

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
 Data File : BP017901.D
 Acq On : 02 Nov 2023 23:25
 Operator : MA/JU
 Sample : 05060-02
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKG0

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-BP101823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017901.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.722	103	108	117	rVV6	14716	38166	1.45%	0.105%
2	3.811	119	123	130	rVV	55288	99098	3.78%	0.272%
3	3.881	131	135	145	rVB3	22811	49260	1.88%	0.135%
4	3.993	152	154	166	rVB7	9878	25532	0.97%	0.070%
5	4.581	250	254	261	rVB7	7602	13285	0.51%	0.036%
6	4.922	308	312	319	rVV	20754	39254	1.50%	0.108%
7	5.069	332	337	340	rBV7	5920	10962	0.42%	0.030%
8	5.840	462	468	472	rBV4	7795	14020	0.53%	0.038%
9	6.475	572	576	580	rBV	14831	22949	0.87%	0.063%
10	7.046	667	673	684	rBV	235988	477269	18.19%	1.308%
11	7.228	696	704	711	rBV	194685	327988	12.50%	0.899%
12	7.399	726	733	744	rBV	261003	450163	17.15%	1.233%
13	7.505	746	751	753	rBV6	7118	10294	0.39%	0.028%
14	7.875	805	814	824	rBV	606669	1017279	38.76%	2.787%
15	8.416	900	906	912	rVB2	23002	38775	1.48%	0.106%
16	8.605	927	938	950	rBV	182127	342614	13.06%	0.939%
17	8.734	955	960	971	rVV	221980	381450	14.54%	1.045%
18	8.887	980	986	995	rVV7	9693	28391	1.08%	0.078%
19	9.087	1013	1020	1031	rBV	244627	415396	15.83%	1.138%
20	9.804	1135	1142	1154	rBV	213425	380848	14.51%	1.044%
21	9.975	1165	1171	1183	rBV	45900	113592	4.33%	0.311%
22	10.175	1199	1205	1211	rVV	28330	48912	1.86%	0.134%
23	10.334	1225	1232	1247	rVB	279179	512044	19.51%	1.403%
24	10.710	1287	1296	1302	rBV	782828	1301022	49.58%	3.565%
25	10.763	1302	1305	1312	rVB	34799	58139	2.22%	0.159%
26	10.893	1318	1327	1339	rBV	93989	193799	7.38%	0.531%
27	11.287	1386	1394	1409	rVV	11338	39103	1.49%	0.107%
28	11.563	1435	1441	1447	rVV	10559	23878	0.91%	0.065%
29	12.246	1553	1557	1563	rVV5	10915	27829	1.06%	0.076%
30	12.381	1575	1580	1585	rVB2	18329	31595	1.20%	0.087%
31	12.598	1612	1617	1623	rVB7	9895	17905	0.68%	0.049%
32	13.410	1745	1755	1762	rBV2	51314	95466	3.64%	0.262%
33	13.528	1768	1775	1783	rBV2	10140	28148	1.07%	0.077%
34	13.828	1823	1826	1831	rVB6	11581	16713	0.64%	0.046%
35	13.881	1831	1835	1840	rBV5	7773	13597	0.52%	0.037%
36	13.992	1848	1854	1870	rVB	552084	771526	29.40%	2.114%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
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37	14.234	1890	1895	1896	rBV5	9360	11924	0.45%	0.033%
38	14.275	1896	1902	1912	rVV2	594741	1031367	39.30%	2.826%
39	14.363	1912	1917	1921	rVV5	12647	23137	0.88%	0.063%
40	14.410	1921	1925	1929	rVB6	13292	19623	0.75%	0.054%

41	14.575	1944	1953	1960	rBV	1068270	1595666	60.80%	4.372%
42	14.640	1960	1964	1970	rVB	70944	105840	4.03%	0.290%
43	14.775	1981	1987	1994	rBV9	13553	24979	0.95%	0.068%
44	14.851	1994	2000	2013	rBV3	66223	180825	6.89%	0.495%
45	14.981	2016	2022	2027	rBV	44063	69430	2.65%	0.190%

46	15.181	2051	2056	2060	rBV7	7685	16306	0.62%	0.045%
47	15.363	2084	2087	2090	rVV5	9040	10293	0.39%	0.028%
48	15.398	2090	2093	2099	rVB7	7796	12336	0.47%	0.034%
49	15.457	2100	2103	2107	rBV5	7180	10119	0.39%	0.028%
50	15.581	2114	2124	2130	rBV	759868	1088401	41.47%	2.982%

51	15.639	2130	2134	2145	rVB	80171	132540	5.05%	0.363%
52	15.739	2146	2151	2159	rBV2	63947	120449	4.59%	0.330%
53	15.928	2179	2183	2186	rBV3	16789	21934	0.84%	0.060%
54	16.069	2203	2207	2211	rBV3	14059	28726	1.09%	0.079%
55	16.175	2222	2225	2228	rVB5	8669	10711	0.41%	0.029%

56	16.245	2230	2237	2240	rVV3	19359	39250	1.50%	0.108%
57	16.281	2240	2243	2250	rVB5	25680	42737	1.63%	0.117%
58	16.351	2251	2255	2261	rBV	29936	45406	1.73%	0.124%
59	16.645	2295	2305	2309	rBV5	28307	69284	2.64%	0.190%
60	16.804	2328	2332	2333	rBV4	8908	11535	0.44%	0.032%

61	16.875	2341	2344	2347	rBV3	14686	16755	0.64%	0.046%
62	17.033	2362	2371	2375	rBV4	37113	98896	3.77%	0.271%
63	17.081	2375	2379	2386	rVB5	32356	59514	2.27%	0.163%
64	17.186	2393	2397	2400	rBV	27877	32225	1.23%	0.088%
65	17.369	2422	2428	2432	rBV	1214534	1697248	64.67%	4.651%

66	17.410	2432	2435	2440	rVV	322577	437731	16.68%	1.199%
67	17.469	2440	2445	2448	rVV	778366	1050071	40.01%	2.877%
68	17.504	2448	2451	2458	rVV	549070	775110	29.54%	2.124%
69	17.586	2461	2465	2468	rVB2	23875	37739	1.44%	0.103%
70	17.804	2498	2502	2508	rVB	88372	132006	5.03%	0.362%

71	17.886	2513	2516	2523	rVB	39639	59278	2.26%	0.162%
72	18.222	2569	2573	2580	rBV3	261648	420300	16.02%	1.152%
73	18.286	2581	2584	2594	rVB2	56867	113375	4.32%	0.311%
74	18.369	2594	2598	2602	rBV	68266	91726	3.50%	0.251%
75	18.433	2604	2609	2613	rBV	173280	273484	10.42%	0.749%

76	18.604	2635	2638	2640	rBV4	31758	42552	1.62%	0.117%
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77	18.728	2656	2659	2663	rVB	99742	123027	4.69%	0.337%
78	18.792	2668	2670	2673	rVB2	57887	67399	2.57%	0.185%
79	18.933	2690	2694	2697	rBV	166451	198044	7.55%	0.543%
80	19.051	2710	2714	2717	rBV2	96966	149484	5.70%	0.410%
81	19.128	2723	2727	2731	rVB3	71255	101606	3.87%	0.278%
82	19.292	2751	2755	2763	rVB	253320	388075	14.79%	1.063%
83	19.398	2768	2773	2781	rVV2	1144369	1980099	75.45%	5.426%
84	19.469	2782	2785	2795	rVV7	103744	200102	7.62%	0.548%
85	19.545	2796	2798	2804	rVB	50717	64220	2.45%	0.176%
86	19.727	2823	2829	2832	rBV2	1040320	1677037	63.90%	4.595%
87	19.763	2832	2835	2841	rVV	1149906	1534832	58.48%	4.205%
88	19.822	2842	2845	2851	rVB2	120300	169147	6.45%	0.463%
89	19.933	2861	2864	2867	rBV	115160	125927	4.80%	0.345%
90	20.145	2895	2900	2904	rBV	971410	1205996	45.95%	3.304%
91	20.339	2930	2933	2936	rBV	171388	239494	9.13%	0.656%
92	20.539	2964	2967	2972	rVB2	169008	194585	7.41%	0.533%
93	20.616	2977	2980	2983	rVV	154656	168812	6.43%	0.463%
94	20.663	2983	2988	2994	rVB	793560	1018127	38.80%	2.790%
95	21.186	3069	3077	3082	rBV5	382996	877209	33.43%	2.404%
96	21.516	3127	3133	3137	rBV2	1329640	2494462	95.05%	6.835%
97	21.557	3137	3140	3144	rVB	789098	967602	36.87%	2.651%
98	22.145	3236	3240	3248	rVB2	553414	1114323	42.46%	3.053%
99	23.286	3429	3434	3448	rVB	730193	2624340	100.00%	7.191%
100	24.080	3564	3569	3575	rVB	666233	1274959	48.58%	3.493%

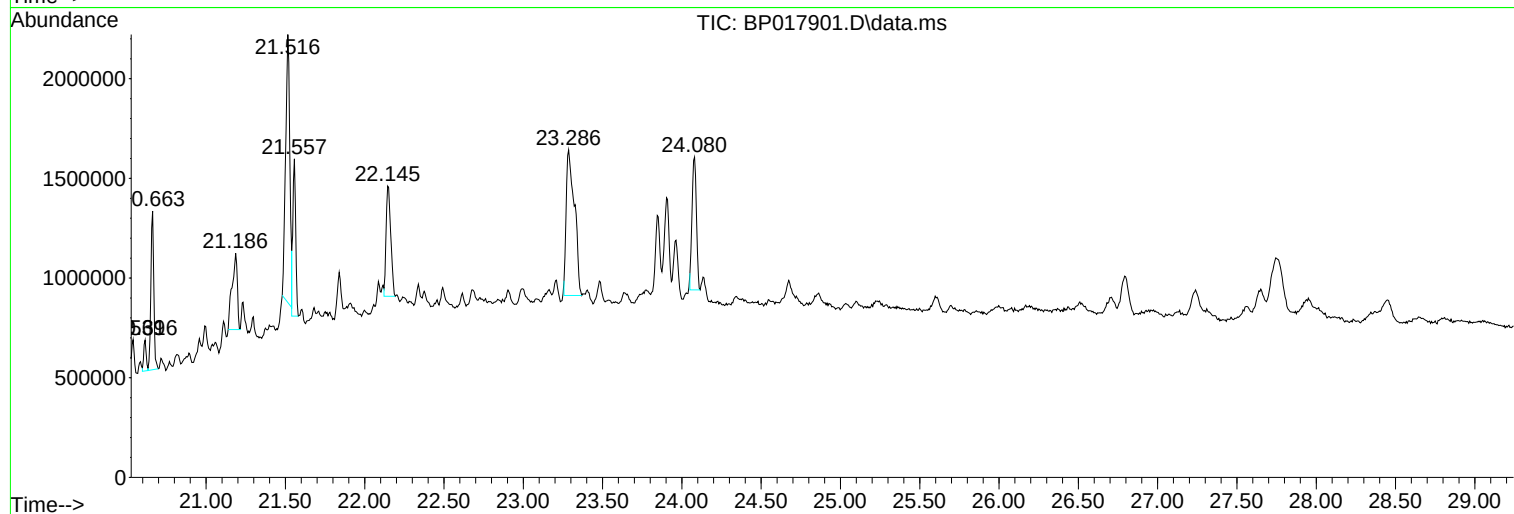
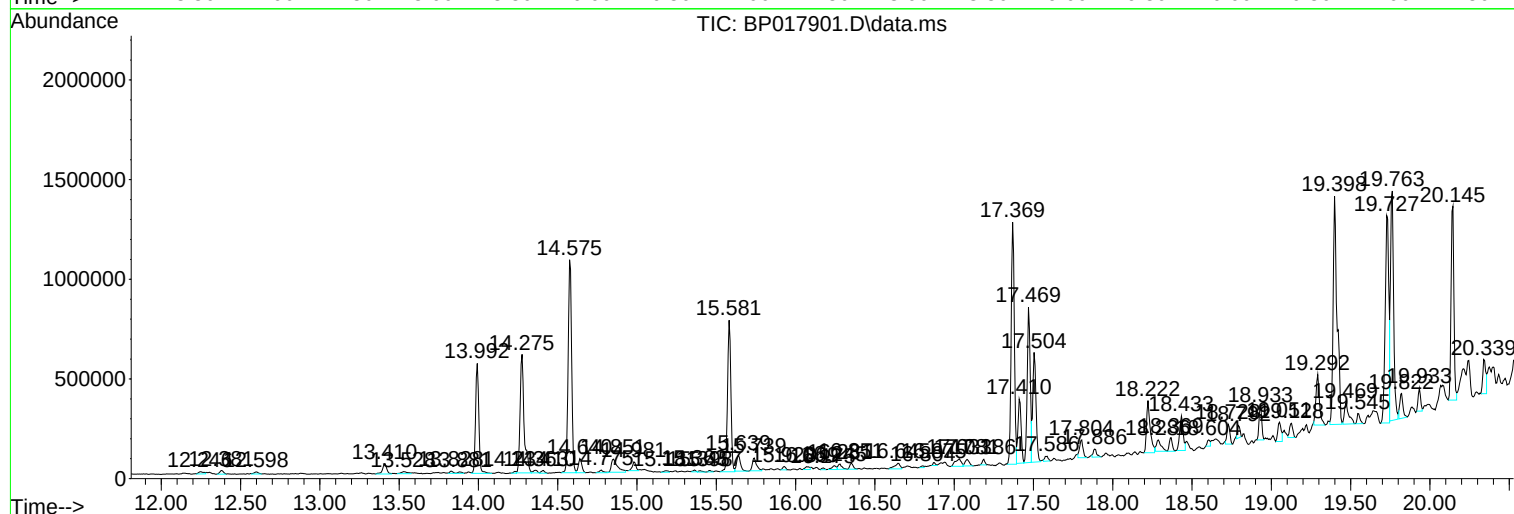
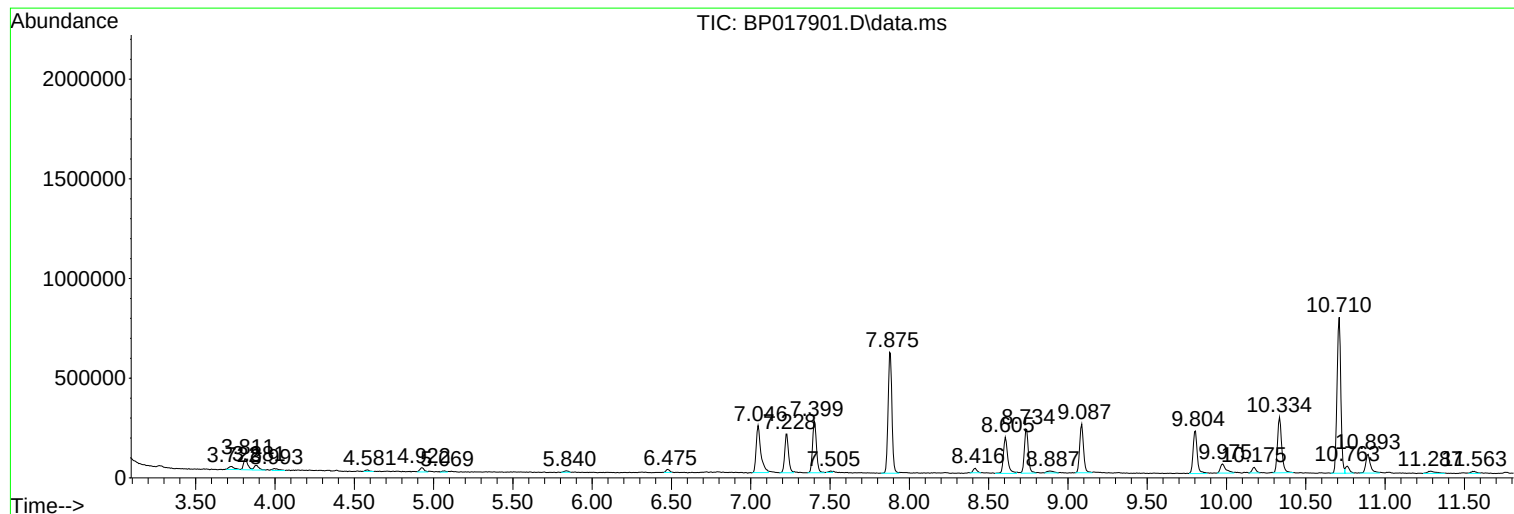
Sum of corrected areas: 36495997

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

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



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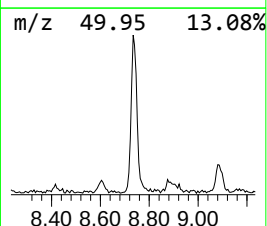
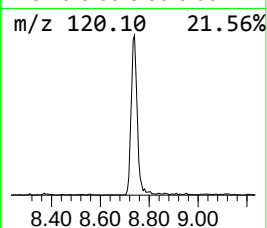
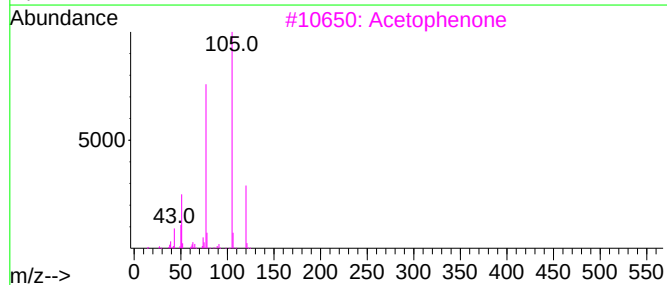
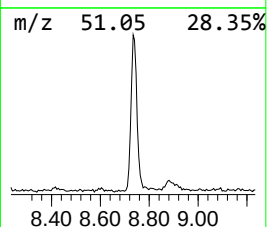
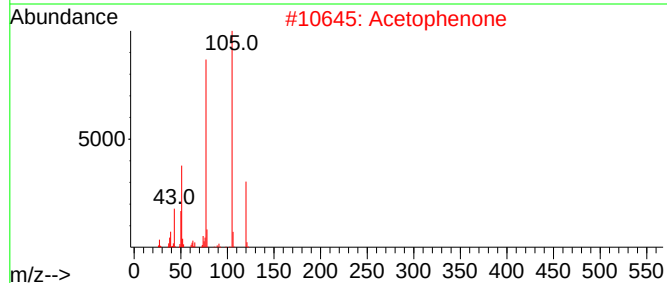
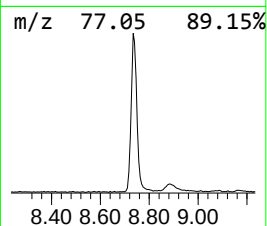
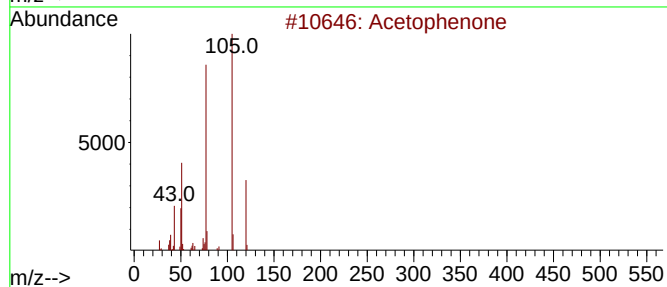
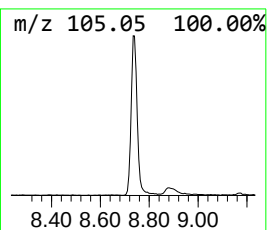
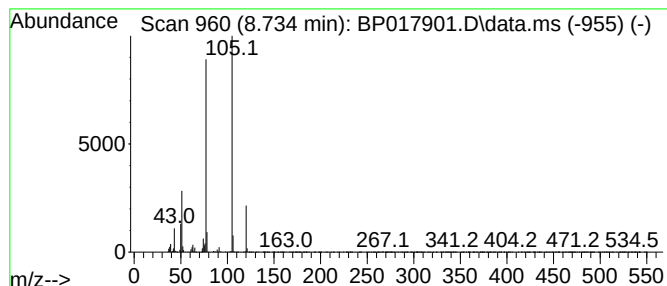
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.734	7.50 ng/ul	381450	1,4-Dichlorobenzene-d4	7.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone			120	C8H8O	000098-86-2	95
2	Acetophenone			120	C8H8O	000098-86-2	94
3	Acetophenone			120	C8H8O	000098-86-2	94
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



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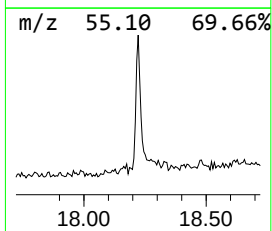
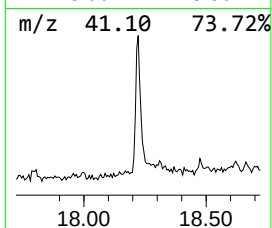
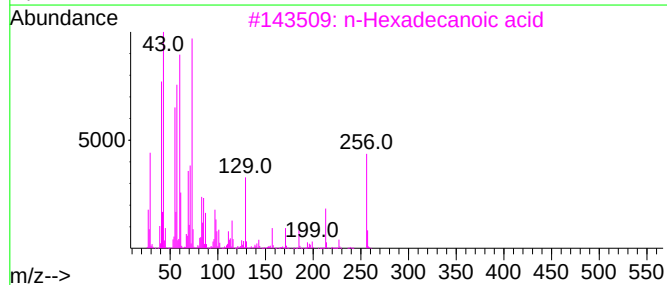
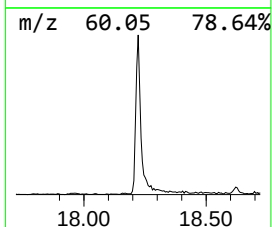
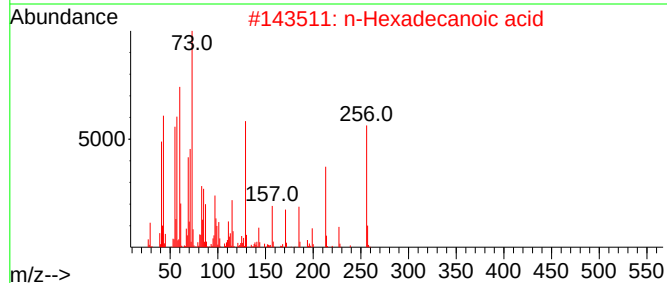
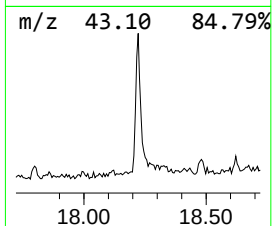
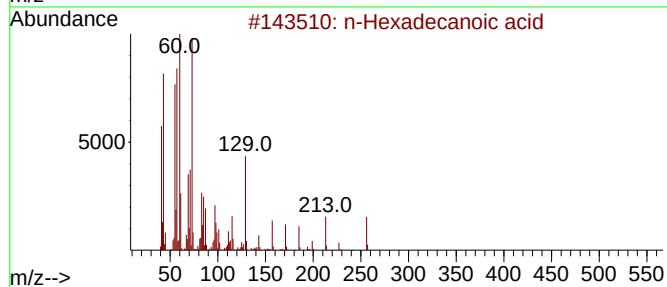
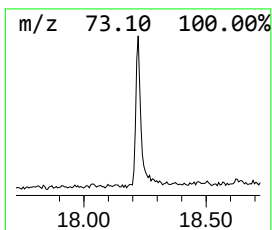
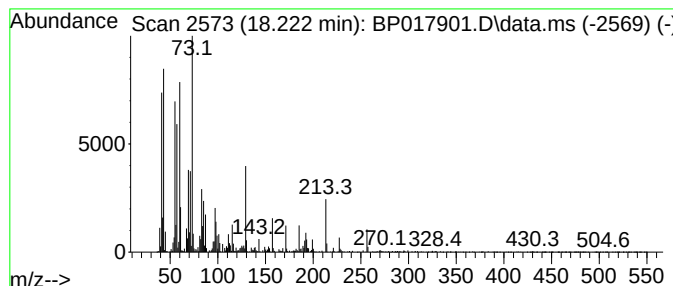
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

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.
18.222	4.95 ng/ul	420300	Phenanthrene-d10	17.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



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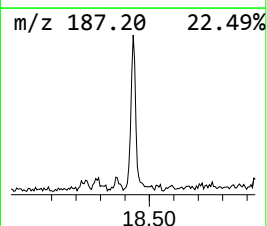
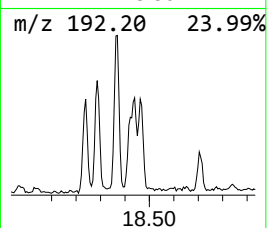
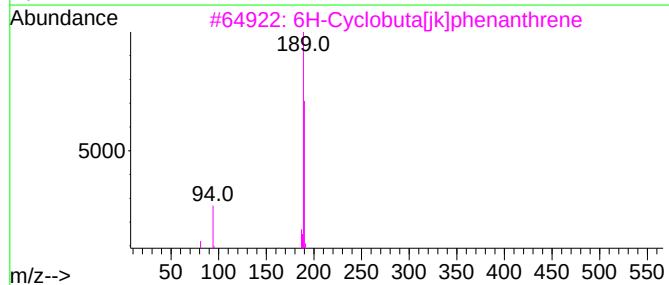
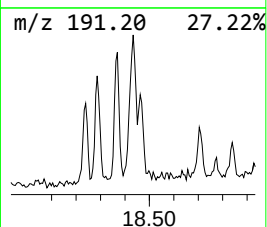
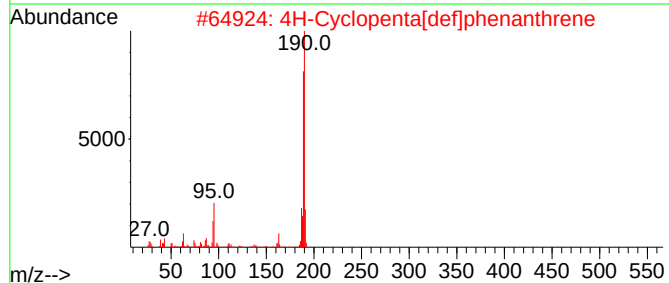
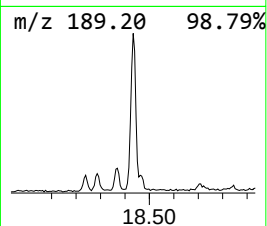
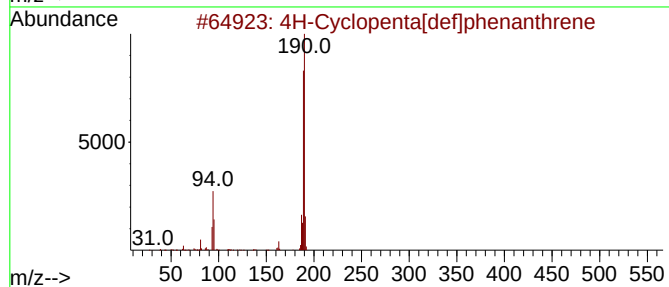
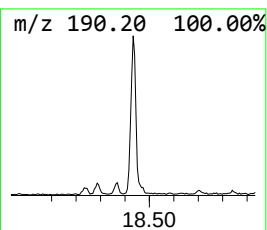
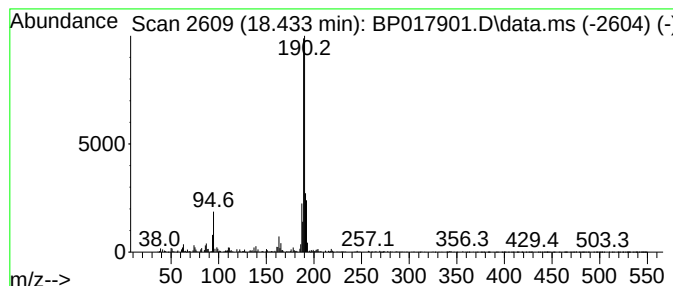
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.433	3.22 ng/ul	273484	Phenanthrene-d10	17.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
4			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
Operator : MA/JU
Sample : 05060-02
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKG0

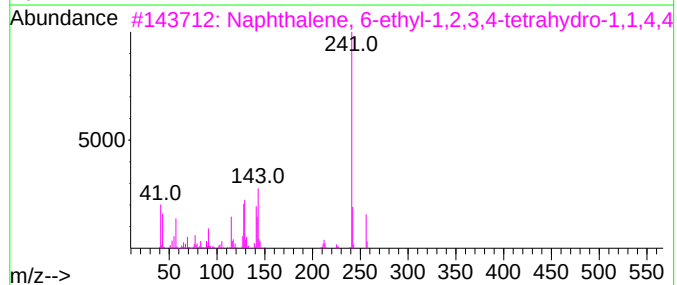
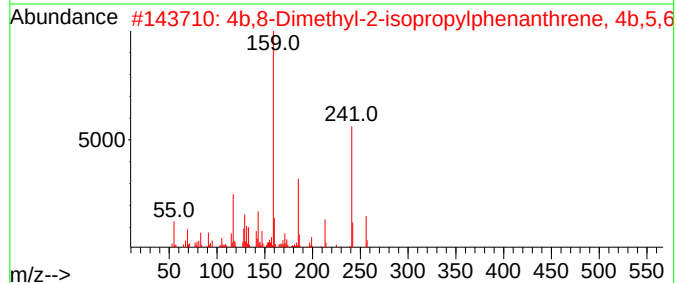
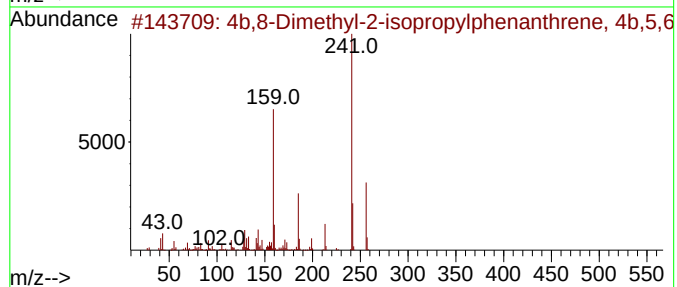
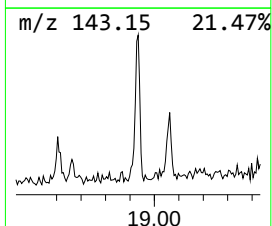
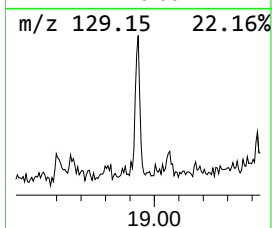
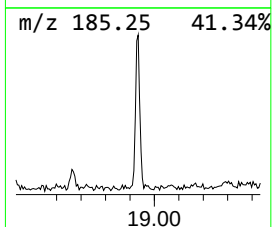
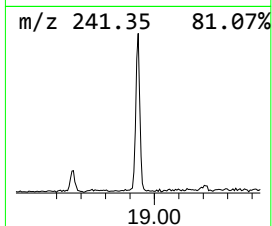
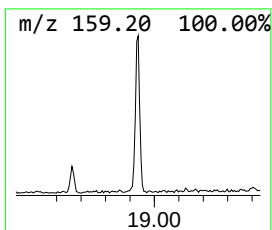
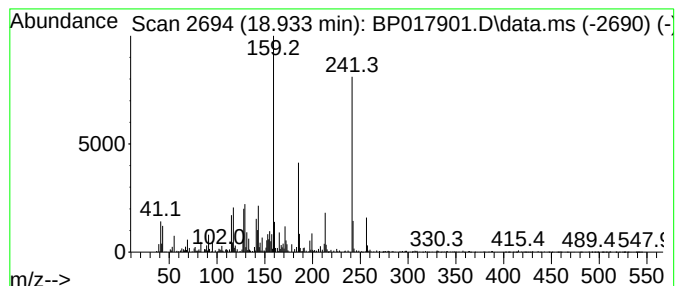
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 4b,8-Dimethyl-2-isopropylph... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.933	2.33 ng/ul	198044	Phenanthrene-d10	17.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	90
2			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	87
3			Naphthalene, 6-ethyl-1,2,3,4-tet...	256	C19H28	301643-35-6	64
4			Decanedioic acid, dibutyl ester	314	C18H34O4	000109-43-3	38
5			Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
Operator : MA/JU
Sample : 05060-02
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKG0

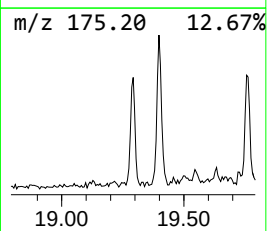
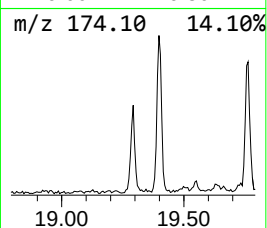
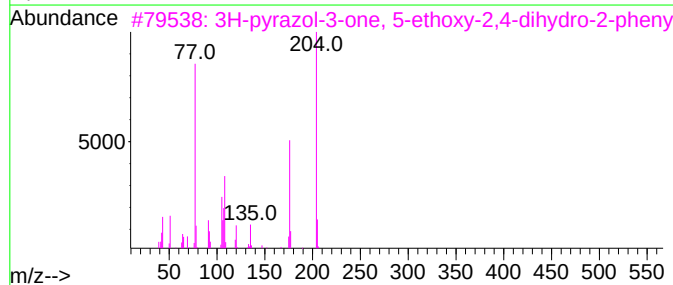
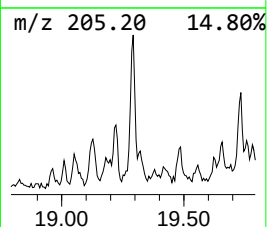
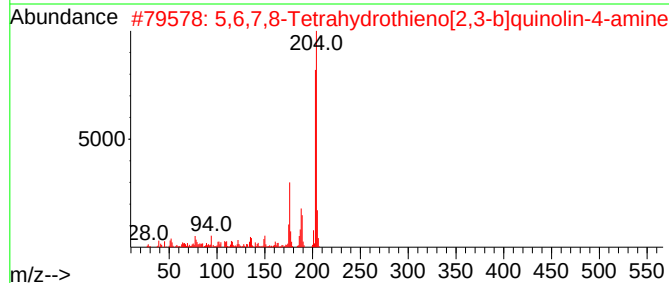
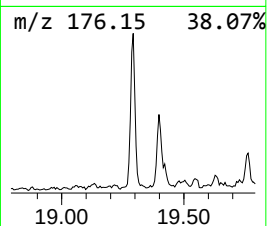
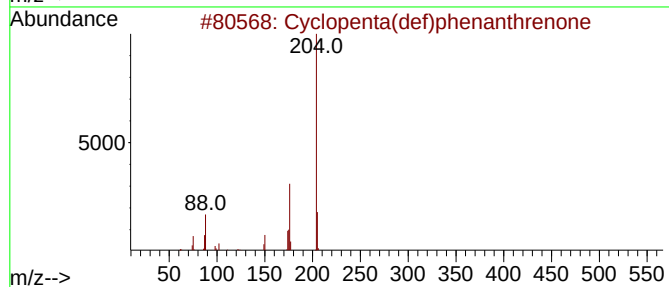
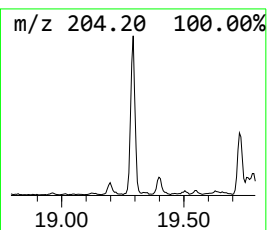
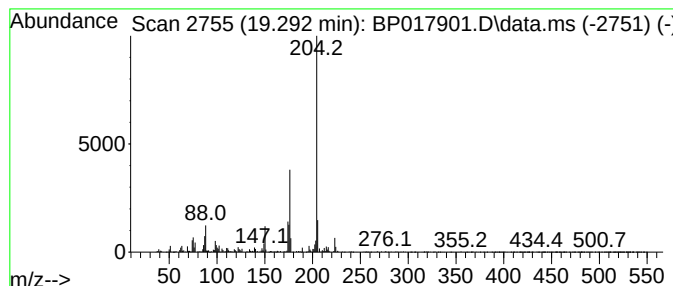
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclopenta(def)phenanthrenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.292	4.57 ng/ul	388075	Phenanthrene-d10	17.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	91
2			5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	64
3			3H-pyrazol-3-one, 5-ethoxy-2,4-d...	204	C11H12N2O2	1000400-47-1	59
4			4-Methyl-6,7-methylenedioxcoumarin	204	C11H8O4	015071-04-2	58
5			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
Operator : MA/JU
Sample : 05060-02
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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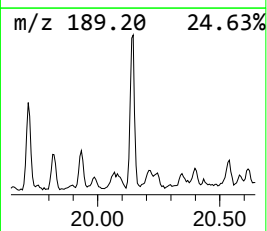
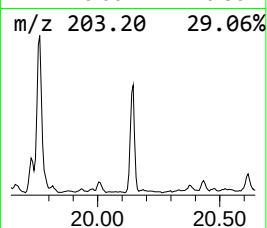
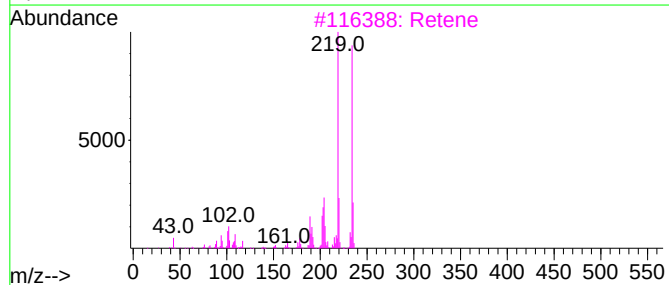
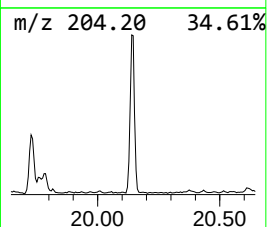
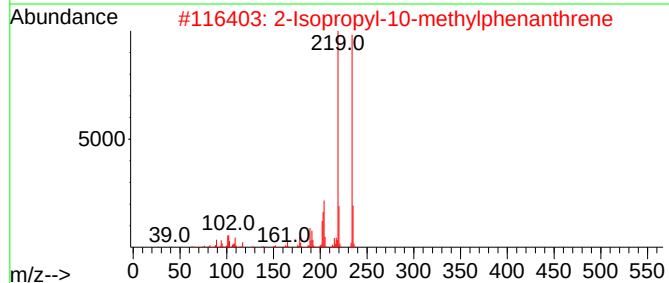
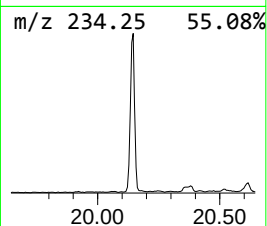
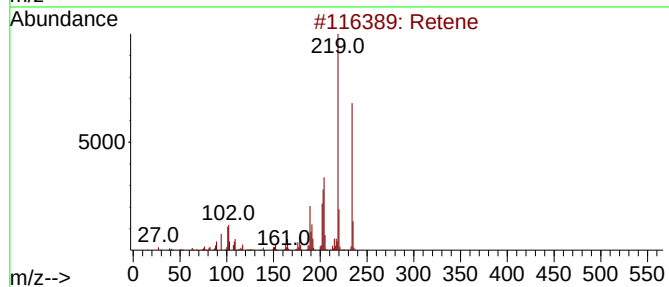
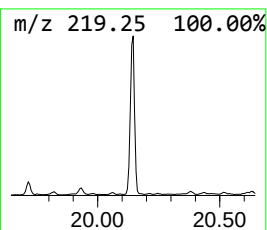
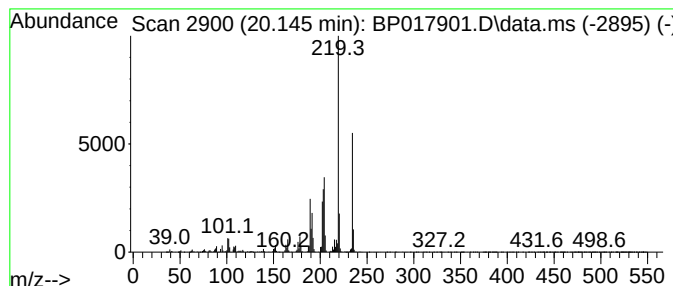
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Retene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.145	9.67 ng/ul	1206000	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Retene			234	C18H18	000483-65-8	98
2	2-Isopropyl-10-methylphenanthrene			234	C18H18	066552-97-4	94
3	Retene			234	C18H18	000483-65-8	94
4	Retene			234	C18H18	000483-65-8	93
5	Phenanthrene, 2,4,5,7-tetramethyl-			234	C18H18	007396-38-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
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Sample : 05060-02
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ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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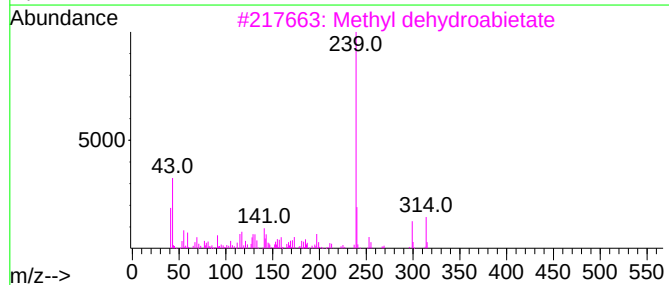
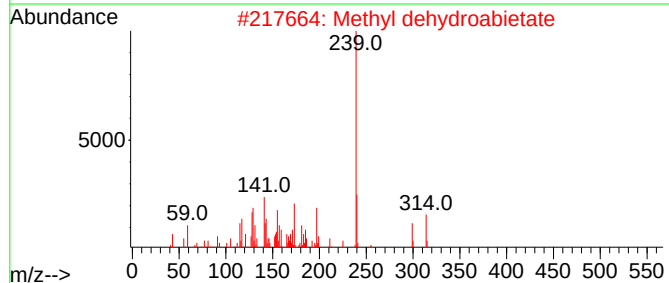
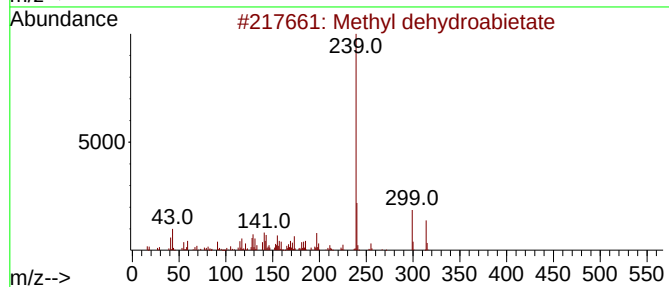
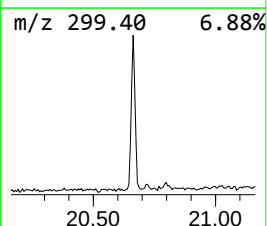
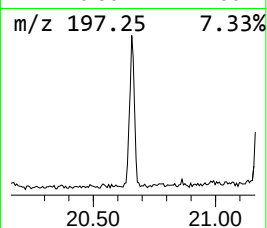
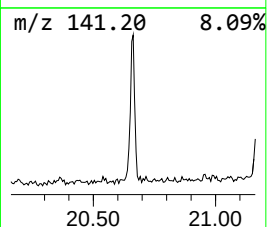
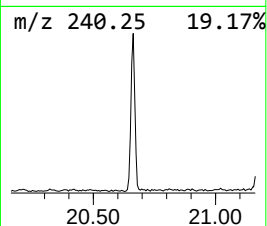
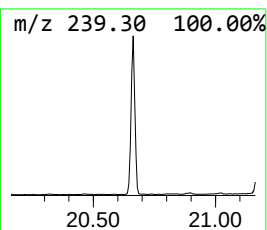
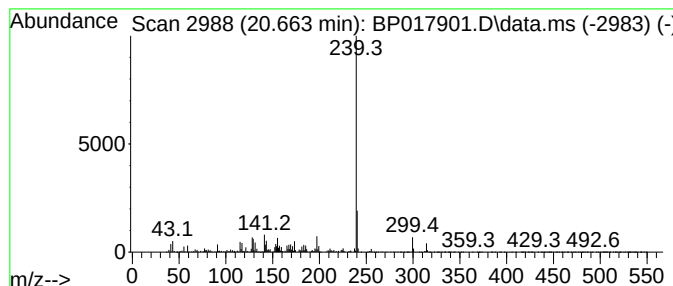
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methyl dehydroabietate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.663	8.16 ng/ul	1018130	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl dehydroabietate	314	C21H30O2	001235-74-1	97
2			Methyl dehydroabietate	314	C21H30O2	001235-74-1	91
3			Methyl dehydroabietate	314	C21H30O2	001235-74-1	89
4			Methyl dehydroabietate	314	C21H30O2	001235-74-1	87
5			Indole, 3-imidazolo[2,1-b]thiazo...	239	C13H9N3S	292855-05-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
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Sample : 05060-02
Misc :
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Instrument :
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ClientSampleId :
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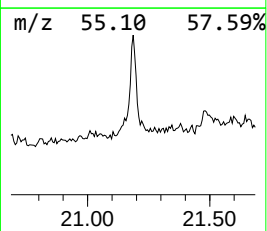
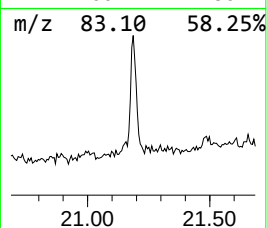
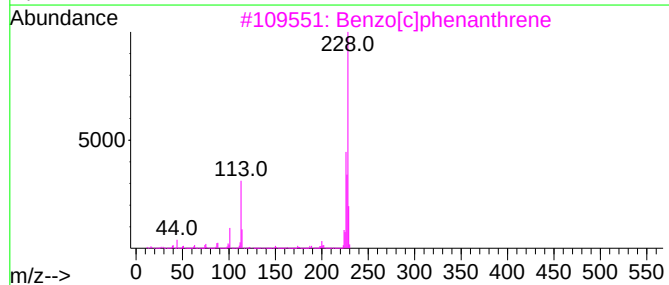
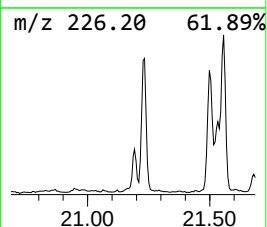
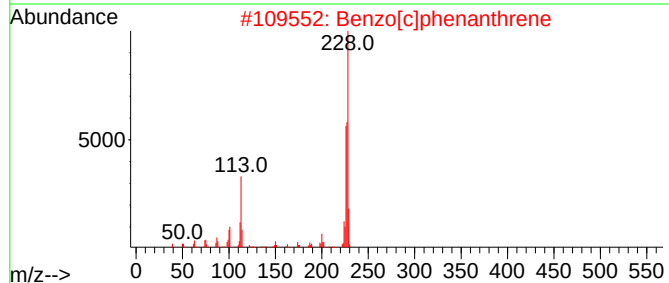
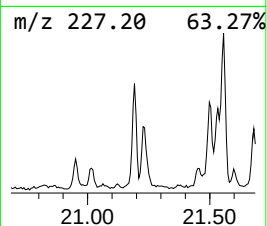
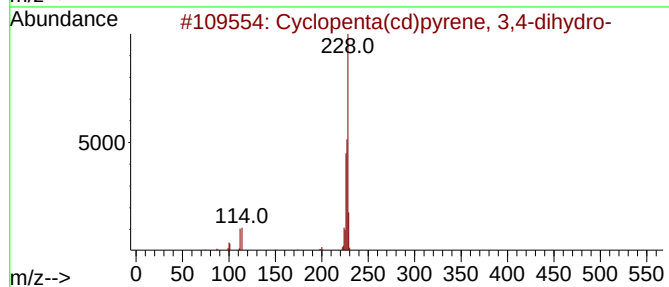
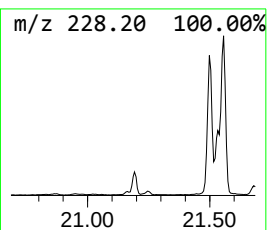
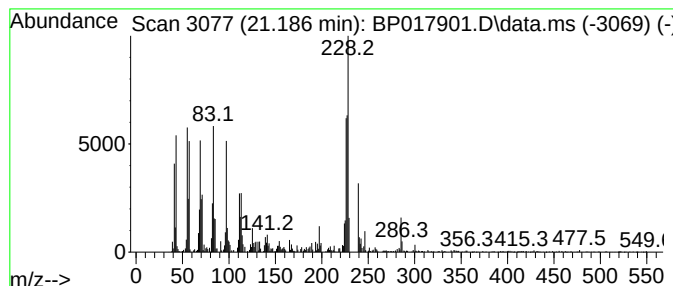
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.186	7.03 ng/ul	877209	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	38
2			Benzo[c]phenanthrene	228	C18H12	000195-19-7	38
3			Benzo[c]phenanthrene	228	C18H12	000195-19-7	30
4			Purine-6(1H)-thione, 3,7-dimethy...	228	C7H8N4Se	023663-58-3	30
5			Benzo[c]phenanthrene	228	C18H12	000195-19-7	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
Operator : MA/JU
Sample : 05060-02
Misc :
ALS Vial : 56 Sample Multiplier: 1

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ClientSampleId :
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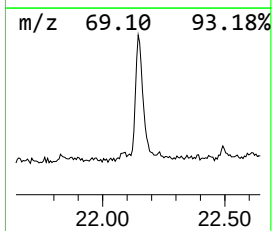
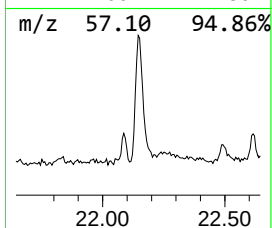
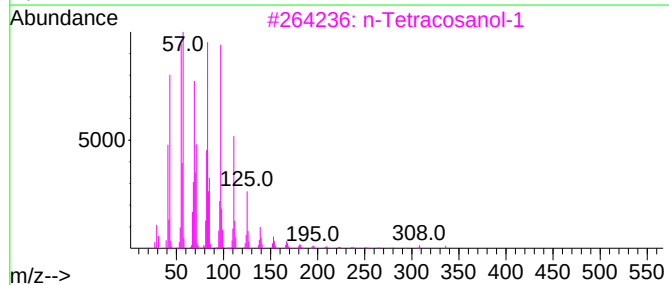
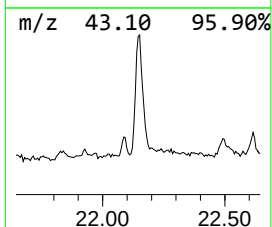
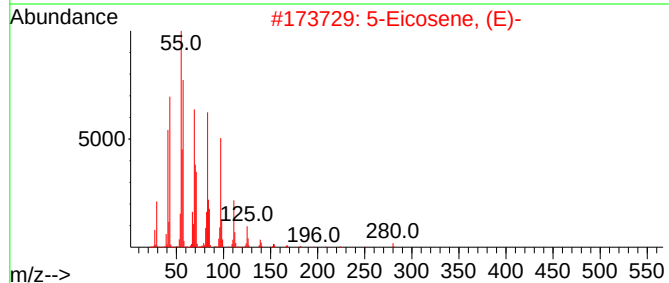
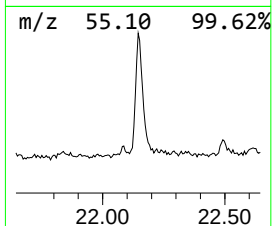
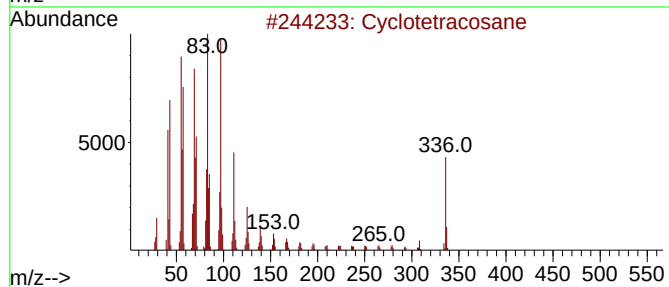
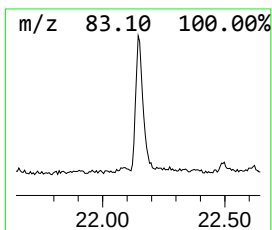
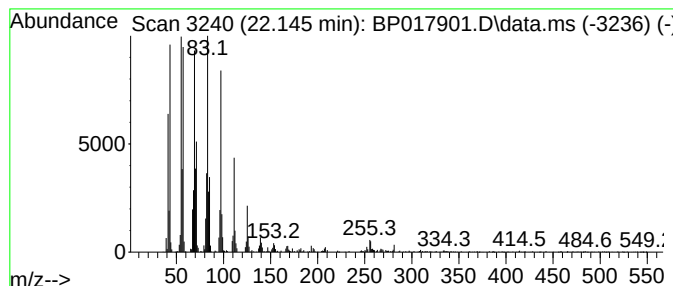
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic22.145 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.145	8.93 ng/ul	1114320	Chrysene-d12	21.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	98
2			5-Eicosene, (E)-	280	C20H40	074685-30-6	98
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110123\
Data File : BP017901.D
Acq On : 02 Nov 2023 23:25
Operator : MA/JU
Sample : 05060-02
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCKG0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.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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n-Hexadecanoic ...	18.222	5.0	ng/ul	420300	4	17.369	1697250	20.0
4H-Cyclopenta[d...]	18.433	3.2	ng/ul	273484	4	17.369	1697250	20.0
4b,8-Dimethyl-2...	18.933	2.3	ng/ul	198044	4	17.369	1697250	20.0
Cyclopenta(def)...	19.292	4.6	ng/ul	388075	4	17.369	1697250	20.0
Retene	20.145	9.7	ng/ul	1206000	5	21.516	2494460	20.0
Methyl dehydroa...	20.663	8.2	ng/ul	1018130	5	21.516	2494460	20.0
unknown-01	21.186	7.0	ng/ul	877209	5	21.516	2494460	20.0
(DEL) Alkane: C...	22.145	8.9	ng/ul	1114320	5	21.516	2494460	20.0