

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019127.D
 Acq On : 28 Dec 2023 16:12
 Operator : MA/JU
 Sample : 05920-15
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCF66

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019127.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	21	24	27	rBV	35941	42606	0.44%	0.037%
2	3.264	27	30	39	rVB	83213	105478	1.08%	0.092%
3	3.570	79	82	89	rVB3	14266	19689	0.20%	0.017%
4	3.693	99	103	108	rVB	79008	93821	0.96%	0.082%
5	4.311	202	208	223	rVB	6943390	9762873	100.00%	8.513%
6	4.917	304	311	327	rBV	3261362	4504301	46.14%	3.928%
7	5.705	440	445	453	rVB	35490	56375	0.58%	0.049%
8	5.993	489	494	503	rVB	51273	76689	0.79%	0.067%
9	6.328	548	551	557	rBV7	6230	12989	0.13%	0.011%
10	6.975	651	661	671	rBV	535285	855405	8.76%	0.746%
11	7.140	683	689	702	rVB	459118	698229	7.15%	0.609%
12	7.334	715	722	729	rBV	542327	858393	8.79%	0.749%
13	7.799	795	801	808	rVB	417266	668996	6.85%	0.583%
14	8.058	840	845	854	rVB	8553	17665	0.18%	0.015%
15	8.340	883	893	899	rBV2	219480	554838	5.68%	0.484%
16	8.387	899	901	908	rVB2	23431	34821	0.36%	0.030%
17	8.522	917	924	938	rBV	606188	993825	10.18%	0.867%
18	8.716	953	957	967	rBV6	9542	20204	0.21%	0.018%
19	8.963	991	999	1003	rBV	481119	797154	8.17%	0.695%
20	9.010	1003	1007	1015	rVB	370227	602193	6.17%	0.525%
21	9.240	1041	1046	1052	rVB	19082	30913	0.32%	0.027%
22	9.375	1063	1069	1080	rVB3	20680	45571	0.47%	0.040%
23	9.687	1115	1122	1125	rBV	381541	660315	6.76%	0.576%
24	9.716	1125	1127	1139	rVB	386598	578205	5.92%	0.504%
25	10.022	1173	1179	1189	rBV	331649	506557	5.19%	0.442%
26	10.228	1207	1214	1217	rBV	655914	1254192	12.85%	1.094%
27	10.599	1265	1277	1282	rBV	535529	943040	9.66%	0.822%
28	10.652	1282	1286	1295	rVV	423103	676433	6.93%	0.590%
29	10.746	1295	1302	1313	rVV	395848	707265	7.24%	0.617%
30	10.934	1330	1334	1339	rVB7	9814	13985	0.14%	0.012%
31	11.899	1489	1498	1511	rBV	1376057	2225222	22.79%	1.940%
32	12.140	1532	1539	1545	rBV2	20731	41675	0.43%	0.036%
33	12.193	1545	1548	1553	rVB2	12441	19292	0.20%	0.017%
34	12.269	1553	1561	1569	rBV	1188417	1843750	18.89%	1.608%
35	12.399	1577	1583	1589	rBV3	34661	63005	0.65%	0.055%
36	12.487	1590	1598	1605	rVB	995208	1543507	15.81%	1.346%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
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37	12.610	1611	1619	1626	rBV	1006027	1480369	15.16%	1.291%
38	12.881	1659	1665	1677	rVB	441997	681240	6.98%	0.594%
39	13.275	1726	1732	1735	rVV2	33974	54118	0.55%	0.047%
40	13.322	1735	1740	1742	rVV	878304	1229922	12.60%	1.072%
41	13.357	1742	1746	1762	rVB	2012329	3209885	32.88%	2.799%
42	13.734	1806	1810	1817	rVB7	8393	15150	0.16%	0.013%
43	13.869	1827	1833	1840	rBV	1205901	1647711	16.88%	1.437%
44	14.040	1856	1862	1869	rVB	303779	404170	4.14%	0.352%
45	14.145	1869	1880	1883	rBV	1321177	1999149	20.48%	1.743%
46	14.451	1925	1932	1938	rBV	790329	1186065	12.15%	1.034%
47	14.516	1938	1943	1951	rVB	738636	1031543	10.57%	0.899%
48	14.669	1962	1969	1979	rBV	453847	741967	7.60%	0.647%
49	14.828	1989	1996	2003	rBV	786669	1069726	10.96%	0.933%
50	15.081	2033	2039	2052	rBV	1221475	1743380	17.86%	1.520%
51	15.245	2063	2067	2073	rBV5	9658	15061	0.15%	0.013%
52	15.310	2073	2078	2087	rVB4	20396	37977	0.39%	0.033%
53	15.451	2093	2102	2106	rBV	1771461	2545376	26.07%	2.220%
54	15.504	2106	2111	2118	rVV	1734603	2388168	24.46%	2.082%
55	15.592	2118	2126	2138	rVV2	2057362	3210219	32.88%	2.799%
56	15.687	2139	2142	2149	rVB3	14389	27109	0.28%	0.024%
57	16.098	2209	2212	2217	rBV5	9144	13973	0.14%	0.012%
58	16.181	2221	2226	2233	rBV5	17284	37215	0.38%	0.032%
59	16.398	2257	2263	2272	rBV	1141234	1566785	16.05%	1.366%
60	16.510	2272	2282	2290	rVB	1379852	1947712	19.95%	1.698%
61	16.581	2290	2294	2298	rBV4	19098	32145	0.33%	0.028%
62	16.687	2308	2312	2318	rVB9	10208	14779	0.15%	0.013%
63	16.863	2339	2342	2350	rVB3	19095	35847	0.37%	0.031%
64	17.210	2395	2401	2405	rBV	1042102	1478938	15.15%	1.290%
65	17.251	2405	2408	2413	rVV	1054159	1430738	14.65%	1.248%
66	17.310	2413	2418	2422	rVV	2116769	3061720	31.36%	2.670%
67	17.345	2422	2424	2436	rVB	682565	804796	8.24%	0.702%
68	17.492	2446	2449	2456	rVB4	12577	17004	0.17%	0.015%
69	17.586	2460	2465	2468	rBV7	8731	17460	0.18%	0.015%
70	17.692	2478	2483	2492	rBV	1288710	1682750	17.24%	1.467%
71	17.892	2512	2517	2521	rBV2	59973	72874	0.75%	0.064%
72	18.104	2549	2553	2561	rBV	212937	312107	3.20%	0.272%
73	18.175	2562	2565	2570	rVB3	34518	51035	0.52%	0.045%
74	18.392	2599	2602	2606	rBV5	12917	14406	0.15%	0.013%
75	18.610	2634	2639	2647	rVB	316276	417833	4.28%	0.364%
76	18.951	2693	2697	2699	rBV5	16032	24062	0.25%	0.021%

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77	18.998	2702	2705	2707	rVV2	51835	62742	0.64%	0.055%
78	19.028	2707	2710	2715	rVB	283651	313906	3.22%	0.274%
79	19.128	2724	2727	2730	rBV3	36995	53289	0.55%	0.046%
80	19.157	2730	2732	2736	rVV	39225	40378	0.41%	0.035%
81	19.233	2736	2745	2753	rVV	3243499	4744585	48.60%	4.137%
82	19.345	2761	2764	2768	rBV	113611	122024	1.25%	0.106%
83	19.557	2794	2800	2803	rBV	3050103	4374753	44.81%	3.815%
84	19.580	2803	2804	2810	rVB	1400643	1447023	14.82%	1.262%
85	20.086	2888	2890	2895	rVB2	57958	63545	0.65%	0.055%
86	20.416	2943	2946	2950	rVB2	54671	54673	0.56%	0.048%
87	20.533	2963	2966	2970	rBV	244992	247337	2.53%	0.216%
88	20.575	2970	2973	2976	rVB	58765	61832	0.63%	0.054%
89	21.039	3047	3052	3059	rBV2	294512	595475	6.10%	0.519%
90	21.098	3059	3062	3069	rVB	139675	160729	1.65%	0.140%
91	21.304	3092	3097	3102	rBV2	2139601	4045994	41.44%	3.528%
92	21.351	3102	3105	3111	rVB	1796674	2092877	21.44%	1.825%
93	22.110	3230	3234	3241	rVB	287094	410722	4.21%	0.358%
94	22.963	3373	3379	3384	rBV	4264904	7680349	78.67%	6.697%
95	23.010	3384	3387	3398	rVB	1433408	2204101	22.58%	1.922%
96	23.533	3469	3476	3481	rBV2	2449971	4829304	49.47%	4.211%
97	23.580	3481	3484	3493	rVB	581170	979352	10.03%	0.854%
98	23.686	3496	3502	3512	rVB	1102547	2171684	22.24%	1.894%
99	26.157	3911	3922	3936	rVB3	1555067	5584335	57.20%	4.869%
100	26.898	4040	4048	4060	rVB	829993	2657158	27.22%	2.317%

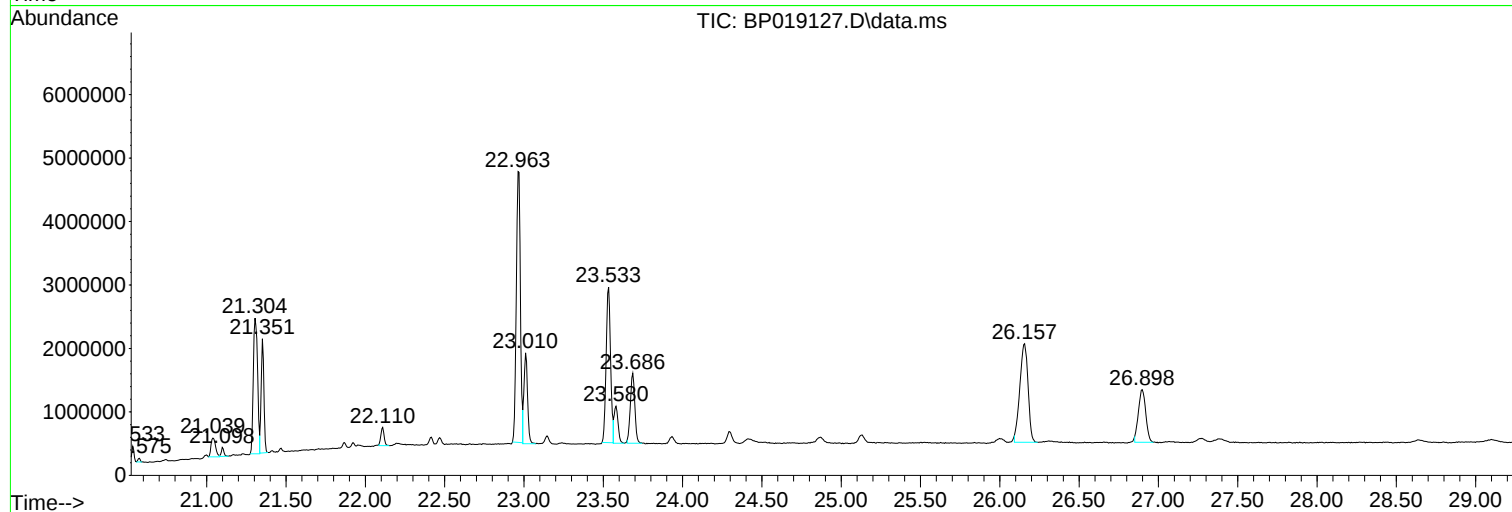
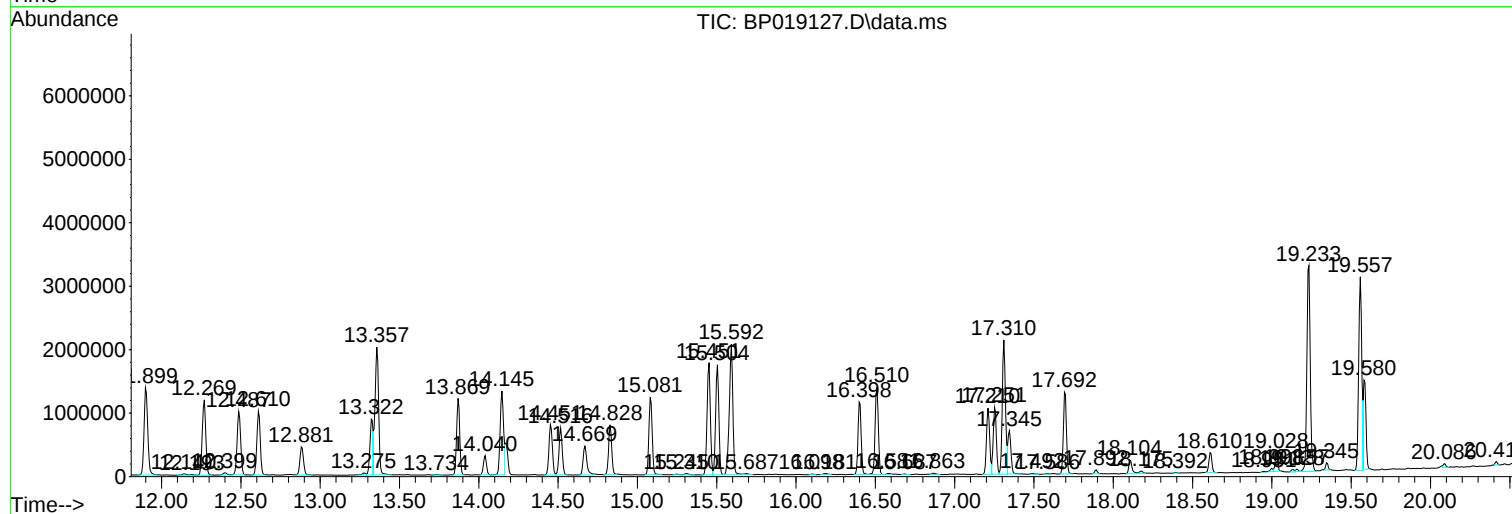
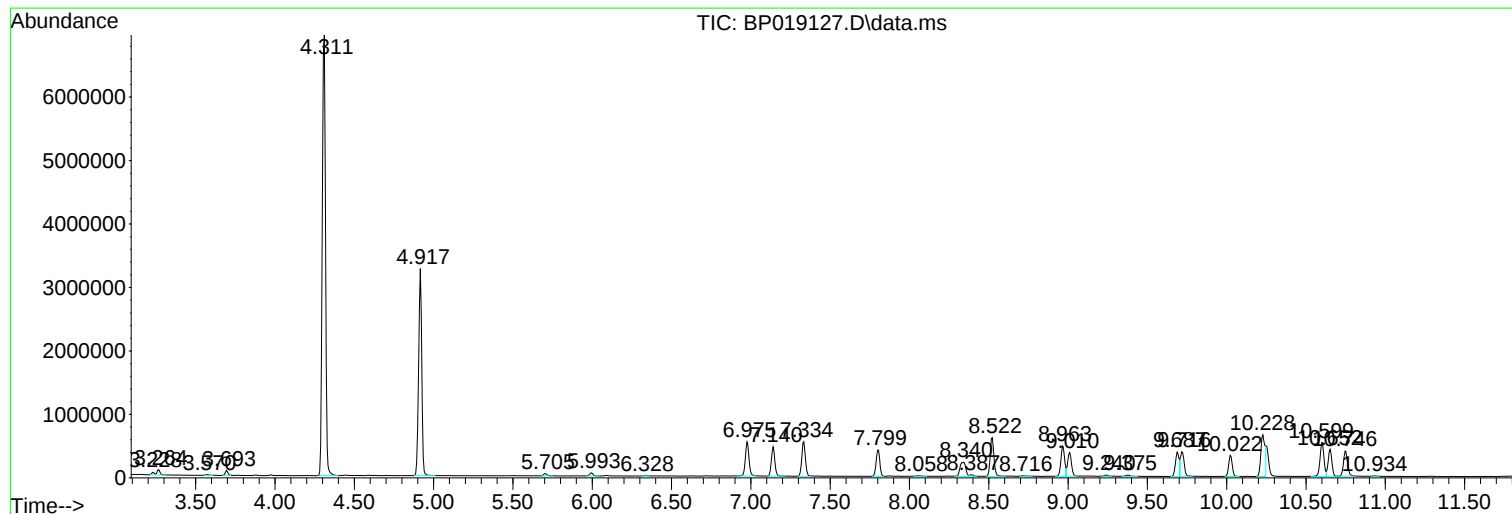
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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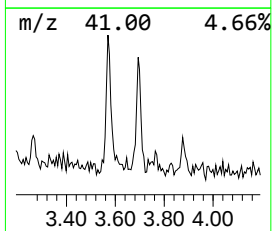
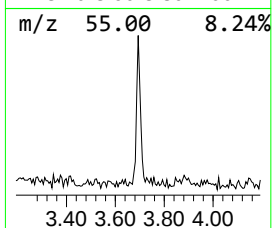
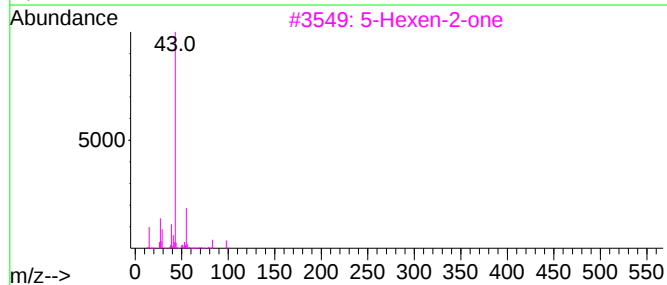
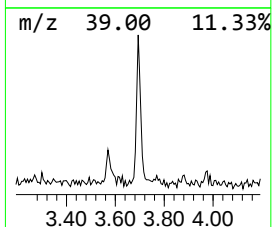
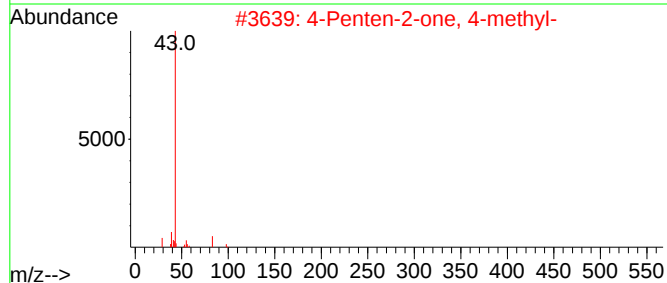
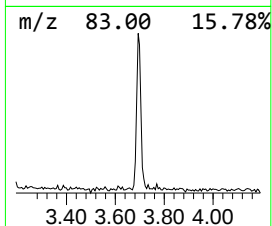
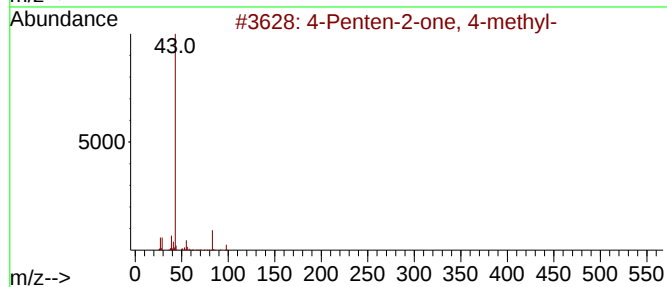
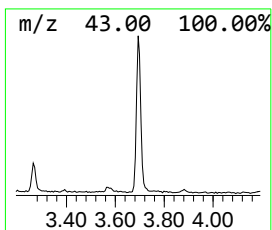
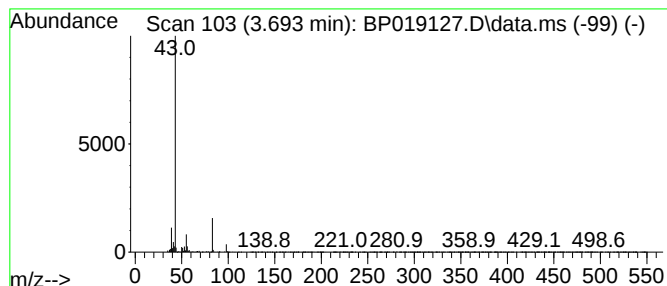
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.693	2.80 ng/ul	93821	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	72
2			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	53
3			5-Hexen-2-one	98	C6H10O	000109-49-9	38
4			5-Hexen-2-one	98	C6H10O	000109-49-9	38
5			2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	37



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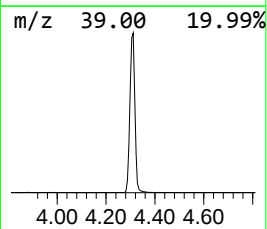
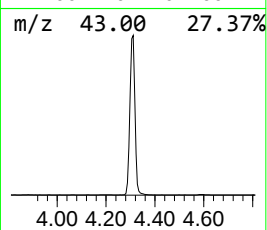
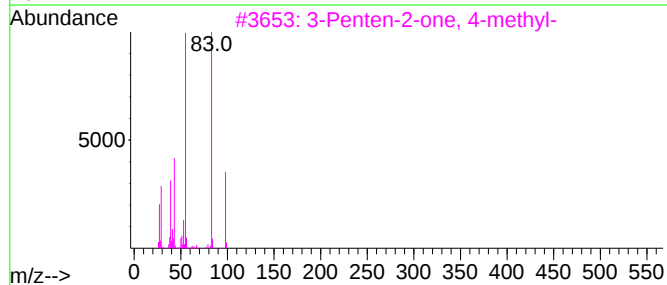
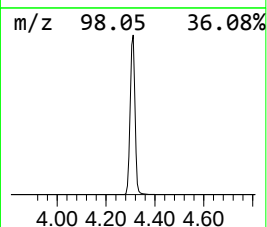
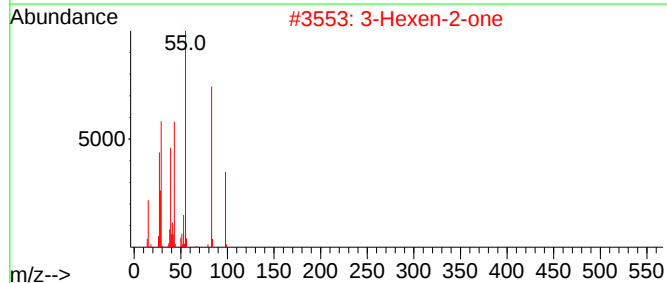
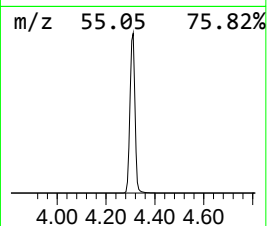
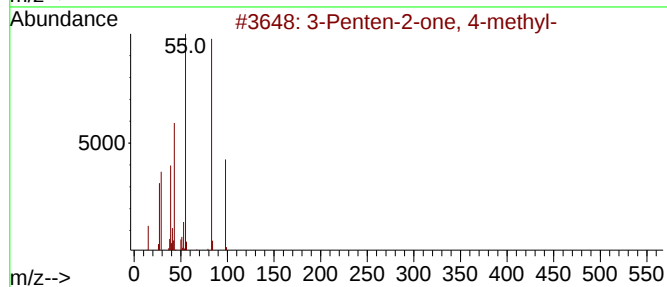
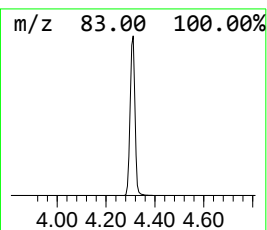
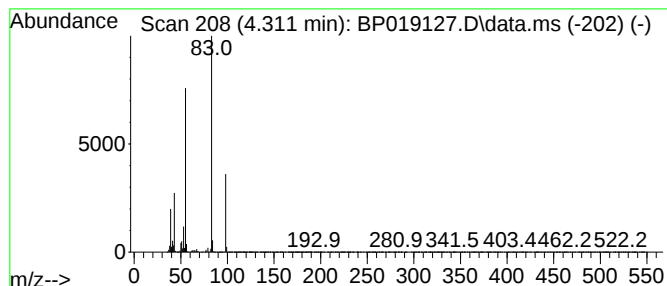
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.311	291.87 ng/ul	9762870	1,4-Dichlorobenzene-d4	7.799

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



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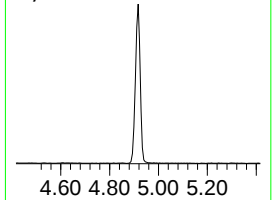
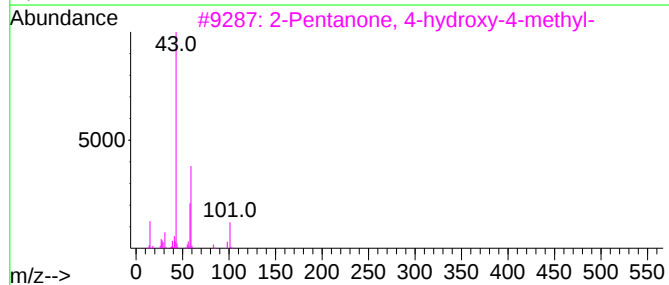
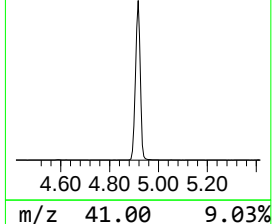
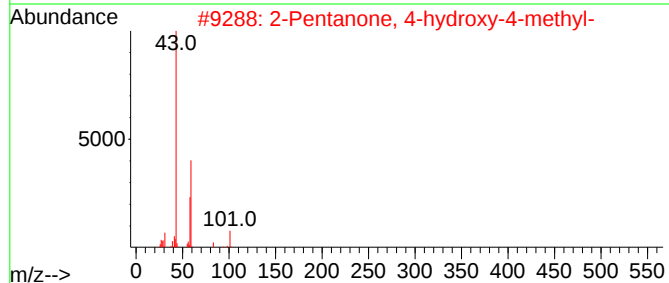
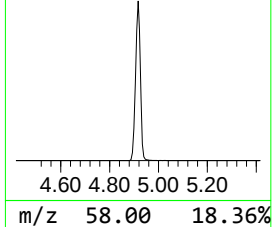
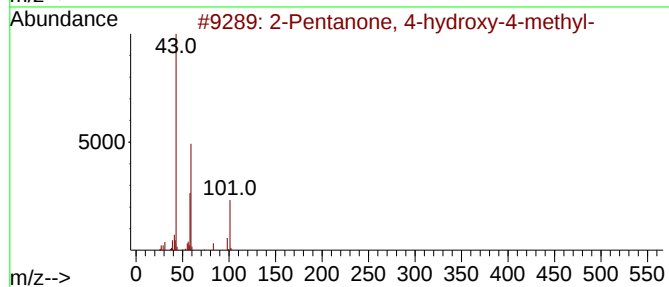
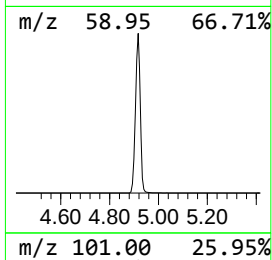
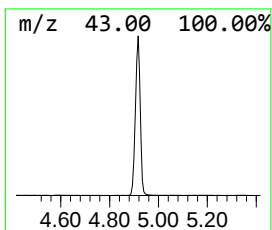
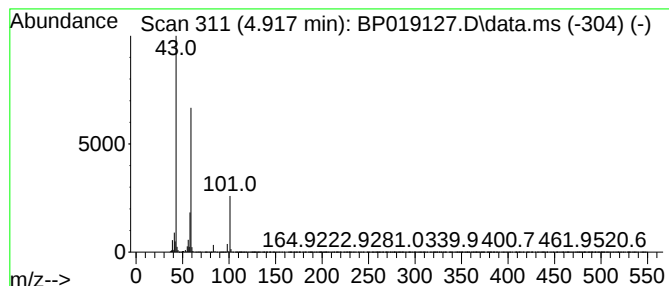
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.917	134.66 ng/ul	4504300	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019127.D
Acq On : 28 Dec 2023 16:12
Operator : MA/JU
Sample : 05920-15
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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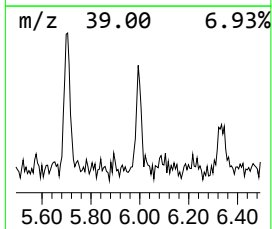
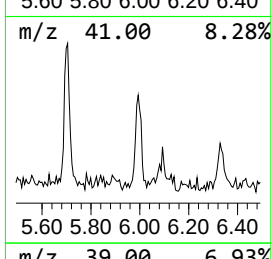
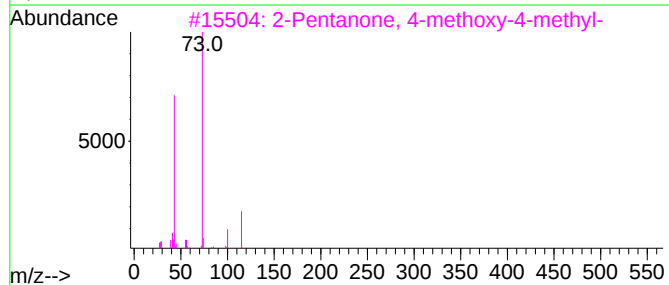
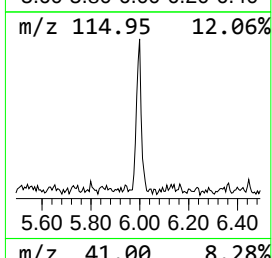
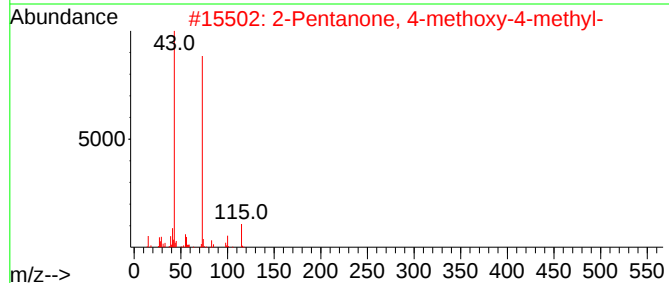
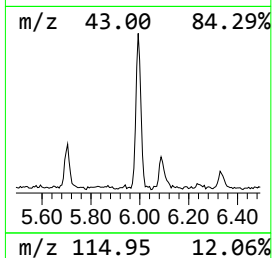
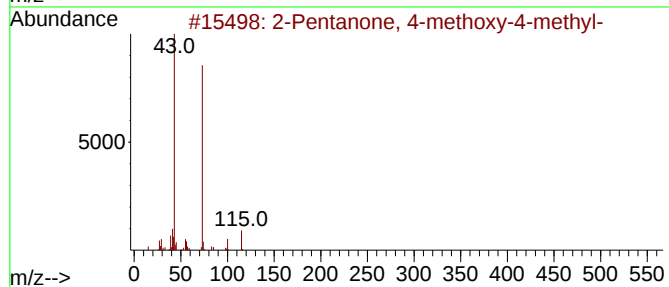
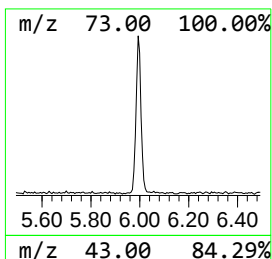
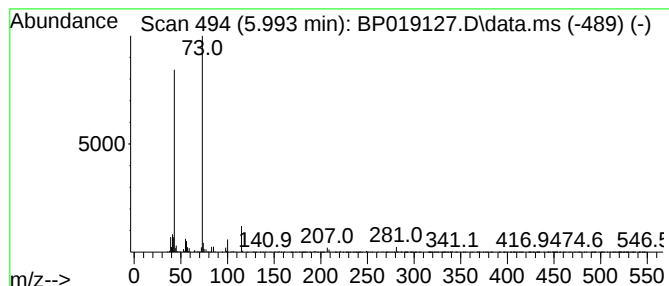
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Pentanone, 4-methoxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.993	2.29 ng/ul	76689	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	59
2			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	59
3			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	59
4			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	43
5			N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
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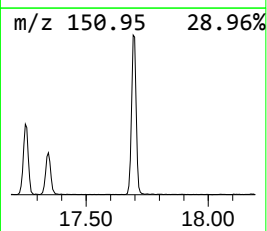
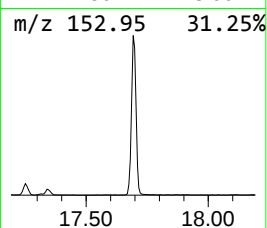
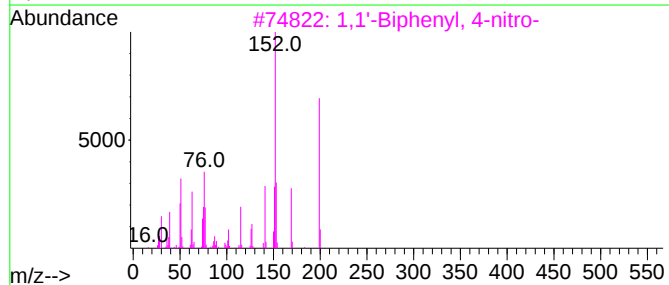
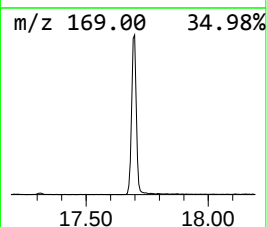
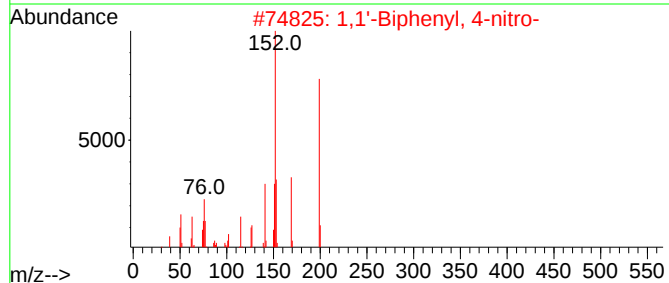
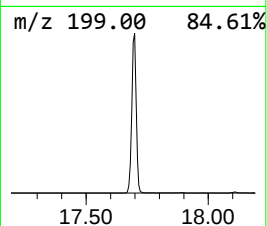
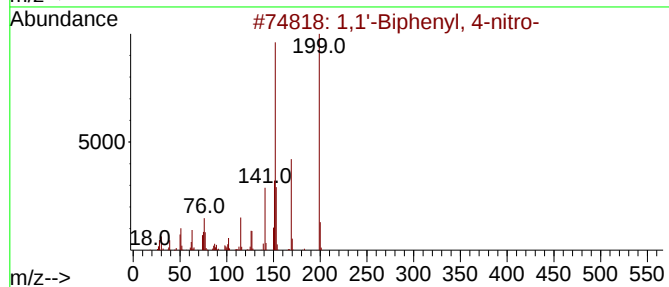
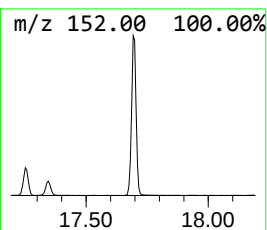
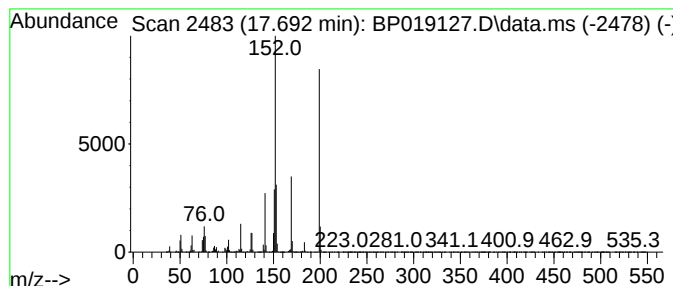
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 4-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.692	22.76 ng/ul	1682750	Phenanthrene-d10	17.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	99
2			1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
3			1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	96
4			Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	94
5			Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	91



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Misc :
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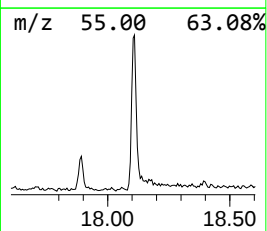
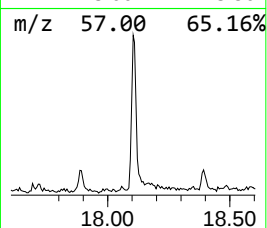
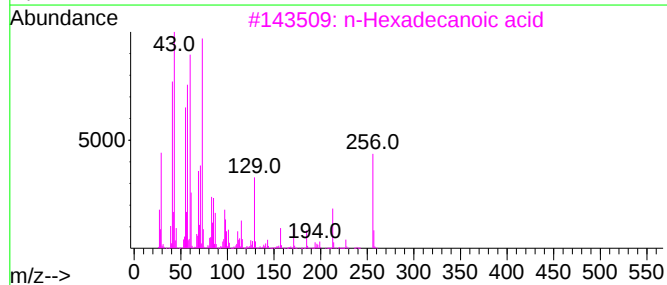
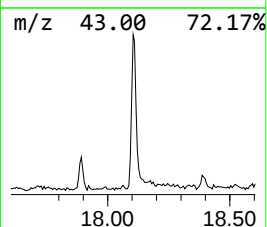
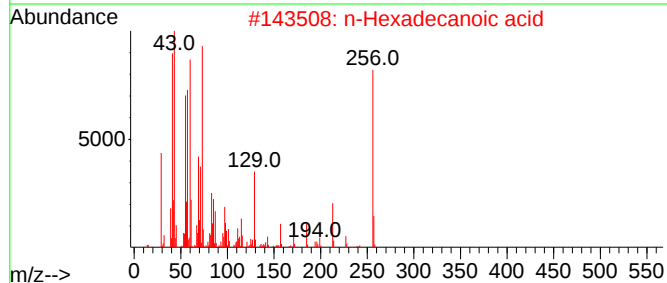
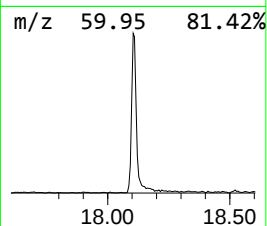
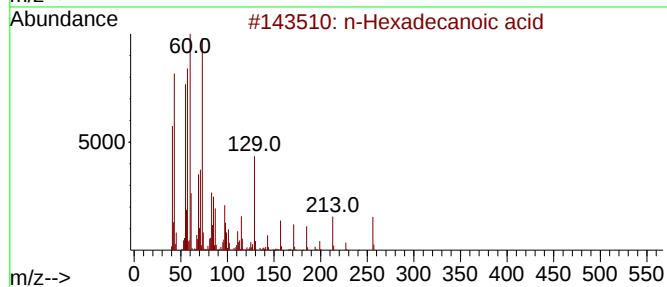
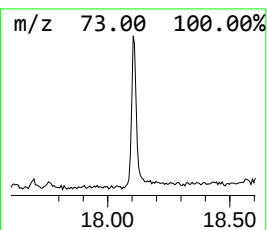
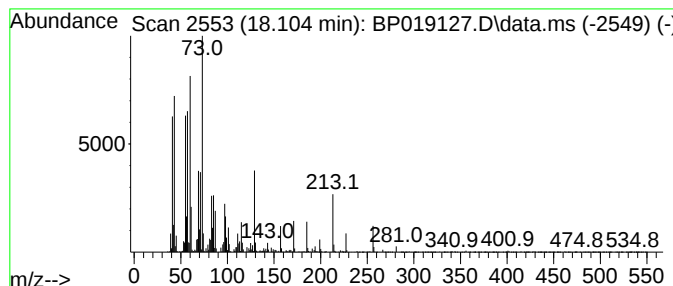
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	4.22 ng/ul	312107	Phenanthrene-d10	17.210

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019127.D
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Sample : 05920-15
Misc :
ALS Vial : 5 Sample Multiplier: 1

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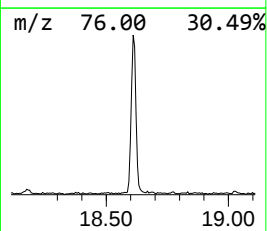
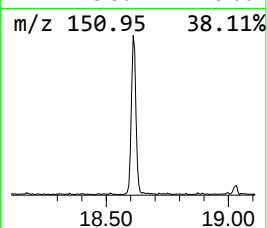
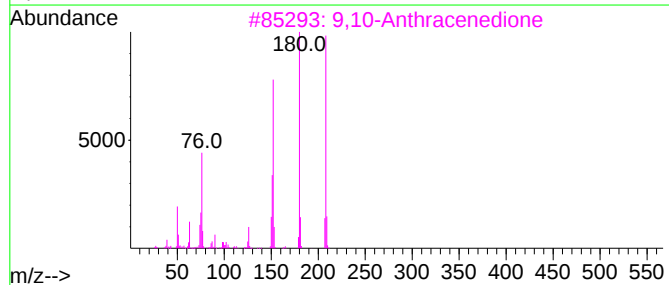
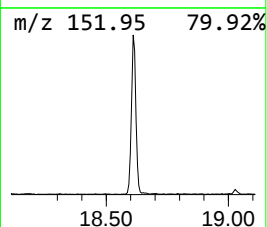
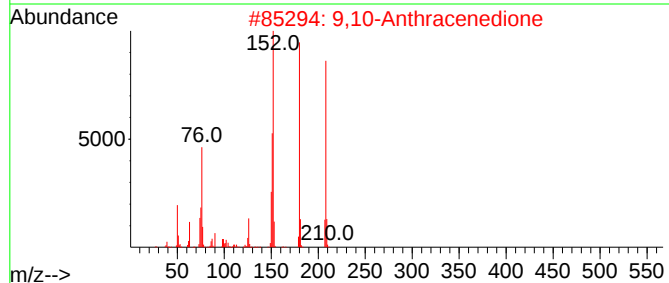
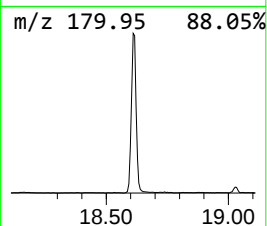
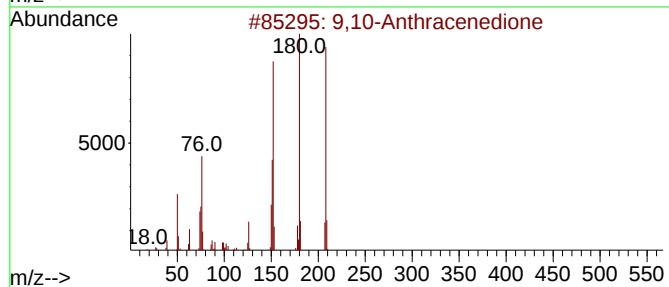
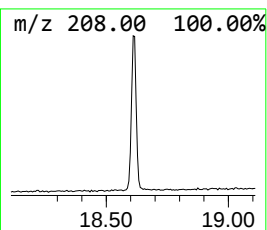
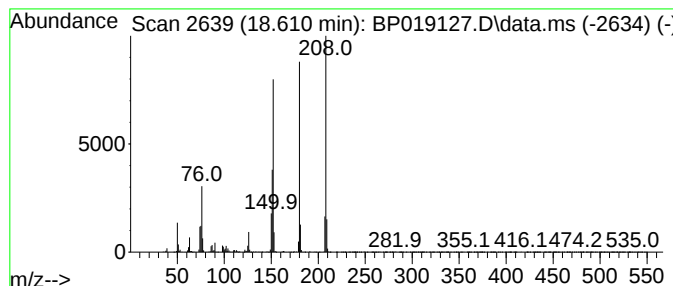
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.610	5.65 ng/ul	417833	Phenanthrene-d10	17.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
5	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
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ALS Vial : 5 Sample Multiplier: 1

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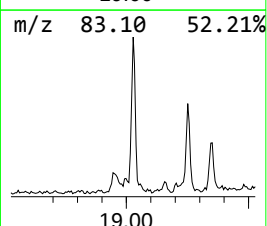
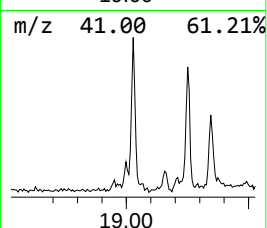
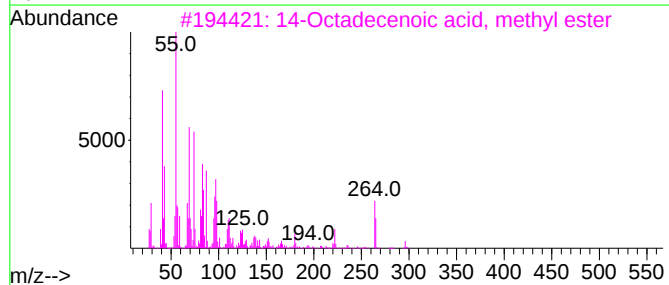
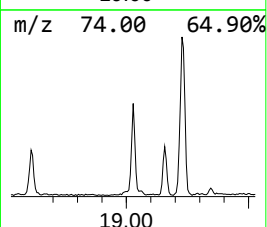
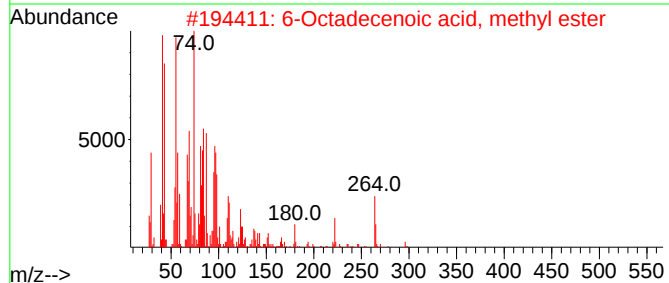
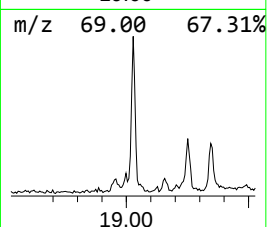
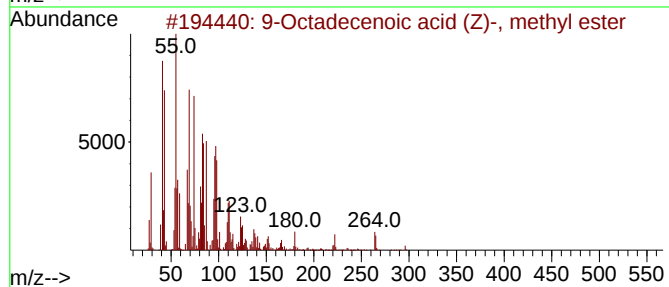
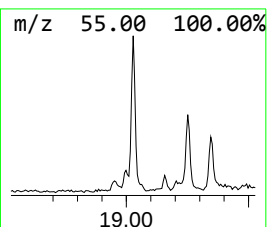
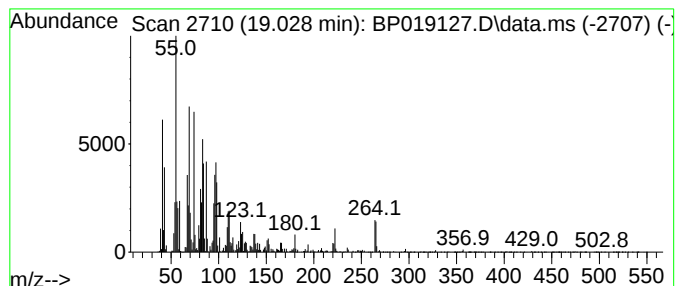
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9-Octadecenoic acid (Z)-, m... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.028	4.25 ng/ul	313906	Phenanthrene-d10	17.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2			6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
3			14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
4			12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
5			10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019127.D
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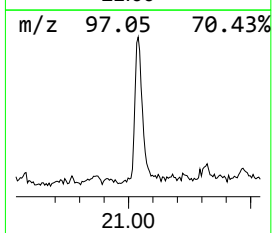
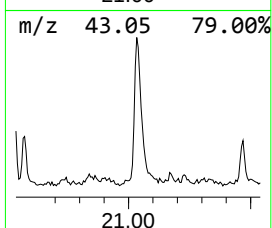
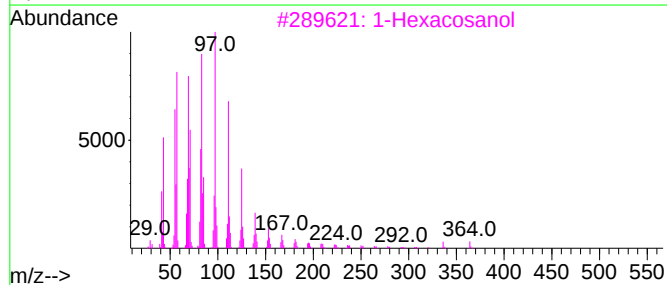
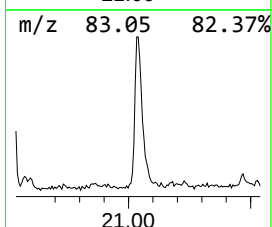
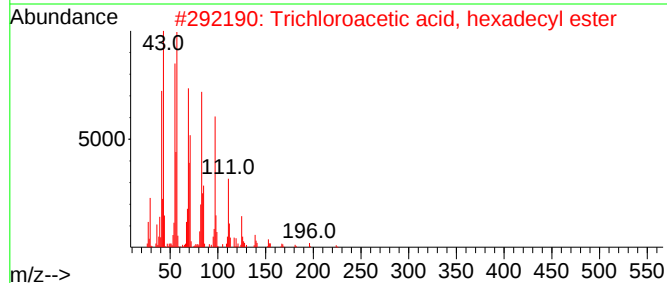
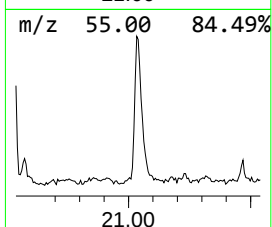
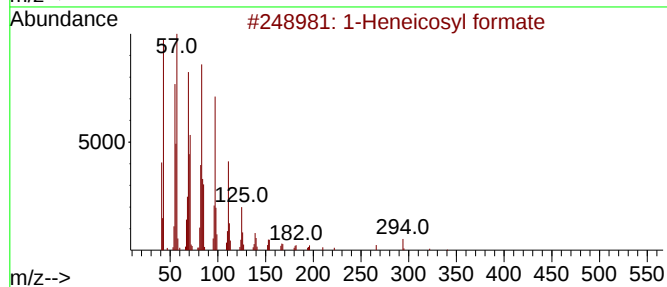
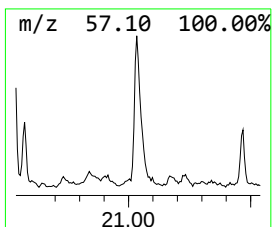
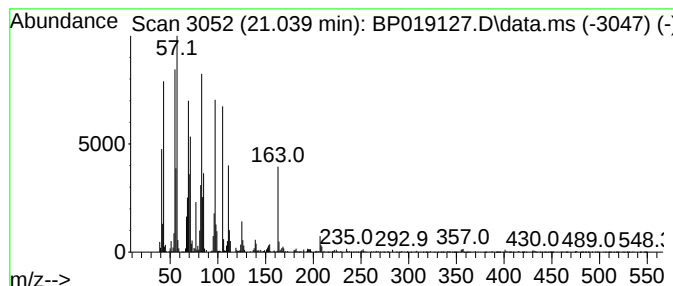
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Heneicosyl formate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.039	2.94 ng/ul	595475	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
3			1-Hexacosanol	382	C26H54O	000506-52-5	87
4			1-Hexacosene	364	C26H52	018835-33-1	87
5			Octacosanol	410	C28H58O	000557-61-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
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ALS Vial : 5 Sample Multiplier: 1

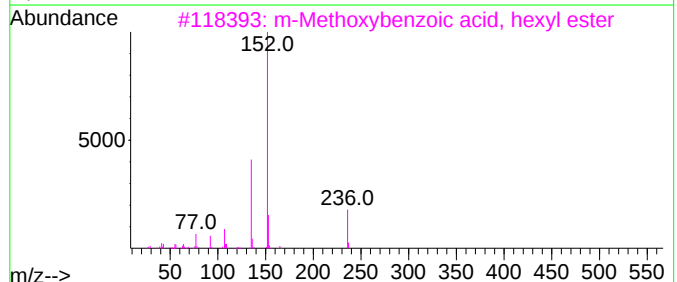
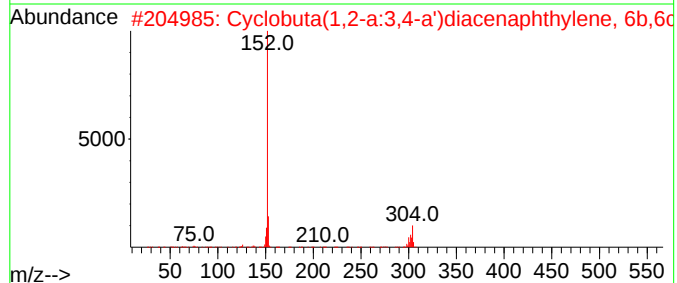
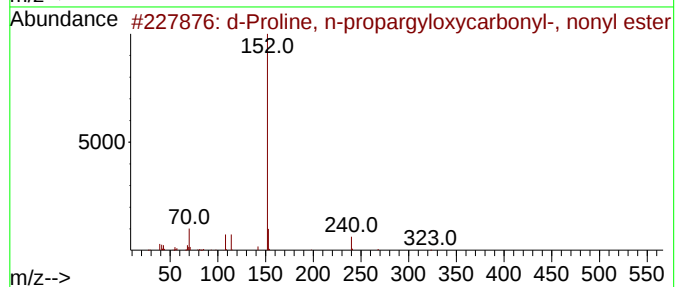
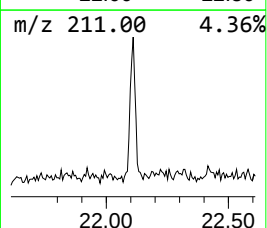
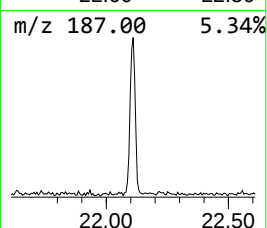
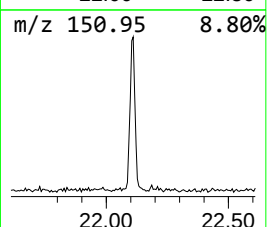
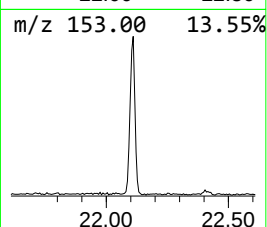
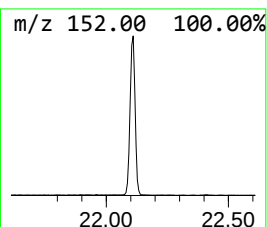
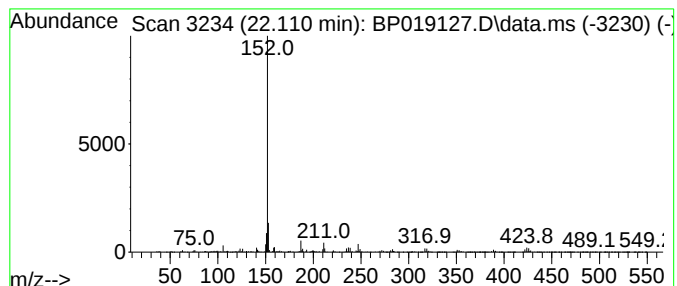
Instrument :
BNA_P
ClientSampleId :
GCF66

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 d-Proline, n-propargyloxycar... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.110	2.03 ng/ul	410722	Chrysene-d12	21.316	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	d-Proline, n-propargyloxycarbony...	323	C18H29NO4	1000328-07-1	64
2	Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	50
3	m-Methoxybenzoic acid, hexyl ester	236	C14H20O3	1000292-36-7	17
4	d-Proline, n-propargyloxycarbony...	281	C15H23NO4	1000328-07-0	9
5	Tisopurine	152	C5H4N4S	005334-23-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019127.D
Acq On : 28 Dec 2023 16:12
Operator : MA/JU
Sample : 05920-15
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF66

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.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
4-Penten-2-one,...	3.693	2.8	ng/ul	93821	1	7.799	668996	20.0
3-Penten-2-one,...	4.311	291.9	ng/ul	9762870	1	7.799	668996	20.0
2-Pentanone, 4-...	4.917	134.7	ng/ul	4504300	1	7.799	668996	20.0
2-Pentanone, 4-...	5.993	2.3	ng/ul	76689	1	7.799	668996	20.0
1,1'-Biphenyl, ...	17.692	22.8	ng/ul	1682750	4	17.210	1478940	20.0
n-Hexadecanoic ...	18.104	4.2	ng/ul	312107	4	17.210	1478940	20.0
9,10-Anthracene...	18.610	5.7	ng/ul	417833	4	17.210	1478940	20.0
9-Octadecenoic ...	19.028	4.3	ng/ul	313906	4	17.210	1478940	20.0
1-Heneicosyl fo...	21.039	2.9	ng/ul	595475	5	21.316	4045990	20.0
d-Proline, n-pr...	22.110	2.0	ng/ul	410722	5	21.316	4045990	20.0