

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011502.D
 Acq On : 19 Aug 2022 11:15
 Operator : CG/JU
 Sample : N4210-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGC48

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-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011502.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	11	15	25	rVB	30438	47525	0.34%	0.076%
2	3.487	82	85	99	rBV	167576	257078	1.86%	0.413%
3	3.693	115	120	129	rVB2	35366	59052	0.43%	0.095%
4	3.787	129	136	149	rVB	111557	167933	1.22%	0.270%
5	4.011	168	174	180	rBV	40757	59968	0.43%	0.096%
6	5.211	373	378	387	rBV	32415	60328	0.44%	0.097%
7	5.575	434	440	446	rVB2	45840	68205	0.49%	0.110%
8	6.740	632	638	650	rVB	450620	716662	5.19%	1.151%
9	6.905	659	666	679	rVB	348196	528453	3.83%	0.849%
10	7.087	691	697	707	rVB	475092	714118	5.18%	1.147%
11	7.175	707	712	714	rVV	47611	67256	0.49%	0.108%
12	7.199	714	716	722	rVB	35464	42161	0.31%	0.068%
13	7.552	766	776	786	rBV	514001	821409	5.95%	1.320%
14	8.275	893	899	909	rBV	473730	746635	5.41%	1.200%
15	8.716	967	974	980	rVV	434003	692596	5.02%	1.113%
16	8.775	980	984	993	rVB	64652	94464	0.68%	0.152%
17	9.434	1087	1096	1107	rBV	364059	608954	4.41%	0.978%
18	9.969	1176	1187	1197	rBV	507873	845039	6.12%	1.358%
19	10.340	1237	1250	1256	rBV	688565	1204011	8.73%	1.935%
20	10.393	1256	1259	1266	rVB	57050	87482	0.63%	0.141%
21	10.499	1270	1277	1285	rBV	286426	524638	3.80%	0.843%
22	11.375	1414	1426	1432	rBV	40173	79176	0.57%	0.127%
23	11.781	1489	1495	1503	rBV	84485	129467	0.94%	0.208%
24	12.022	1530	1536	1542	rVB	89870	152260	1.10%	0.245%
25	12.163	1554	1560	1567	rBV2	17198	44852	0.33%	0.072%
26	12.240	1567	1573	1583	rVB3	69998	119654	0.87%	0.192%
27	12.757	1656	1661	1665	rVB	32435	42546	0.31%	0.068%
28	12.910	1680	1687	1691	rBV	39962	55863	0.40%	0.090%
29	13.051	1706	1711	1716	rBV	113672	148083	1.07%	0.238%
30	13.387	1762	1768	1769	rBV3	35785	56038	0.41%	0.090%
31	13.545	1789	1795	1800	rBV	56770	95970	0.70%	0.154%
32	13.598	1800	1804	1807	rVV	52586	74113	0.54%	0.119%
33	13.651	1807	1813	1819	rVV	859559	1133745	8.22%	1.822%
34	13.716	1819	1824	1828	rVV3	59903	91961	0.67%	0.148%
35	13.781	1828	1835	1838	rVV5	37809	77636	0.56%	0.125%
36	13.916	1852	1858	1872	rBV	911906	1428596	10.35%	2.295%

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Stop Thrs : 0

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
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37	14.140	1889	1896	1900	rBV	110310	143564	1.04%	0.231%
38	14.228	1904	1911	1917	rBV	922890	1315673	9.53%	2.114%
39	14.287	1917	1921	1925	rBV2	33701	49083	0.36%	0.079%
40	14.457	1943	1950	1960	rBV	298050	511564	3.71%	0.822%
41	14.634	1975	1980	1984	rBV	49975	73171	0.53%	0.118%
42	14.710	1990	1993	1998	rVB	29611	38058	0.28%	0.061%
43	14.763	1998	2002	2010	rVB5	27411	47709	0.35%	0.077%
44	14.857	2011	2018	2022	rBV3	28082	57367	0.42%	0.092%
45	15.022	2040	2046	2049	rBV4	31463	60926	0.44%	0.098%
46	15.098	2055	2059	2068	rVB	91055	116801	0.85%	0.188%
47	15.234	2076	2082	2087	rBV	1128651	1612183	11.68%	2.590%
48	15.281	2087	2090	2095	rVB2	66595	96812	0.70%	0.156%
49	15.369	2099	2105	2111	rBV	336882	451669	3.27%	0.726%
50	15.510	2124	2129	2134	rVB	48144	78350	0.57%	0.126%
51	15.587	2138	2142	2151	rVB	35826	66444	0.48%	0.107%
52	15.734	2159	2167	2175	rVV5	42745	102909	0.75%	0.165%
53	15.816	2177	2181	2188	rVB5	19578	39801	0.29%	0.064%
54	15.975	2202	2208	2209	rBV2	133560	177821	1.29%	0.286%
55	16.534	2299	2303	2307	rBV3	34419	63205	0.46%	0.102%
56	16.657	2320	2324	2330	rBV2	43376	72490	0.53%	0.116%
57	16.728	2332	2336	2340	rBV2	29111	52982	0.38%	0.085%
58	16.775	2340	2344	2348	rVV	102628	132617	0.96%	0.213%
59	16.822	2348	2352	2360	rVB3	62964	112836	0.82%	0.181%
60	16.998	2376	2382	2386	rBV	1083918	1528969	11.08%	2.457%
61	17.039	2386	2389	2394	rVV	694526	901126	6.53%	1.448%
62	17.098	2394	2399	2403	rVV	1381032	1876179	13.60%	3.014%
63	17.134	2403	2405	2411	rVV	173779	205942	1.49%	0.331%
64	17.198	2411	2416	2421	rVB4	29597	56717	0.41%	0.091%
65	17.416	2449	2453	2463	rVV	86164	167229	1.21%	0.269%
66	17.522	2467	2471	2476	rBV2	107349	134482	0.97%	0.216%
67	17.698	2497	2501	2505	rBV	130164	148967	1.08%	0.239%
68	17.804	2515	2519	2523	rBV4	33941	52991	0.38%	0.085%
69	17.869	2525	2530	2535	rVV3	218335	413887	3.00%	0.665%
70	17.922	2535	2539	2547	rVV	620097	909932	6.59%	1.462%
71	17.992	2547	2551	2558	rVV2	117272	263946	1.91%	0.424%
72	18.063	2559	2563	2568	rVV2	185041	354357	2.57%	0.569%
73	18.104	2568	2570	2576	rVB	100076	119827	0.87%	0.193%
74	18.204	2584	2587	2593	rVB	113004	138265	1.00%	0.222%
75	18.375	2612	2616	2619	rBV	89960	110074	0.80%	0.177%
76	18.416	2619	2623	2626	rVB	81641	101927	0.74%	0.164%

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77	18.780	2681	2685	2689	rBV2	93398	154031	1.12%	0.247%
78	18.822	2689	2692	2694	rBV2	128414	153577	1.11%	0.247%
79	18.845	2694	2696	2699	rVB2	142722	132331	0.96%	0.213%
80	18.916	2705	2708	2715	rVB2	76796	147375	1.07%	0.237%
81	18.980	2716	2719	2722	rBV	91666	111467	0.81%	0.179%
82	19.033	2723	2728	2733	rBV	1567236	1971894	14.29%	3.168%
83	19.175	2748	2752	2757	rBV2	216990	311256	2.26%	0.500%
84	19.363	2779	2784	2787	rBV	1859727	2643214	19.16%	4.247%
85	19.392	2787	2789	2798	rVB	1384174	1488908	10.79%	2.392%
86	19.916	2875	2878	2882	rVB	102070	105024	0.76%	0.169%
87	19.980	2886	2889	2892	rVB	125121	150471	1.09%	0.242%
88	20.869	3035	3040	3044	rBV4	414405	770322	5.58%	1.238%
89	21.063	3068	3073	3078	rBV	13061130	13798806	100.00%	22.171%
90	21.127	3078	3084	3088	rVV2	1666041	3027245	21.94%	4.864%
91	21.169	3088	3091	3100	rVB	710016	949363	6.88%	1.525%
92	21.733	3185	3187	3192	rVB2	242501	273841	1.98%	0.440%
93	22.457	3306	3310	3319	rVB2	494976	817627	5.93%	1.314%
94	22.739	3352	3358	3362	rBV2	719062	1614624	11.70%	2.594%
95	22.816	3367	3371	3381	rVB2	789921	1696156	12.29%	2.725%
96	22.998	3399	3402	3418	rVB	598781	1401742	10.16%	2.252%
97	23.280	3445	3450	3454	rBV2	962481	1646468	11.93%	2.645%
98	23.327	3454	3458	3467	rVB	670717	1344740	9.75%	2.161%
99	23.427	3470	3475	3481	rVB2	781359	1358828	9.85%	2.183%
100	24.345	3626	3631	3641	rVB	486463	1066425	7.73%	1.713%

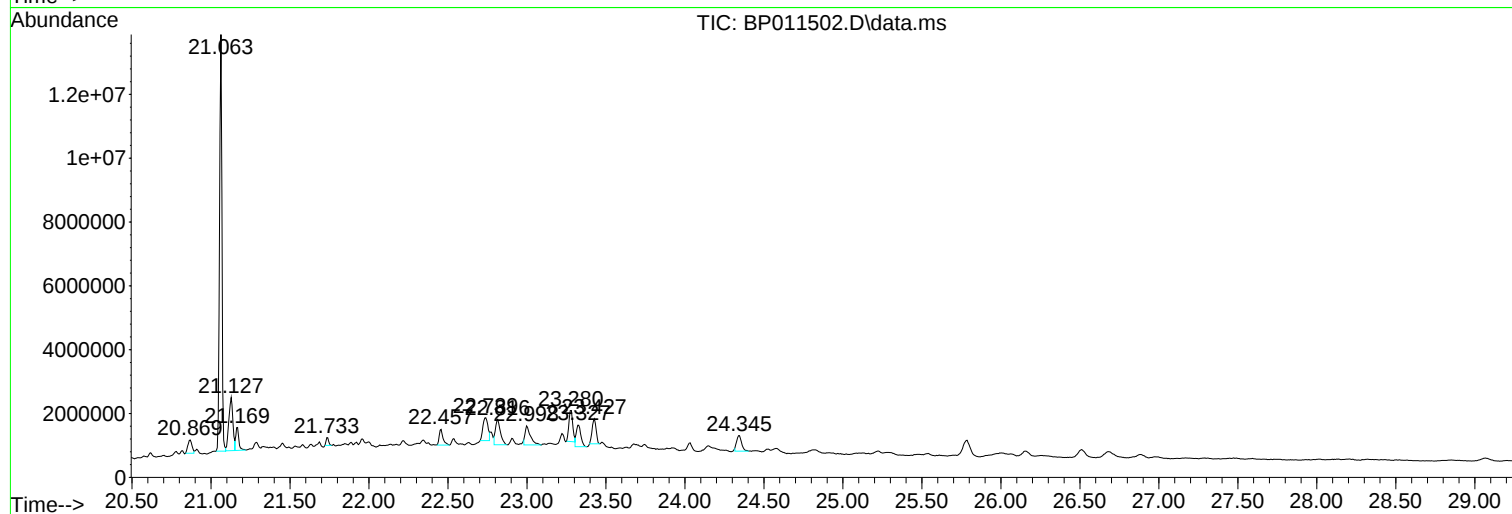
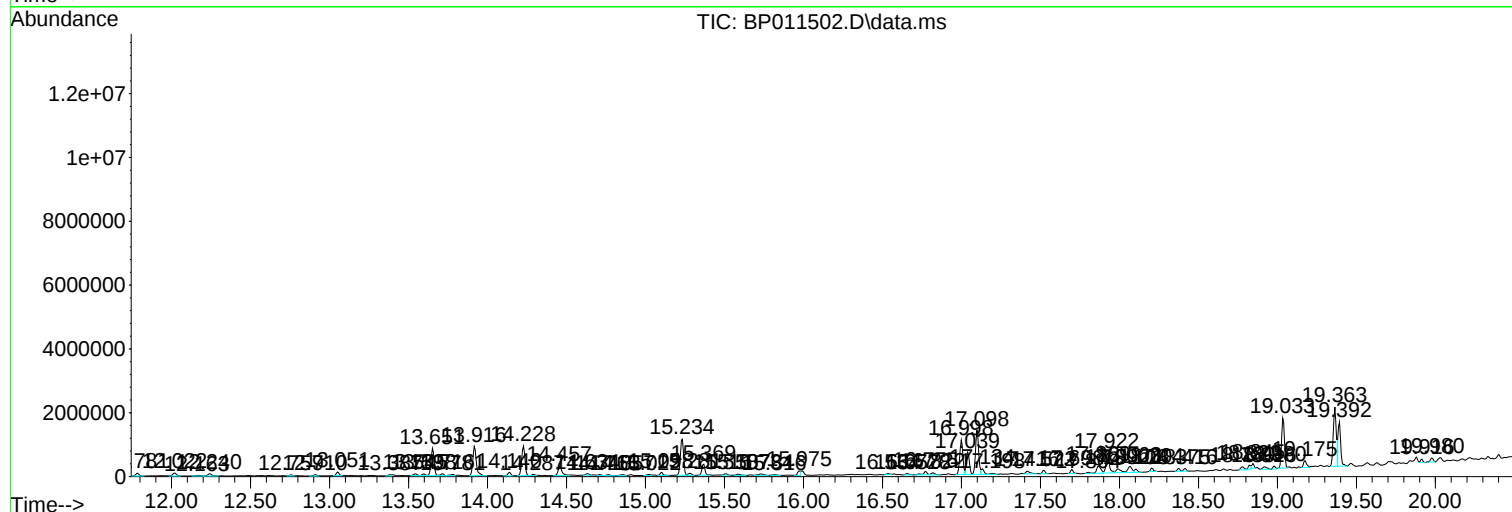
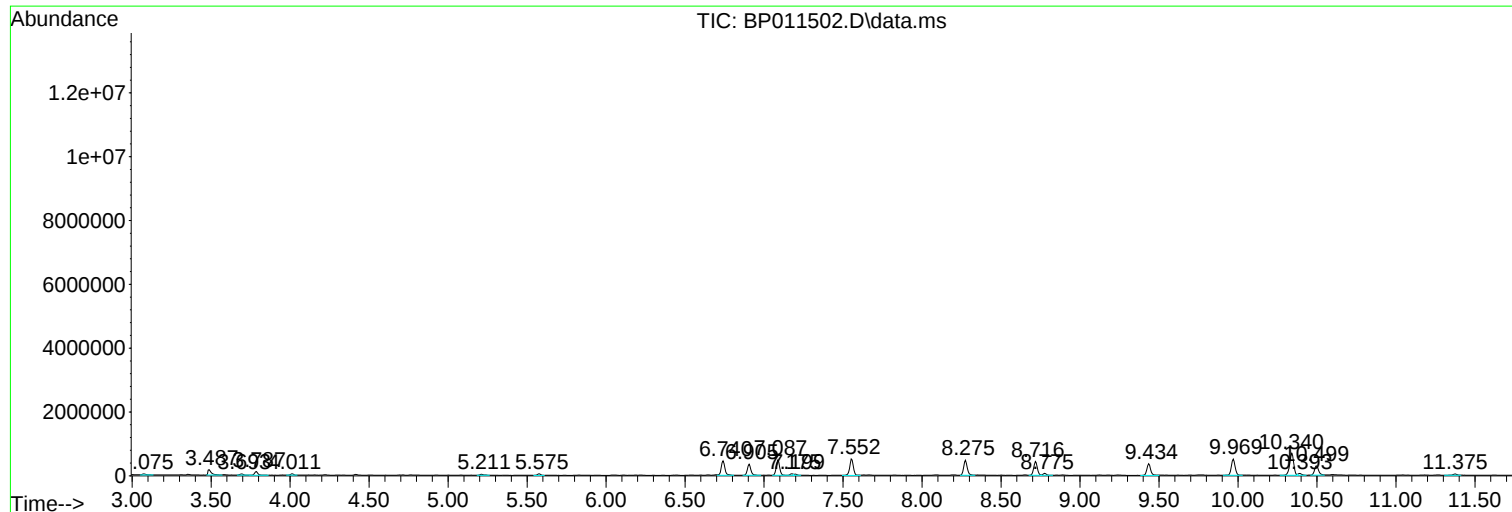
Sum of corrected areas: 62238514

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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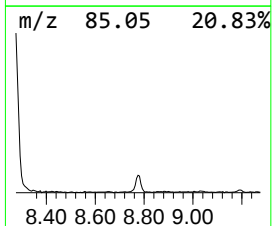
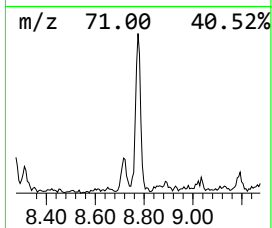
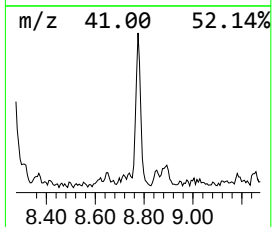
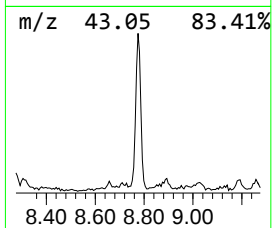
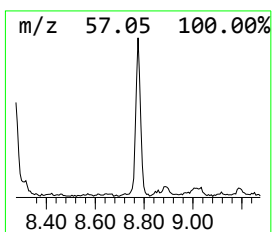
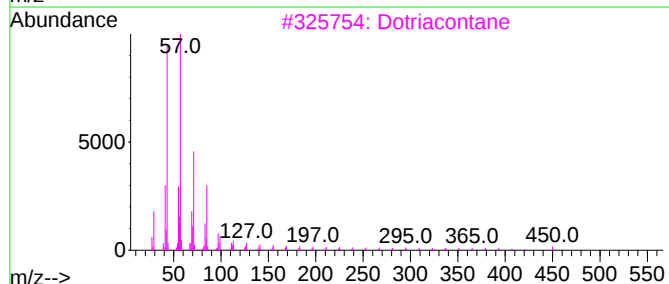
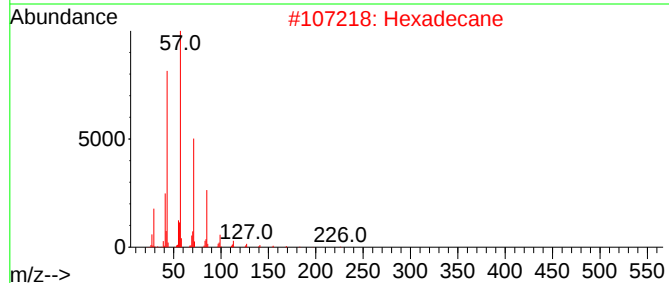
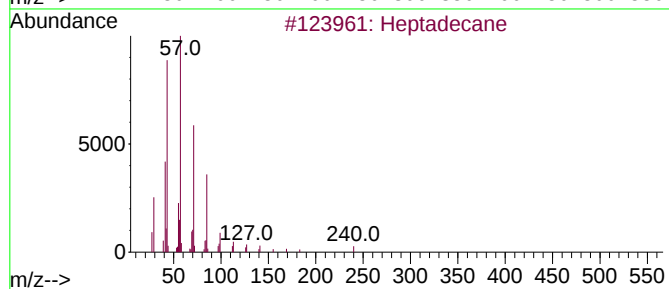
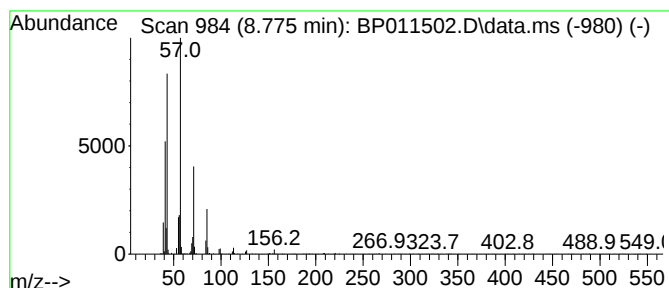
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.775	2.30 ng/ul	94464	1,4-Dichlorobenzene-d4	7.552

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	91
2		Hexadecane	226	C16H34	000544-76-3	90
3		Dotriacontane	451	C32H66	000544-85-4	83
4		Tetradecane	198	C14H30	000629-59-4	83
5		Tridecane	184	C13H28	000629-50-5	72



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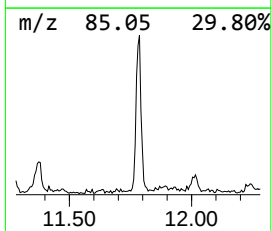
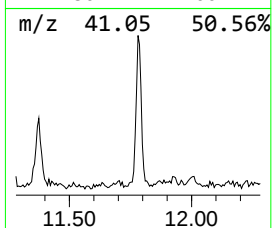
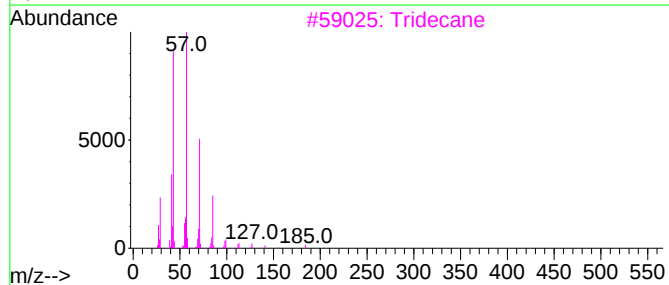
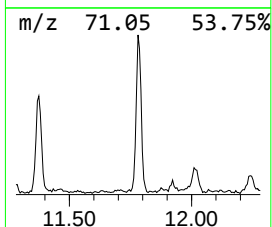
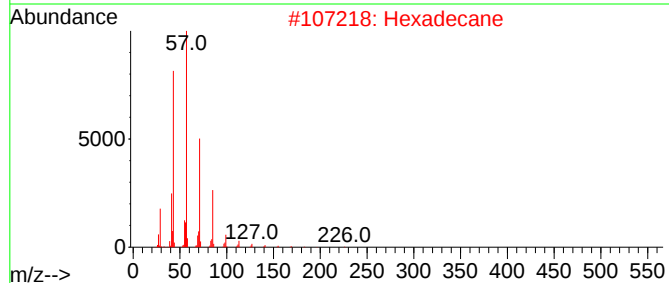
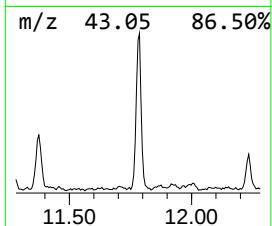
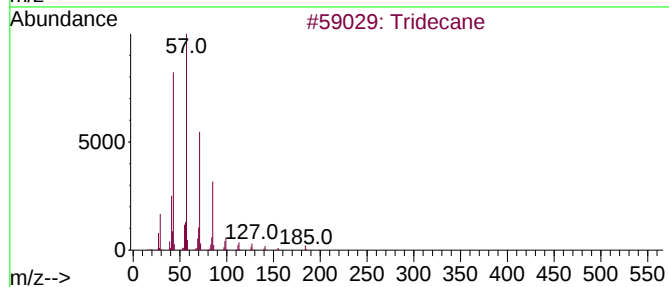
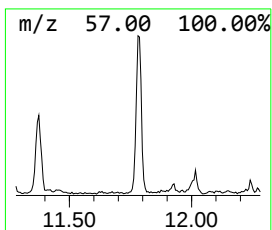
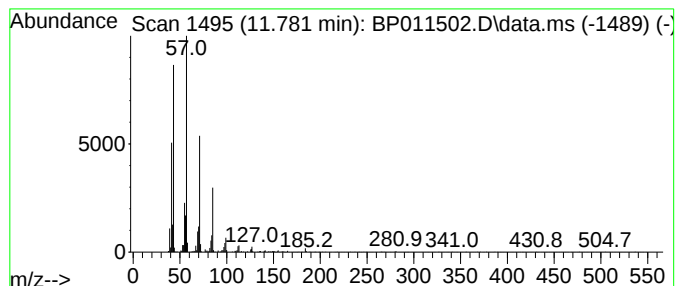
Instrument :
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.781	2.15 ng/ul	129467	Naphthalene-d8	10.340	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	93
2	Hexadecane	226	C16H34	000544-76-3	91
3	Tridecane	184	C13H28	000629-50-5	90
4	Tetradecane	198	C14H30	000629-59-4	90
5	Nonadecane	268	C19H40	000629-92-5	90



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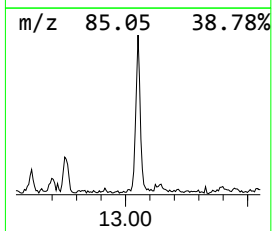
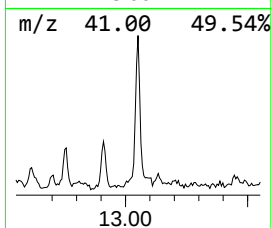
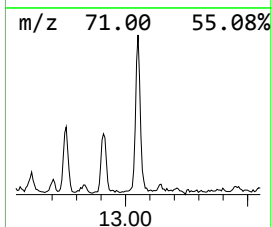
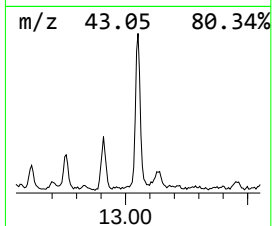
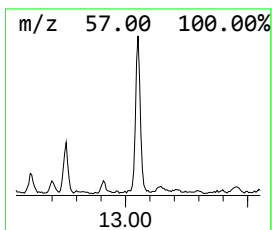
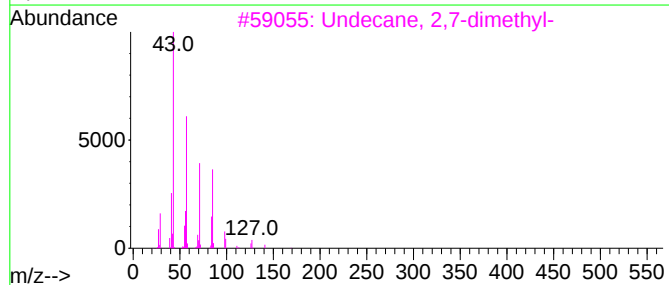
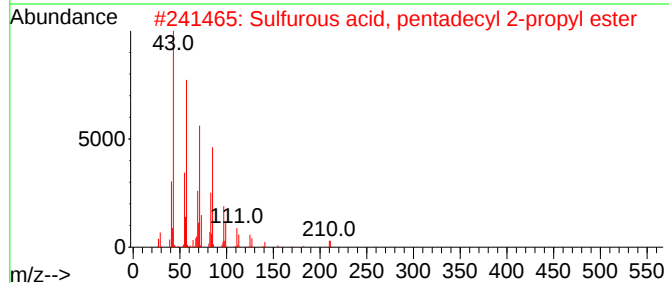
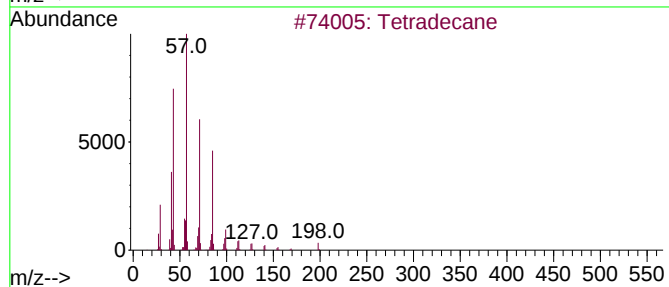
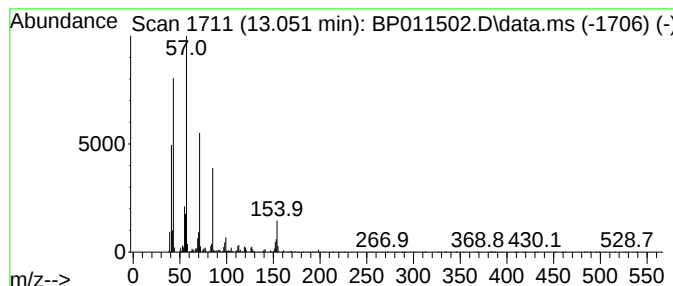
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.051	2.25 ng/ul	148083	Acenaphthene-d10	14.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	90
2			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	59
3			Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	53
4			Undecane, 5-methyl-	170	C12H26	001632-70-8	53
5			Hexadecane	226	C16H34	000544-76-3	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
Operator : CG/JU
Sample : N4210-05
Misc :
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Instrument :
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ClientSampleId :
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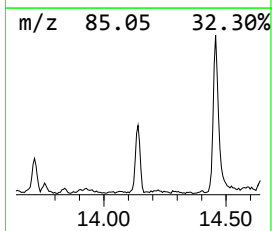
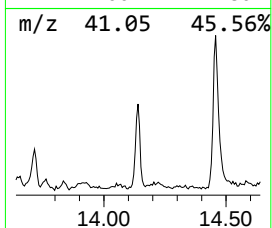
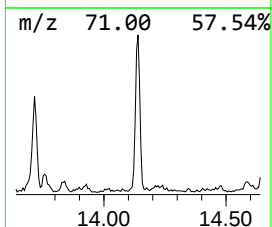
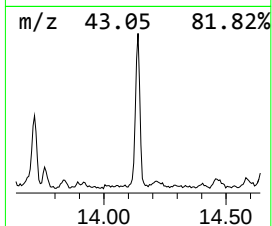
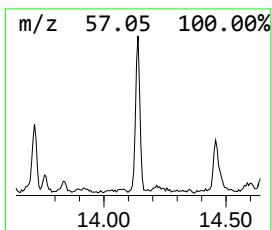
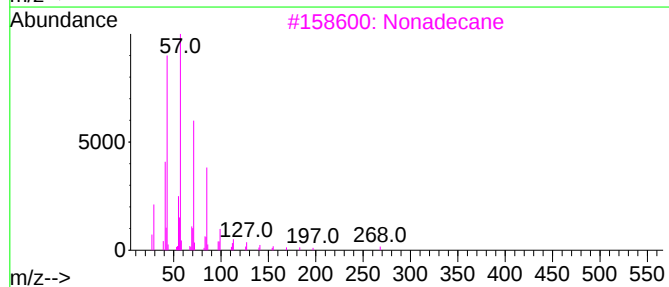
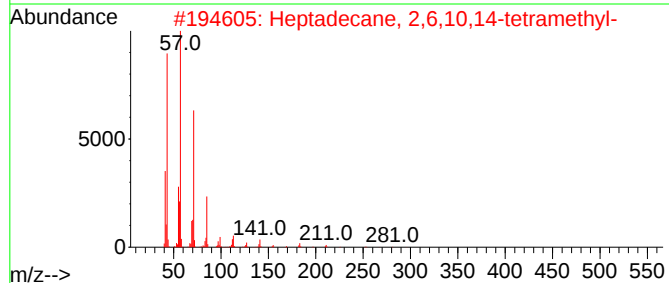
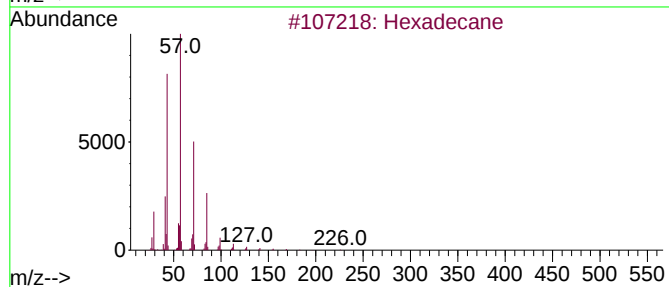
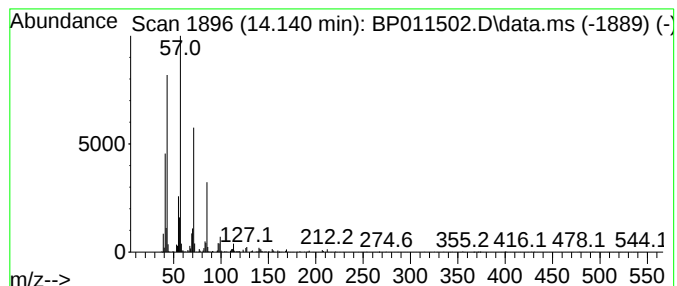
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.140	2.18 ng/ul	143564	Acenaphthene-d10	14.228

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	90
3			Nonadecane	268	C19H40	000629-92-5	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Pentadecane	212	C15H32	000629-62-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
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ClientSampleId :
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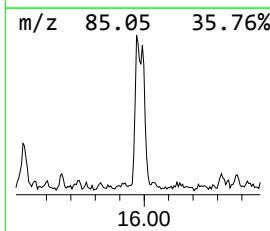
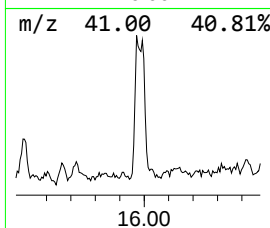
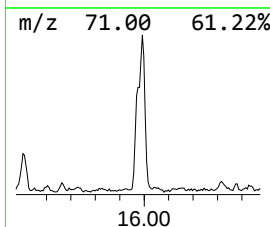
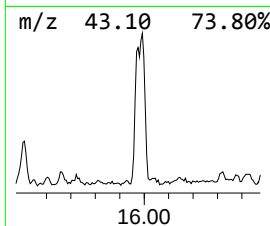
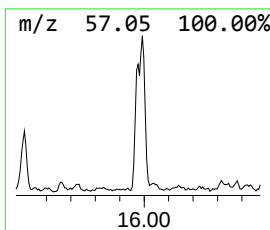
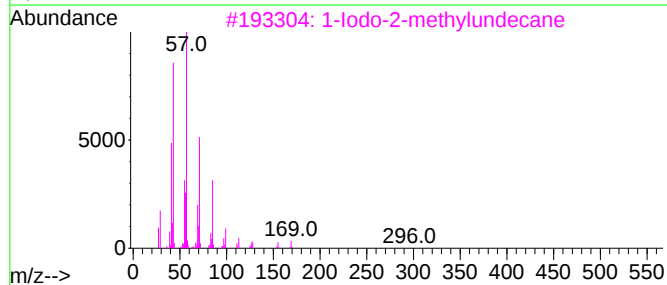
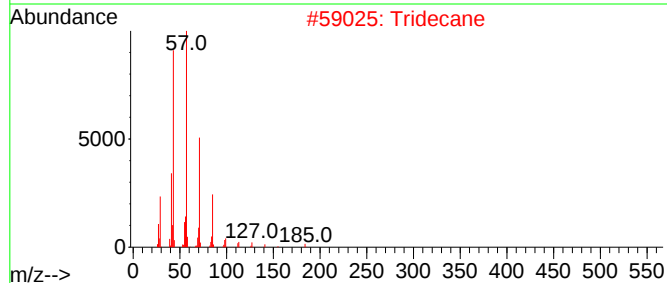
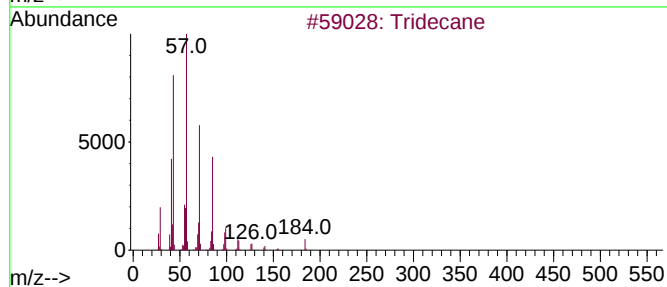
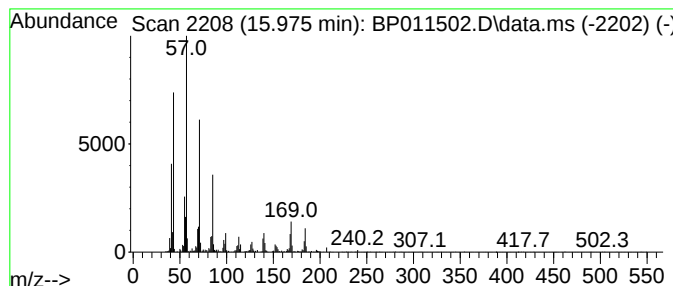
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.975	2.33 ng/ul	177821	Phenanthrene-d10	16.998

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane	184	C13H28	000629-50-5	89
2		Tridecane	184	C13H28	000629-50-5	68
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	64
4		Heptadecane	240	C17H36	000629-78-7	62
5		Pentadecane	212	C15H32	000629-62-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
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Sample : N4210-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

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ClientSampleId :
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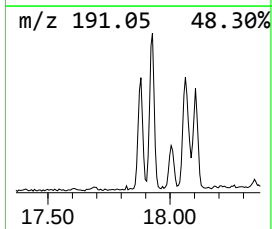
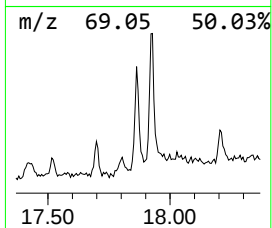
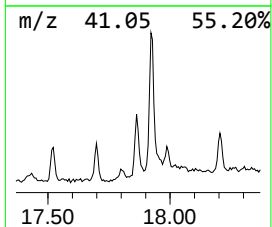
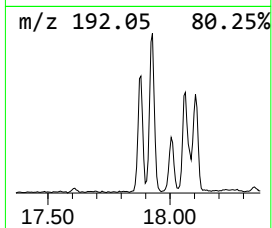
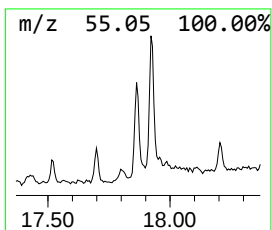
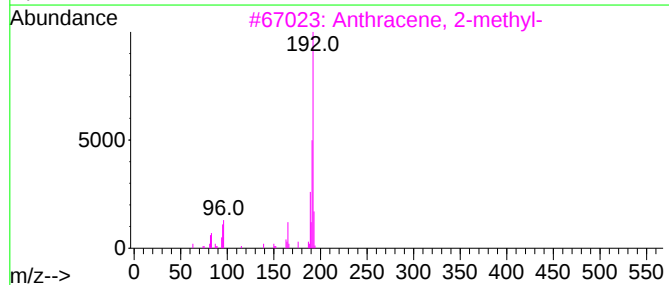
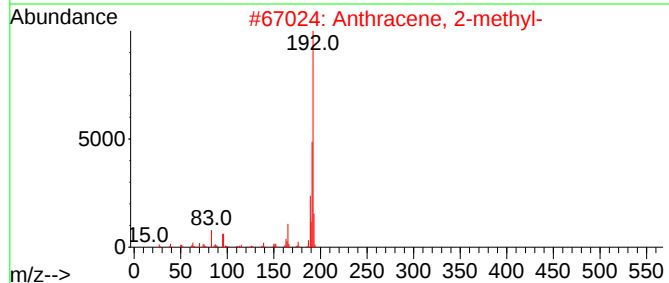
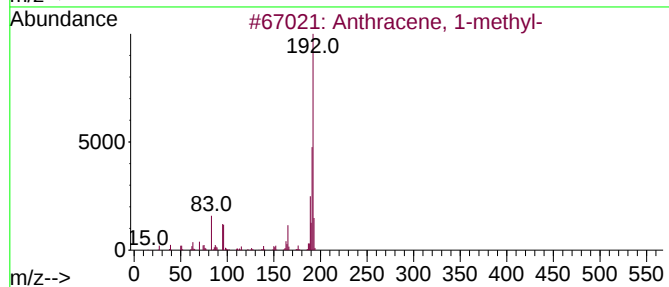
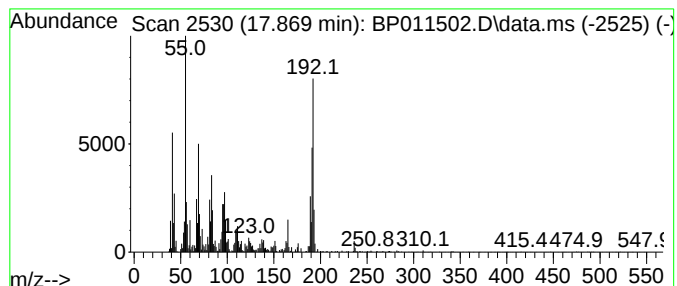
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Anthracene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.869	5.41 ng/ul	413887	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	91
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
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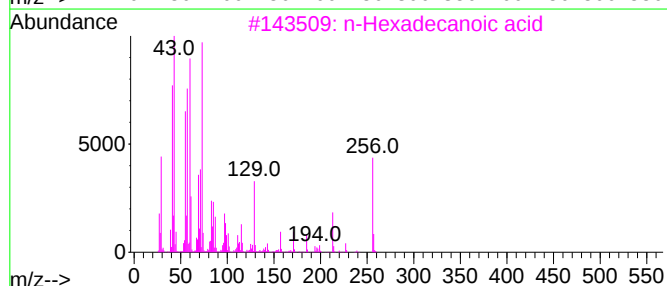
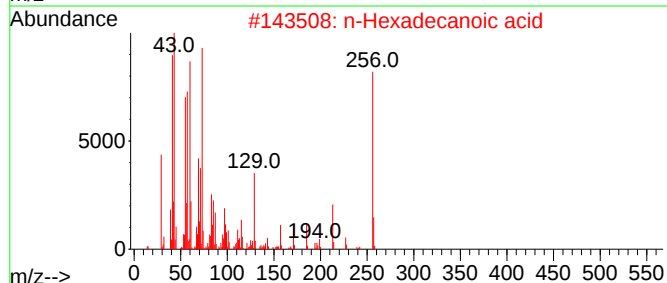
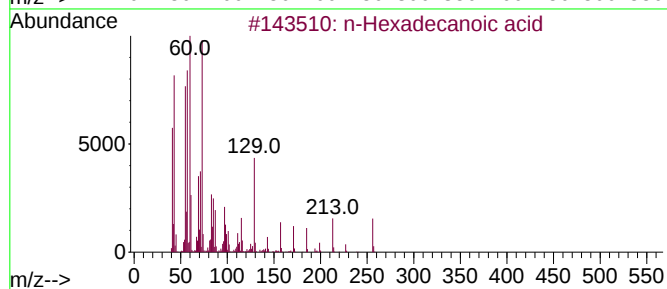
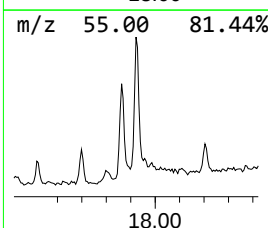
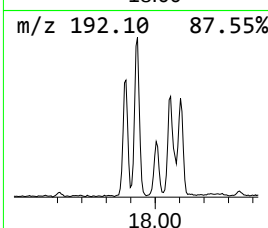
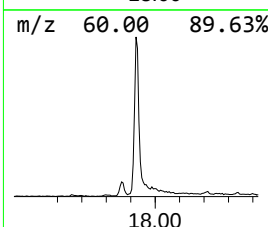
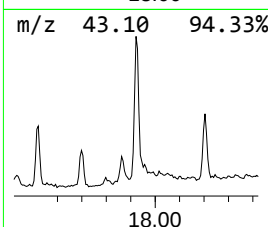
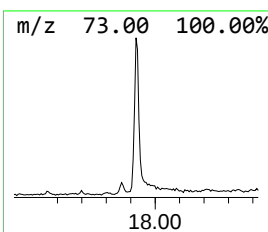
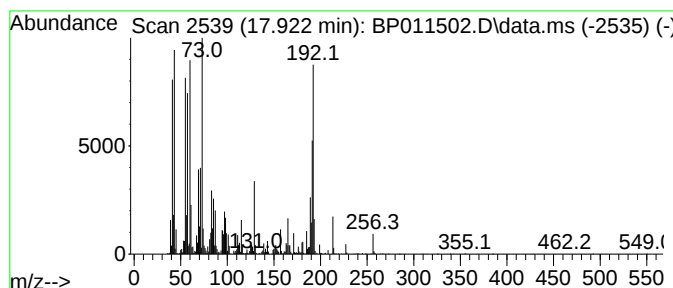
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.922	11.90 ng/ul	909932	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
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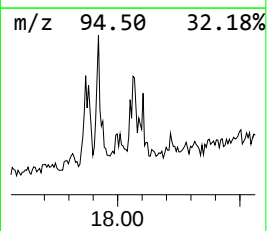
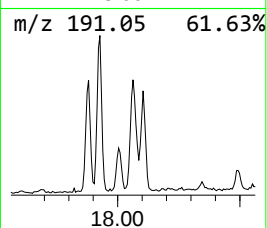
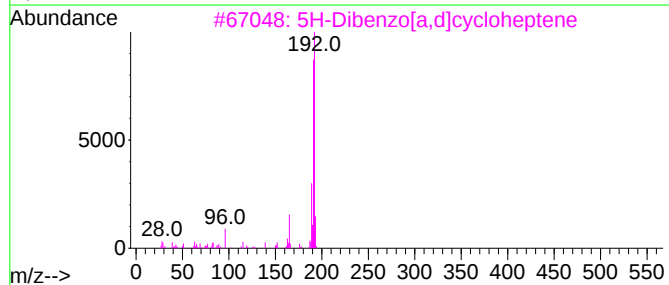
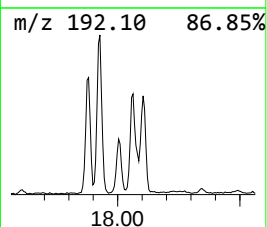
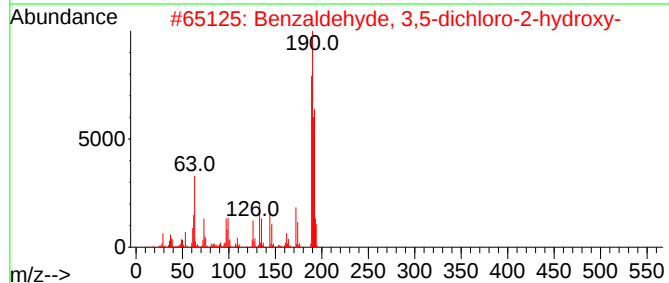
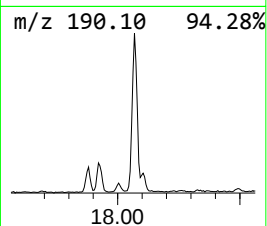
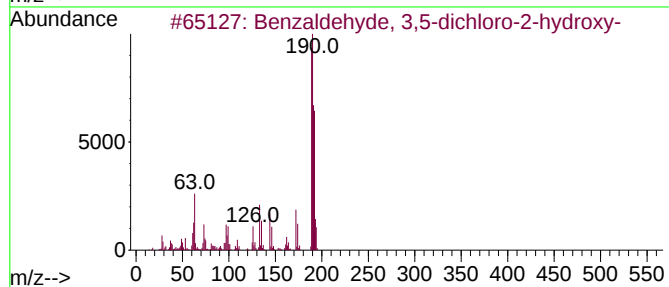
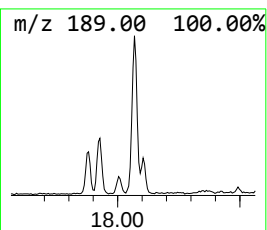
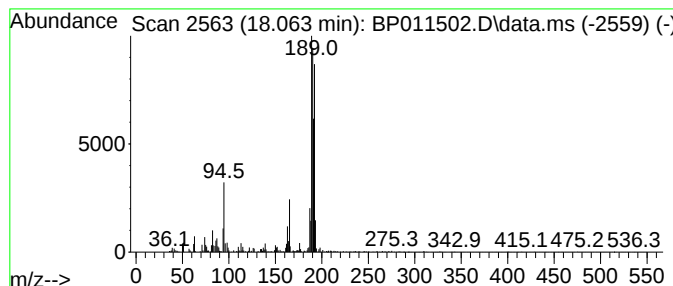
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzaldehyde, 3,5-dichloro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.063	4.64 ng/ul	354357	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
3			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	55
4			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	46
5			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	45



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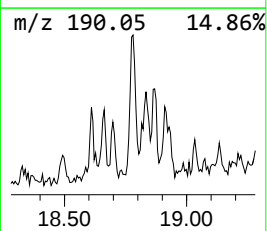
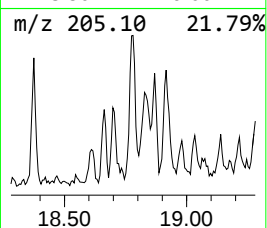
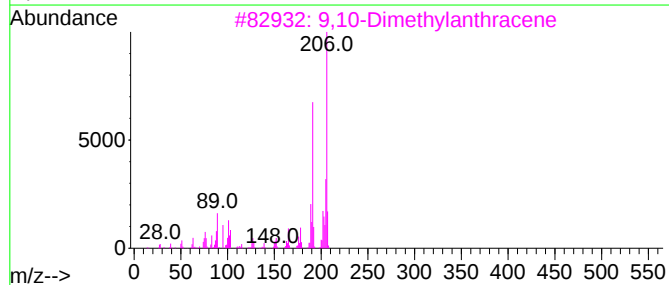
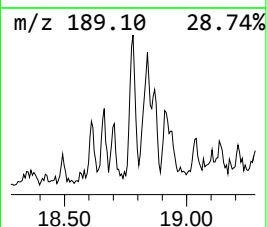
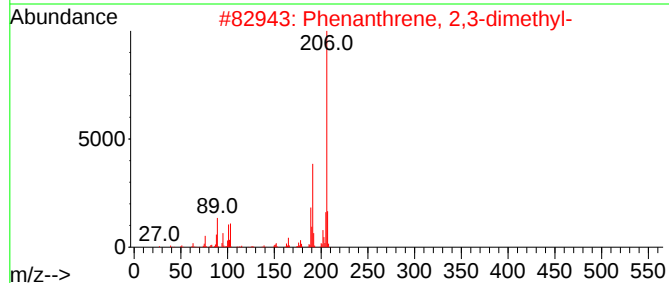
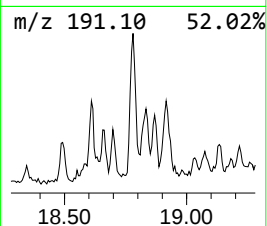
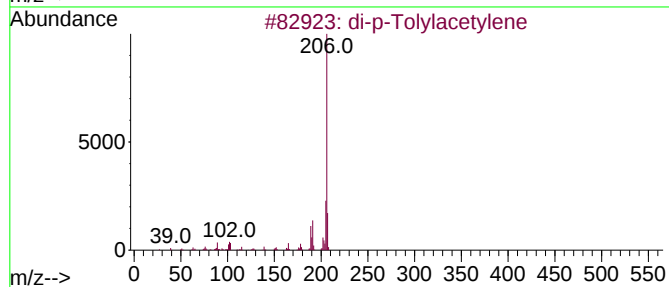
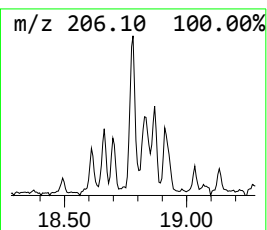
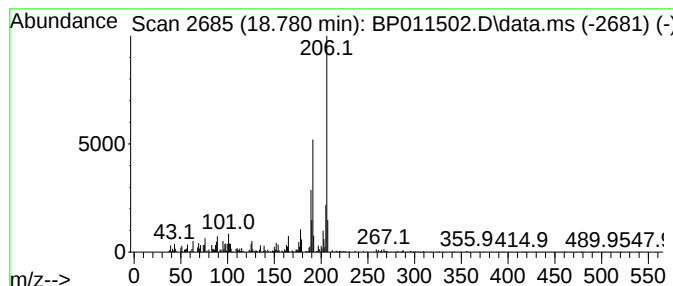
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 di-p-Tolylacetylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.781	2.01 ng/ul	154031	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
3			9,10-Dimethylantracene	206	C16H14	000781-43-1	90
4			9,10-Dimethylantracene	206	C16H14	000781-43-1	87
5			9,10-Dimethylantracene	206	C16H14	000781-43-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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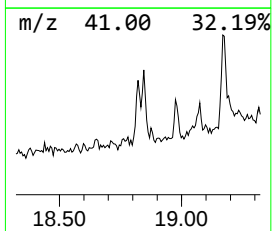
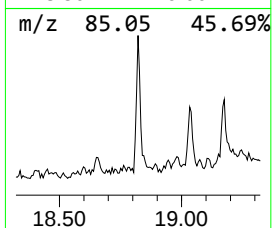
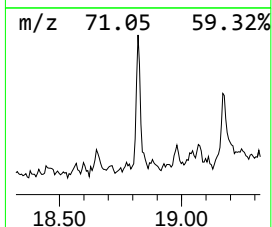
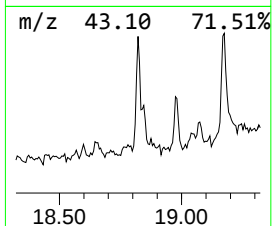
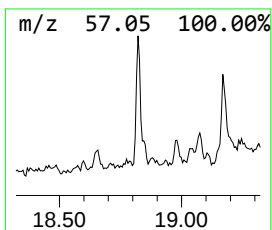
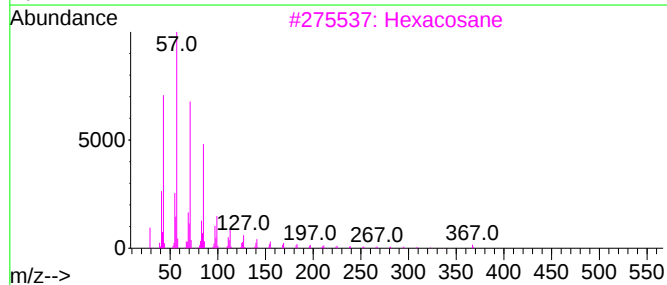
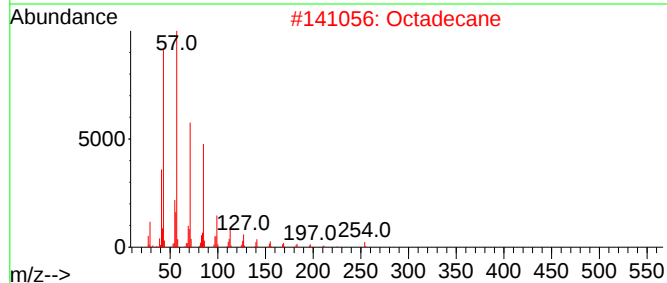
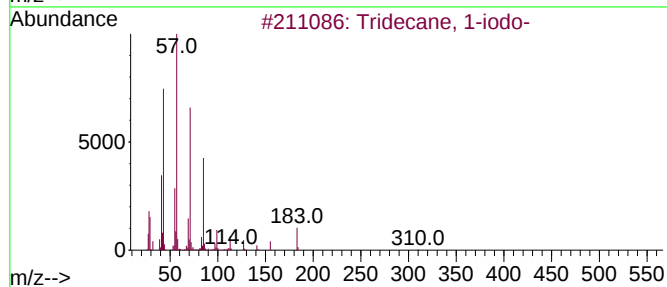
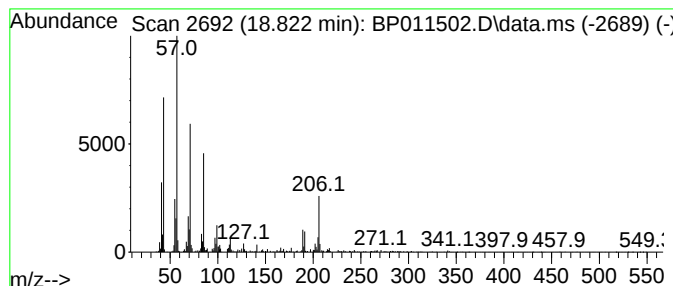
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Tridecane, 1-iodo- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.822	2.01 ng/ul	153577	Phenanthrene-d10	16.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
2			Octadecane	254	C18H38	000593-45-3	58
3			Hexacosane	366	C26H54	000630-01-3	52
4			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	52
5			2-Bromo dodecane	248	C12H25Br	013187-99-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
Operator : CG/JU
Sample : N4210-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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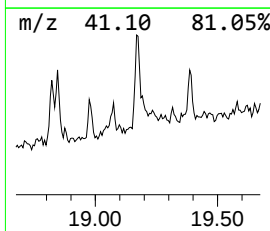
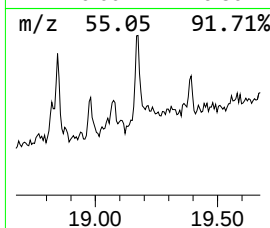
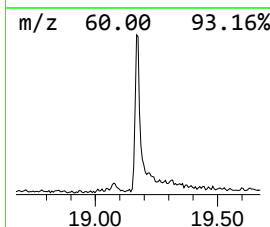
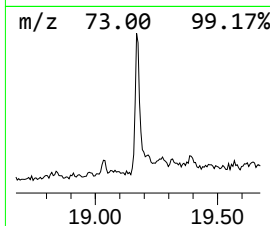
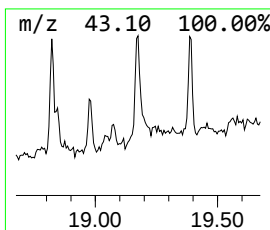
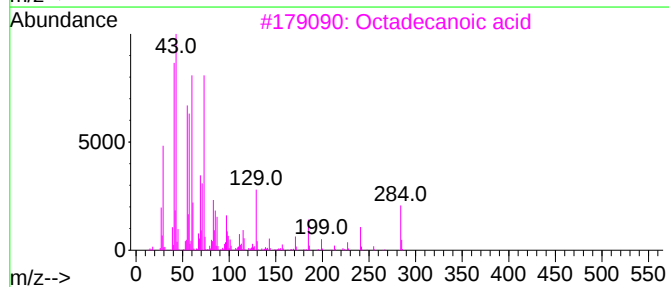
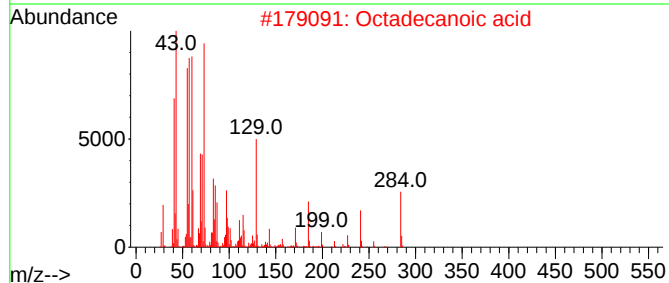
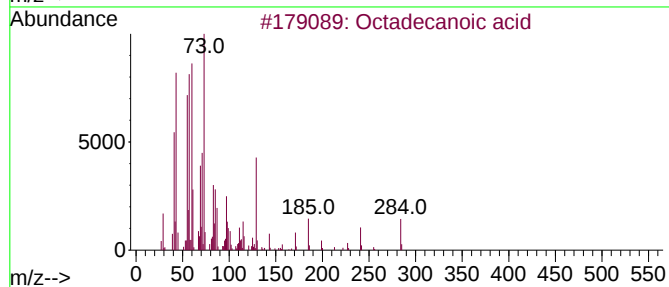
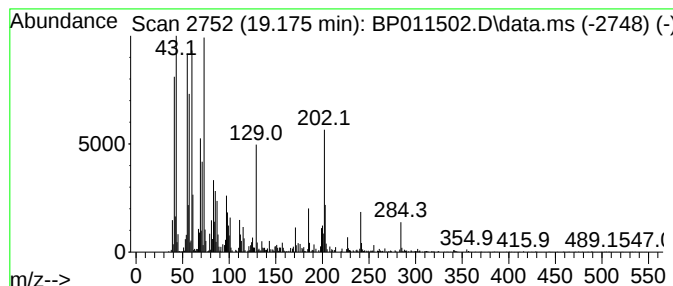
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.175	2.06 ng/ul	311256	Chrysene-d12	21.127

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	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
Operator : CG/JU
Sample : N4210-05
Misc :
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Instrument :
BNA_P
ClientSampleId :
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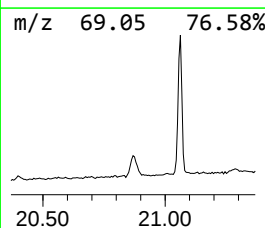
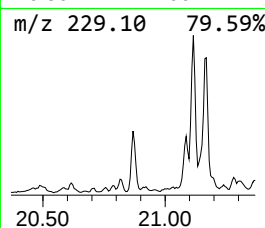
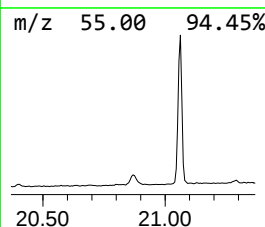
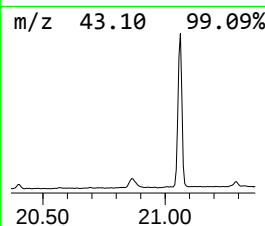
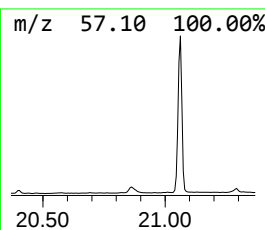
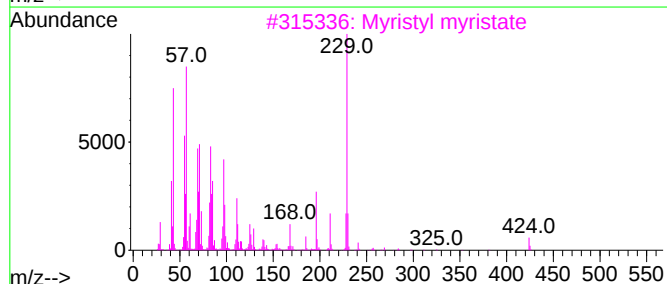
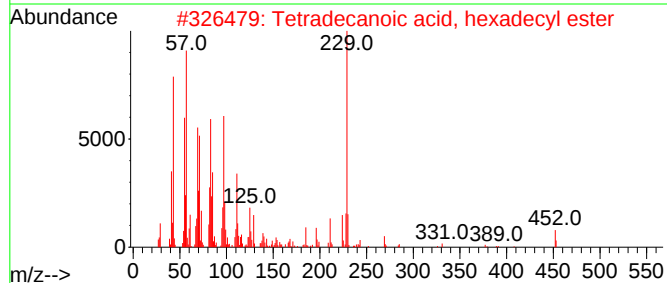
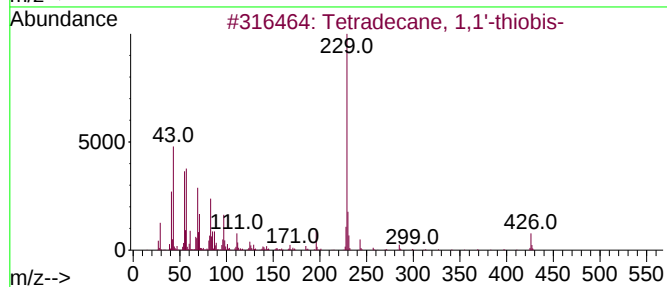
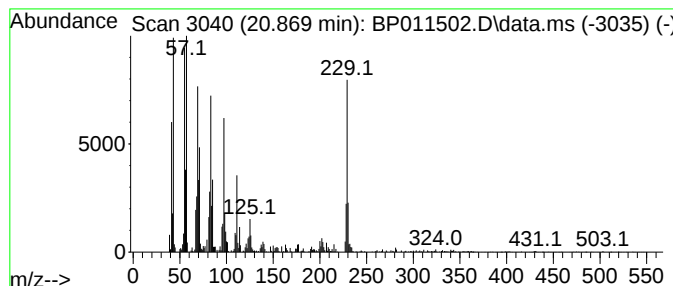
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Tetradecane, 1,1'-thiobis- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.869	5.09 ng/ul	770322	Chrysene-d12	21.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 1,1'-thiobis-	426	C28H58S	035599-83-8	64
2			Tetradecanoic acid, hexadecyl ester	452	C30H60O2	002599-01-1	52
3			Myristyl myristate	424	C28H56O2	003234-85-3	43
4			Tetradecanoic acid, hexadecyl ester	452	C30H60O2	002599-01-1	40
5			Hexadecanoic acid, 2-hydroxy-, m...	286	C17H34O3	016742-51-1	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
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Sample : N4210-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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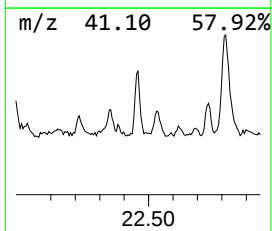
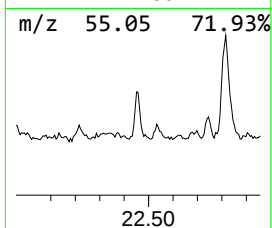
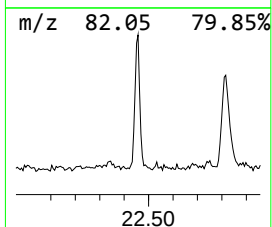
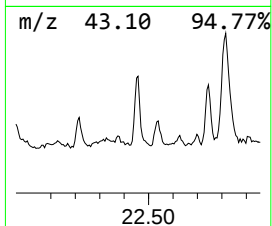
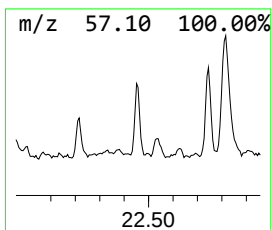
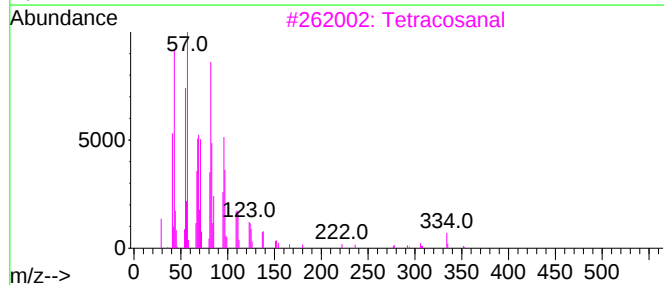
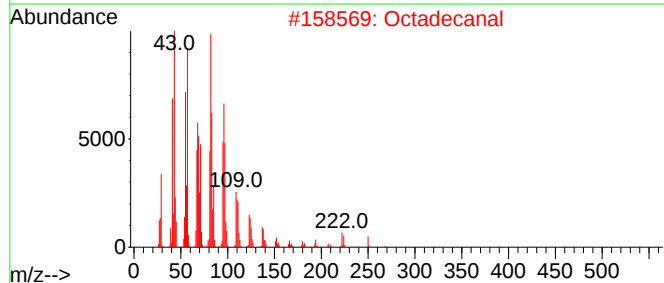
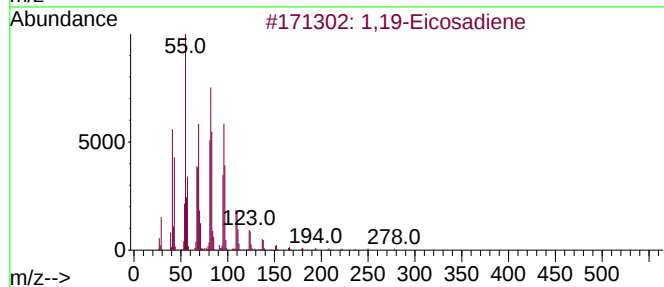
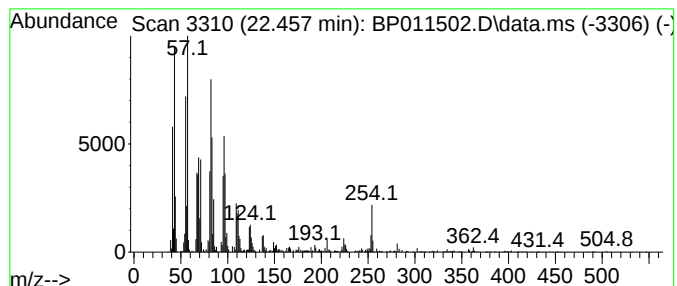
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1,19-Eicosadiene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.457	12.03 ng/ul	817627	Perylene-d12	23.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,19-Eicosadiene	278	C20H38	014811-95-1	94
2			Octadecanal	268	C18H36O	000638-66-4	90
3			Tetracosanal	352	C24H48O	057866-08-7	89
4			Tetracosanal	352	C24H48O	057866-08-7	87
5			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
Operator : CG/JU
Sample : N4210-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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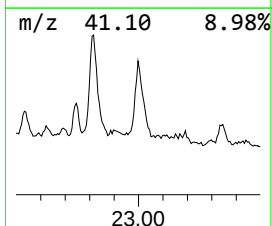
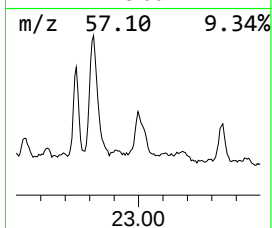
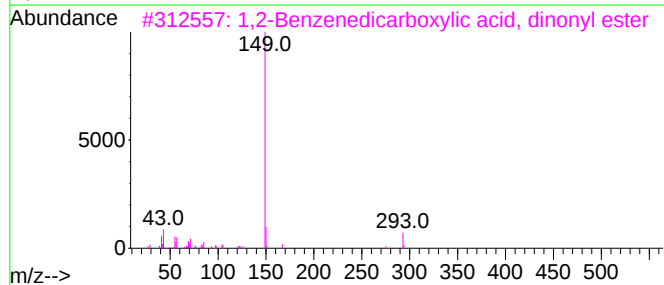
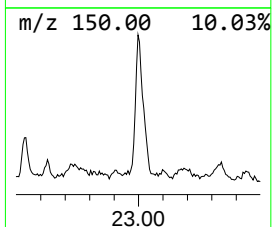
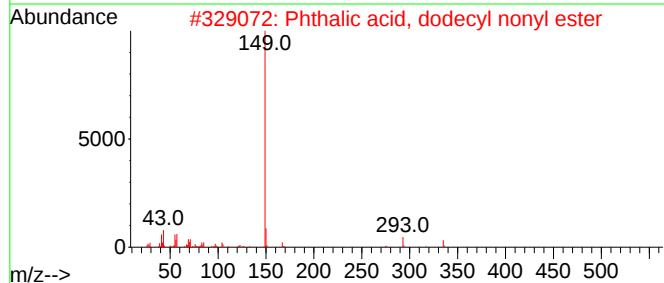
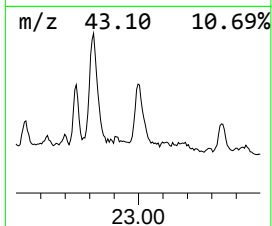
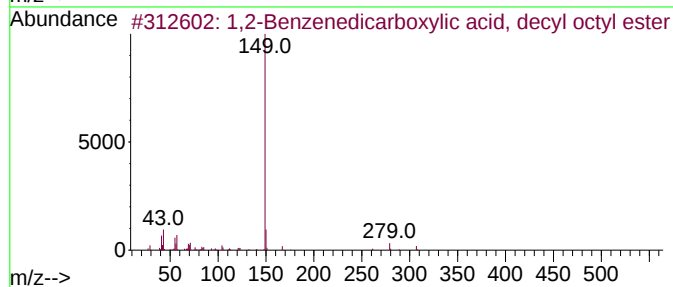
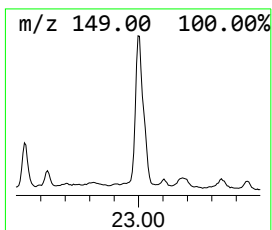
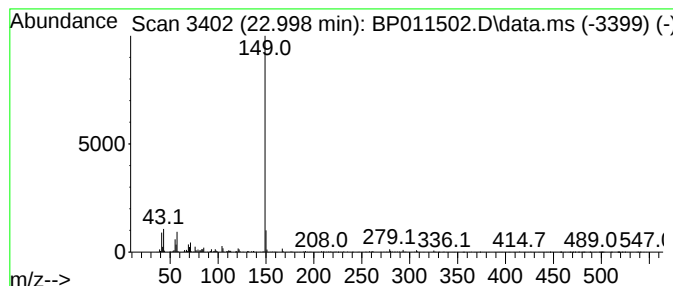
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,2-Benzenedicarboxylic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.998	20.63 ng/ul	1401740	Perylene-d12	23.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	87
2			Phthalic acid, dodecyl nonyl ester	460	C29H48O4	1000308-92-6	80
3			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	80
4			Di-n-octyl phthalate	390	C24H38O4	000117-84-0	80
5			1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011502.D
Acq On : 19 Aug 2022 11:15
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Sample : N4210-05
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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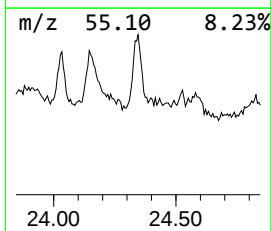
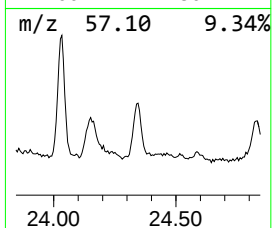
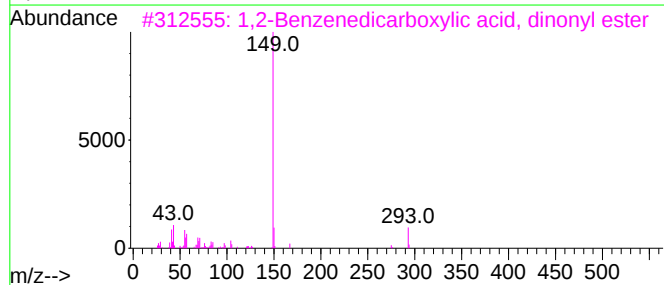
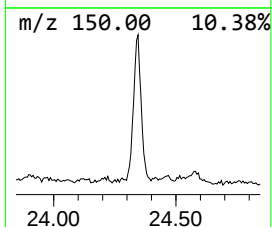
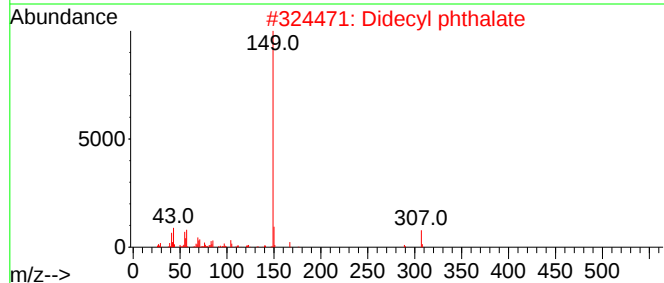
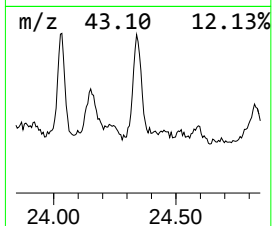
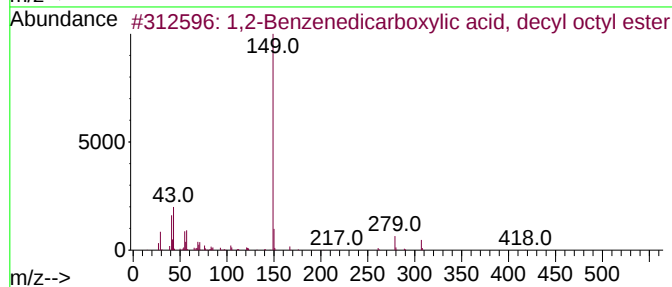
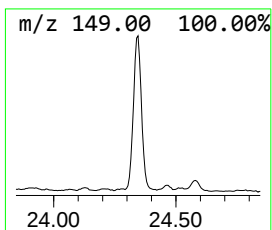
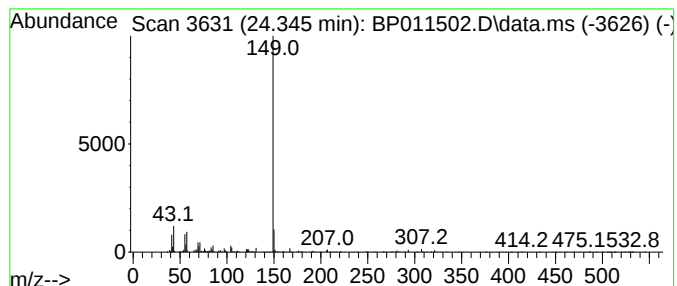
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Didecyl phthalate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.345	15.70 ng/ul	1066430	Perylene-d12	23.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	83
2			Didecyl phthalate	446	C28H46O4	000084-77-5	83
3			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	80
4			Di-n-octyl phthalate	390	C24H38O4	000117-84-0	80
5			1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011502.D
 Acq On : 19 Aug 2022 11:15
 Operator : CG/JU
 Sample : N4210-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.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
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(DEL) Alkane: S...	11.781	2.1	ng/ul	129467	2	10.340	1204010	20.0
(DEL) Alkane: S...	13.051	2.3	ng/ul	148083	3	14.228	1315670	20.0
(DEL) Alkane: S...	14.140	2.2	ng/ul	143564	3	14.228	1315670	20.0
(DEL) Alkane: S...	15.975	2.3	ng/ul	177821	4	16.998	1528970	20.0
Anthracene, 1-m...	17.869	5.4	ng/ul	413887	4	16.998	1528970	20.0
n-Hexadecanoic ...	17.922	11.9	ng/ul	909932	4	16.998	1528970	20.0
Benzaldehyde, 3...	18.063	4.6	ng/ul	354357	4	16.998	1528970	20.0
di-p-Tolylacety...	18.781	2.0	ng/ul	154031	4	16.998	1528970	20.0
Tridecane, 1-iodo-	18.822	2.0	ng/ul	153577	4	16.998	1528970	20.0
Octadecanoic acid	19.175	2.1	ng/ul	311256	5	21.127	3027250	20.0
Tetradecane, 1,...	20.869	5.1	ng/ul	770322	5	21.127	3027250	20.0
1,19-Eicosadiene	22.457	12.0	ng/ul	817627	6	23.427	1358830	20.0
1,2-Benzenedica...	22.998	20.6	ng/ul	1401740	6	23.427	1358830	20.0
Didecyl phthalate	24.345	15.7	ng/ul	1066430	6	23.427	1358830	20.0