexacosanal	24.356	10.6	ng/ul	2973120	6	24.145	5594820	20.0
(DEL) Alkane: S...	24.750	14.4	ng/ul	4035500	6	24.145	5594820	20.0
1-Heptacosanol	24.880	13.5	ng/ul	3765850	6	24.145	5594820	20.0