

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
 Data File : BP020702.D
 Acq On : 28 Jun 2024 21:58
 Operator : MA/JU
 Sample : P2934-12
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CP8

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-BP062524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020702.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.011	3	4	11	rVB	3911500	3649683	100.00%	7.110%
2	3.275	45	49	55	rVB	43758	56783	1.56%	0.111%
3	3.546	91	95	104	rBV	63893	93726	2.57%	0.183%
4	3.693	117	120	131	rVV	74013	103031	2.82%	0.201%
5	3.787	132	136	141	rVB	45040	55134	1.51%	0.107%
6	3.999	169	172	179	rVB	15968	23379	0.64%	0.046%
7	4.158	195	199	204	rBV5	9403	13412	0.37%	0.026%
8	4.364	232	234	239	rVB5	8106	10626	0.29%	0.021%
9	4.499	253	257	261	rBV2	11623	15277	0.42%	0.030%
10	4.546	261	265	270	rVV	34061	51178	1.40%	0.100%
11	4.599	270	274	284	rVB	24751	35805	0.98%	0.070%
12	4.793	302	307	314	rVB8	7002	15575	0.43%	0.030%
13	4.899	319	325	334	rVB	237902	350590	9.61%	0.683%
14	5.005	339	343	349	rVB5	9637	16554	0.45%	0.032%
15	5.858	482	488	496	rVV	35675	62663	1.72%	0.122%
16	6.775	638	644	652	rVB8	6859	14938	0.41%	0.029%
17	6.928	663	670	683	rVB	695190	1073676	29.42%	2.092%
18	7.087	690	697	706	rBV	553984	866939	23.75%	1.689%
19	7.287	724	731	739	rBV	734545	1150903	31.53%	2.242%
20	7.358	739	743	753	rVB2	21166	44192	1.21%	0.086%
21	7.740	799	808	815	rBV	755906	1212229	33.21%	2.361%
22	7.810	815	820	828	rVB3	24294	42302	1.16%	0.082%
23	8.446	920	928	941	rBV	760846	1196150	32.77%	2.330%
24	8.557	941	947	954	rVB2	20304	40605	1.11%	0.079%
25	8.893	996	1004	1012	rBV	617414	1073532	29.41%	2.091%
26	9.275	1064	1069	1076	rVB7	8036	14854	0.41%	0.029%
27	9.446	1088	1098	1112	rVB2	101153	168705	4.62%	0.329%
28	9.599	1115	1124	1138	rBV	520813	910911	24.96%	1.775%
29	9.846	1161	1166	1171	rVB9	5663	12836	0.35%	0.025%
30	10.146	1209	1217	1230	rBV	755467	1348660	36.95%	2.627%
31	10.516	1263	1280	1292	rBV	1013308	1640942	44.96%	3.197%
32	10.651	1296	1303	1320	rBV	472813	842048	23.07%	1.640%
33	11.051	1364	1371	1377	rBV3	34139	70001	1.92%	0.136%
34	11.257	1399	1406	1424	rVV	139184	270629	7.42%	0.527%
35	12.093	1543	1548	1554	rVV2	13842	26272	0.72%	0.051%
36	12.398	1596	1600	1607	rVB6	5915	10221	0.28%	0.020%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	12.716	1648	1654	1657	rVV8	5577	9158	0.25%	0.018%
38	13.181	1729	1733	1740	rVB9	6416	13753	0.38%	0.027%
39	13.687	1814	1819	1826	rBV9	6511	14429	0.40%	0.028%
40	13.787	1826	1836	1845	rVV	1292754	1873545	51.33%	3.650%
41	14.022	1870	1876	1878	rBV6	13643	24144	0.66%	0.047%
42	14.069	1878	1884	1901	rVV	1506183	2294306	62.86%	4.469%
43	14.198	1901	1906	1911	rVV9	5541	12485	0.34%	0.024%
44	14.381	1927	1937	1944	rVV	1354910	1987383	54.45%	3.872%
45	14.440	1944	1947	1955	rVB	30650	44540	1.22%	0.087%
46	14.510	1955	1959	1962	rBV6	7506	11888	0.33%	0.023%
47	14.598	1967	1974	1994	rBV	439681	836258	22.91%	1.629%
48	14.745	1995	1999	2001	rVV4	17807	26542	0.73%	0.052%
49	14.804	2003	2009	2015	rVB5	24390	52286	1.43%	0.102%
50	15.051	2049	2051	2057	rVB7	7579	10467	0.29%	0.020%
51	15.187	2068	2074	2078	rBV7	11994	25899	0.71%	0.050%
52	15.239	2080	2083	2087	rVB6	10273	16600	0.45%	0.032%
53	15.387	2101	2108	2114	rBV	1778961	2665785	73.04%	5.193%
54	15.445	2114	2118	2124	rVB3	43400	65731	1.80%	0.128%
55	15.516	2124	2130	2137	rBV	417651	612924	16.79%	1.194%
56	15.651	2149	2153	2159	rVB3	34956	58544	1.60%	0.114%
57	15.745	2164	2169	2172	rBV4	9634	15500	0.42%	0.030%
58	15.904	2192	2196	2201	rVB4	9264	15135	0.41%	0.029%
59	15.969	2203	2207	2210	rBV6	8915	15542	0.43%	0.030%
60	16.134	2229	2235	2241	rVB6	12711	31422	0.86%	0.061%
61	16.281	2256	2260	2263	rBV6	7117	12026	0.33%	0.023%
62	16.398	2274	2280	2284	rBV2	28062	38724	1.06%	0.075%
63	16.475	2290	2293	2297	rVB5	12225	16005	0.44%	0.031%
64	16.581	2307	2311	2315	rBV7	12533	21907	0.60%	0.043%
65	16.834	2350	2354	2360	rVB3	15899	23886	0.65%	0.047%
66	16.910	2362	2367	2368	rBV5	9784	16643	0.46%	0.032%
67	16.998	2378	2382	2389	rVB	28052	46974	1.29%	0.092%
68	17.098	2395	2399	2405	rVB6	16515	25948	0.71%	0.051%
69	17.186	2407	2414	2418	rBV	1433465	2152054	58.97%	4.192%
70	17.228	2418	2421	2425	rVV	456708	587667	16.10%	1.145%
71	17.286	2425	2431	2442	rVB2	1755672	2793060	76.53%	5.441%
72	17.375	2442	2446	2448	rBV3	19468	26553	0.73%	0.052%
73	17.569	2477	2479	2482	rBV3	7830	10495	0.29%	0.020%
74	17.610	2483	2486	2491	rVV2	25678	36220	0.99%	0.071%
75	17.792	2513	2517	2525	rVB6	25824	61697	1.69%	0.120%
76	17.998	2547	2552	2554	rBV6	12066	21513	0.59%	0.042%

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77	18.098	2565	2569	2570	rBV	60991	68862	1.89%	0.134%
78	18.128	2570	2574	2584	rVV2	362832	656991	18.00%	1.280%
79	18.222	2587	2590	2595	rVB	22381	32627	0.89%	0.064%
80	18.286	2597	2601	2607	rVV2	100757	171229	4.69%	0.334%
81	18.463	2628	2631	2632	rBV2	24729	22401	0.61%	0.044%
82	18.616	2653	2657	2661	rVB	43275	58496	1.60%	0.114%
83	18.710	2670	2673	2677	rVB5	32967	37465	1.03%	0.073%
84	19.322	2772	2777	2783	rVB	848585	1135176	31.10%	2.211%
85	19.457	2796	2800	2809	rBV3	147345	246726	6.76%	0.481%
86	19.669	2831	2836	2840	rBV	2005479	3087305	84.59%	6.014%
87	20.216	2926	2929	2934	rVB	95689	144959	3.97%	0.282%
88	20.263	2935	2937	2940	rVB2	81489	76828	2.11%	0.150%
89	20.327	2944	2948	2952	rBV2	95080	128397	3.52%	0.250%
90	20.839	3033	3035	3041	rVB5	121546	134866	3.70%	0.263%
91	21.163	3086	3090	3094	rBV2	303727	427800	11.72%	0.833%
92	21.210	3094	3098	3102	rVB	265257	368550	10.10%	0.718%
93	21.316	3112	3116	3123	rVB	487030	673479	18.45%	1.312%
94	21.480	3138	3144	3149	rBV2	923770	1542258	42.26%	3.004%
95	21.533	3149	3153	3161	rVV2	480282	926065	25.37%	1.804%
96	21.645	3165	3172	3177	rVV3	1366998	2947447	80.76%	5.742%
97	21.692	3177	3180	3185	rVB	283764	436575	11.96%	0.850%
98	22.545	3322	3325	3335	rBV5	144033	337056	9.24%	0.657%
99	24.774	3697	3704	3712	rVB2	711722	1961759	53.75%	3.822%
100	25.010	3737	3744	3758	rVB	868778	2499249	68.48%	4.869%

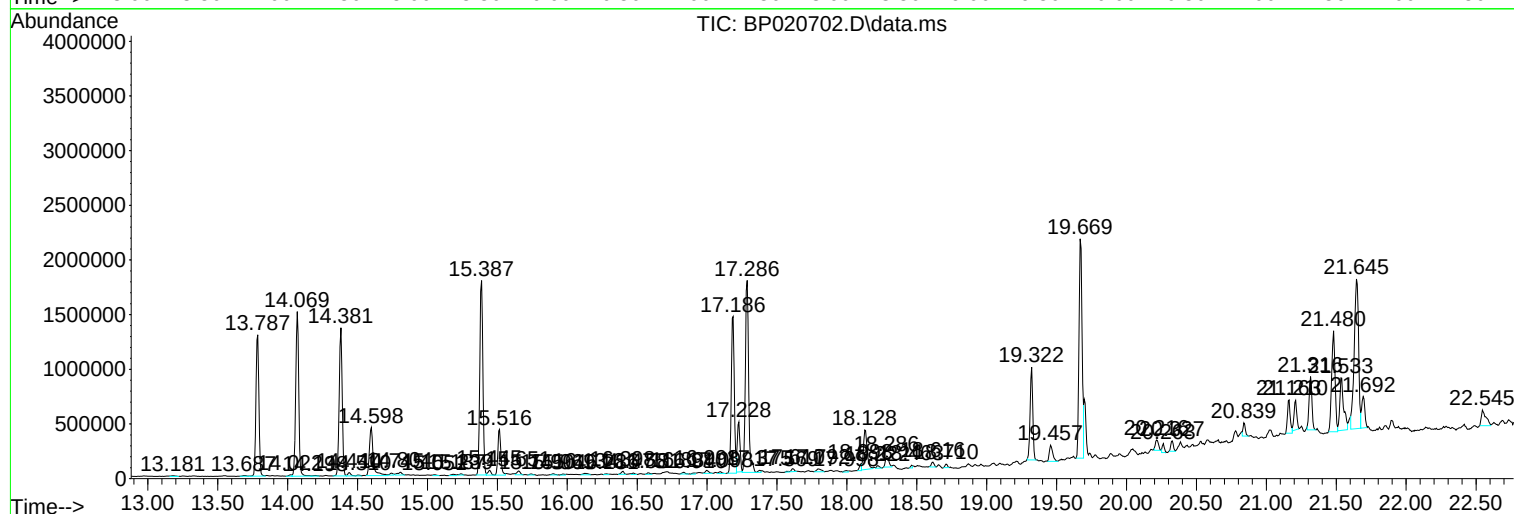
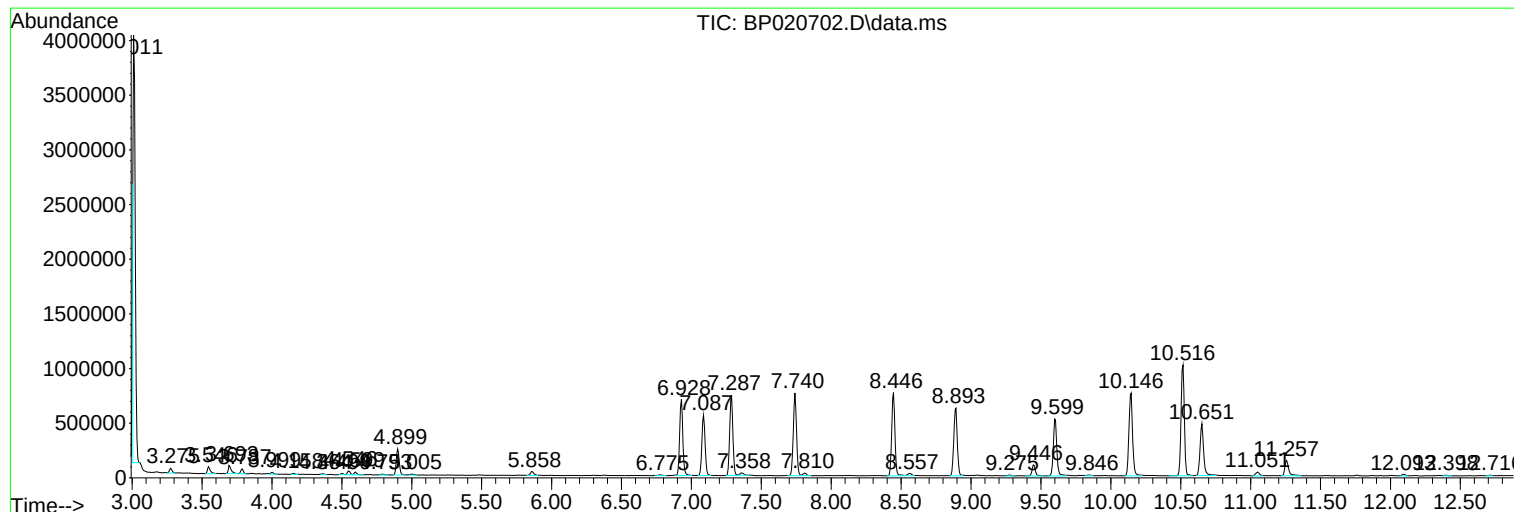
Sum of corrected areas: 51333165

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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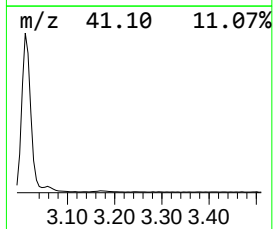
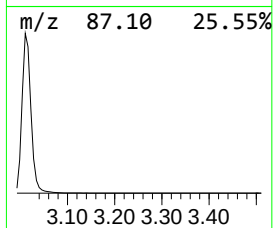
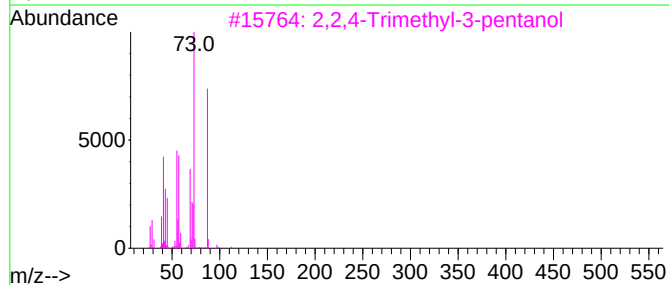
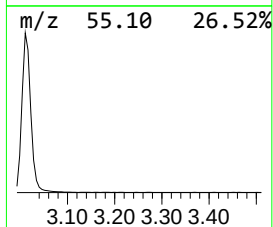
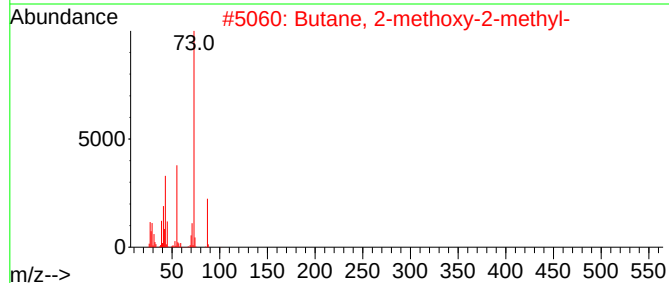
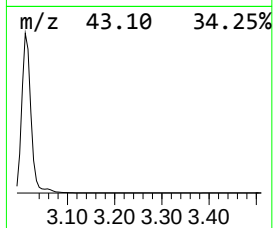
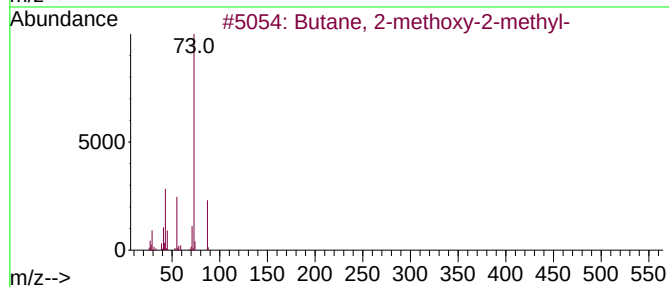
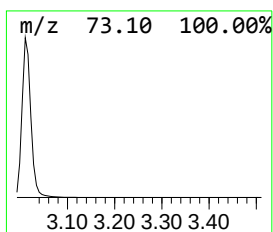
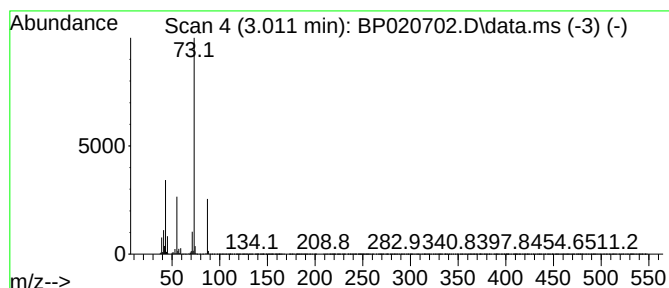
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.011	60.21 ng/ul	3649680	1,4-Dichlorobenzene-d4	7.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	45		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	10		



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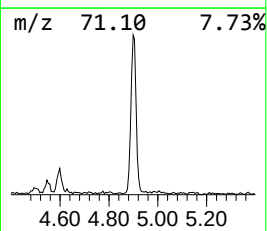
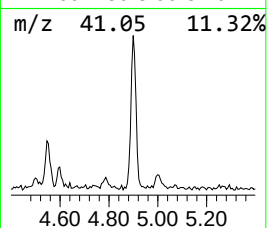
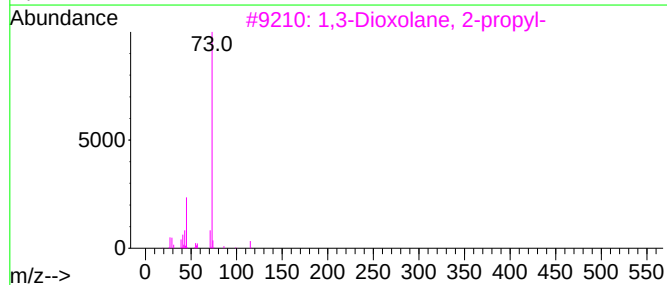
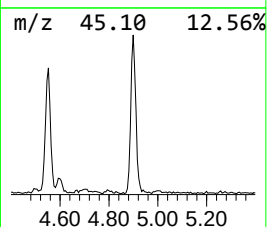
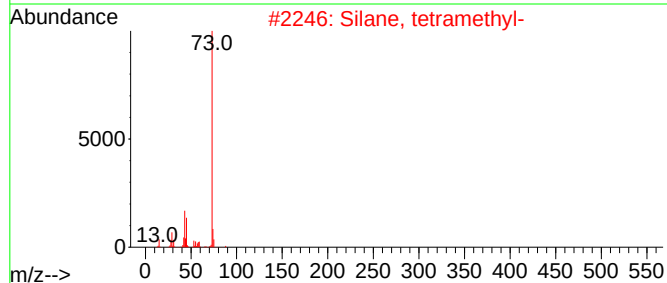
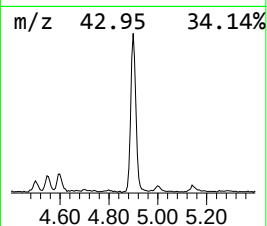
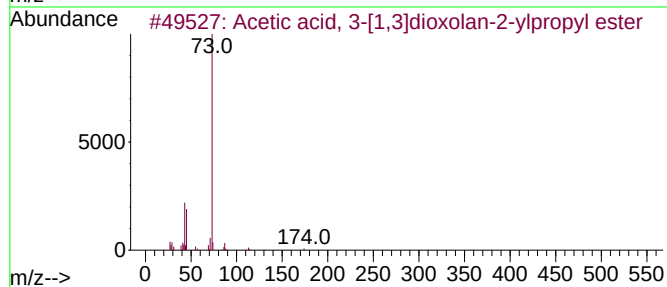
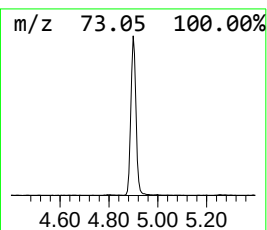
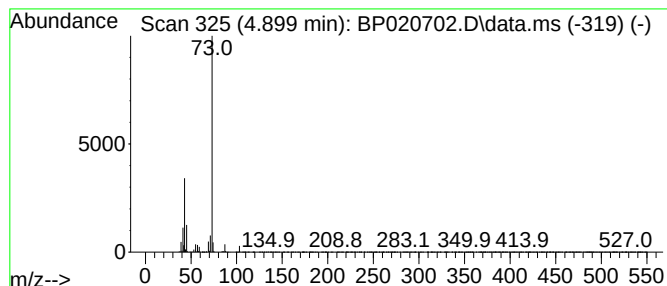
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.899	5.78 ng/ul	350590	1,4-Dichlorobenzene-d4	7.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	42
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	28
4			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	28
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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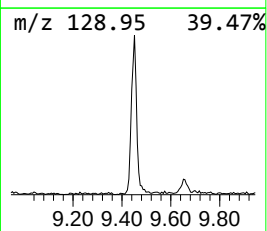
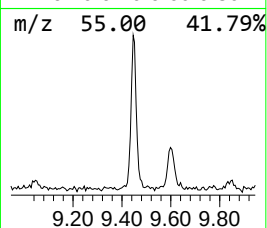
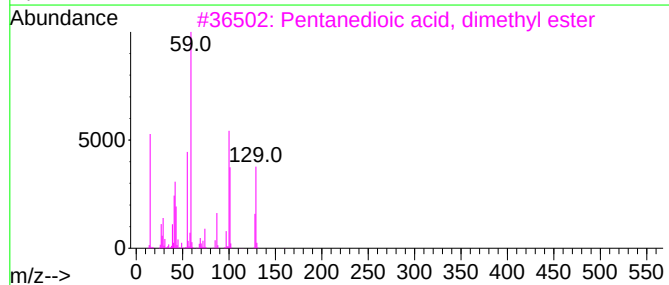
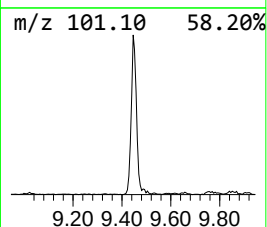
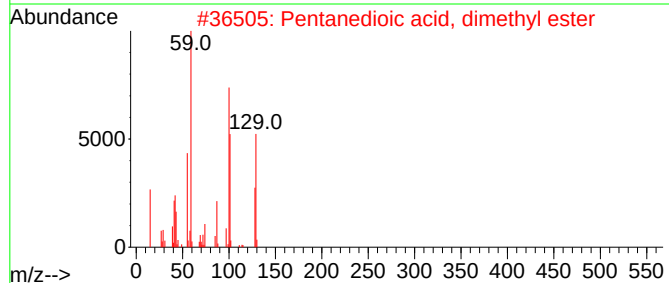
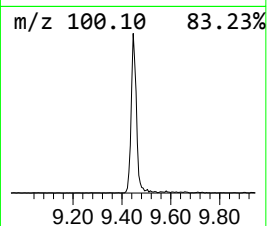
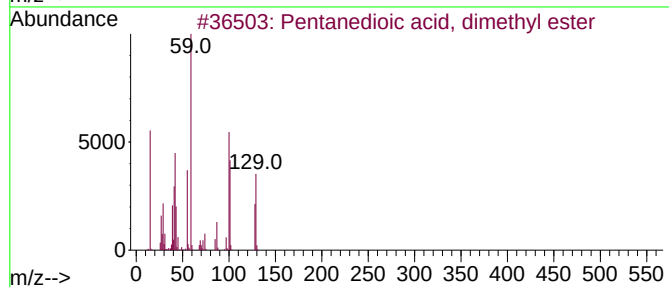
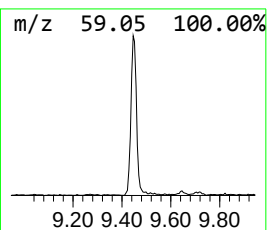
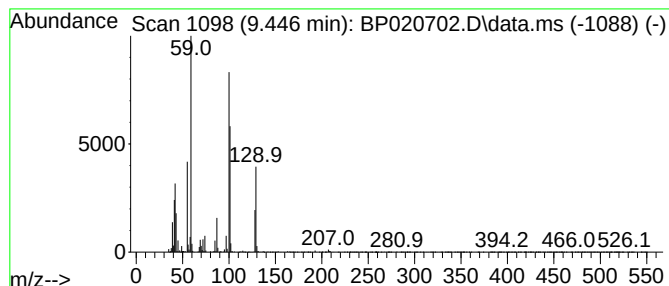
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanedioic acid, dimethyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.446	2.06 ng/ul	168705	Naphthalene-d8	10.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	83
2			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	64
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	59
4			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	50
5			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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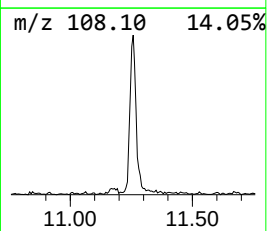
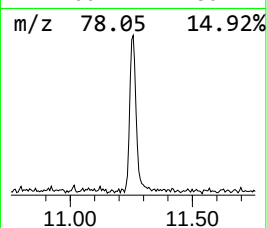
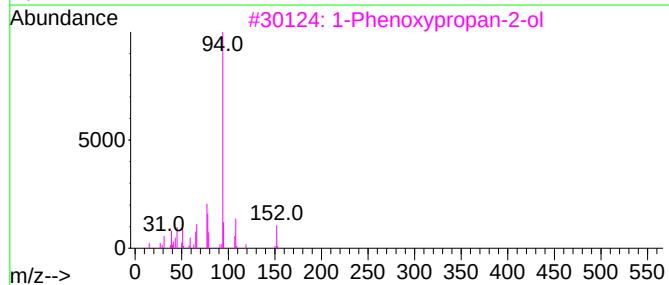
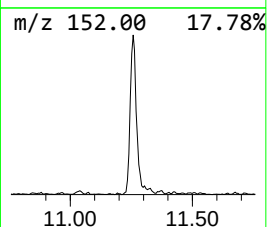
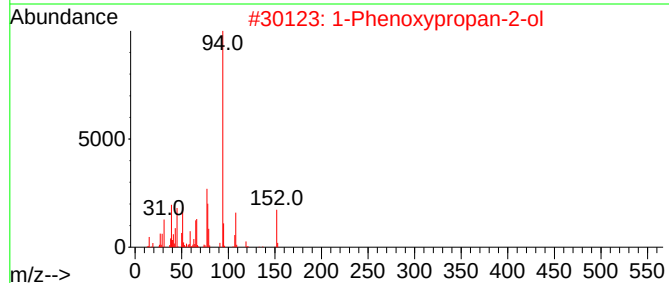
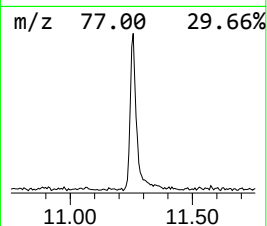
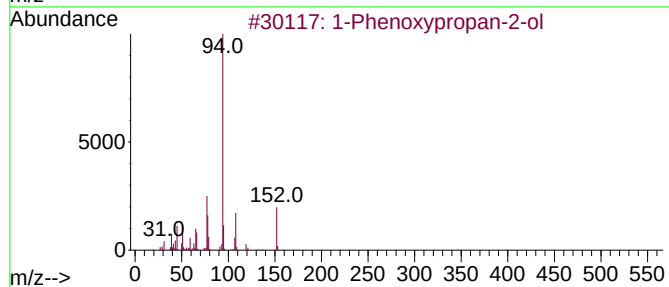
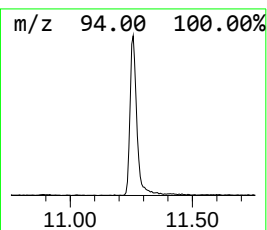
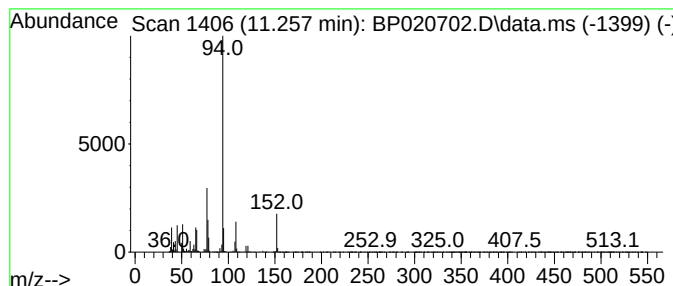
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.257	3.30 ng/ul	270629	Naphthalene-d8	10.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
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Instrument :
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ClientSampleId :
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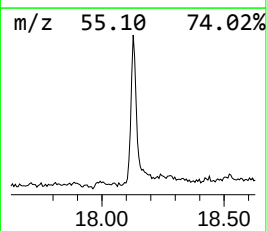
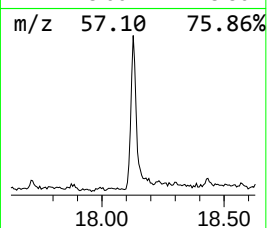
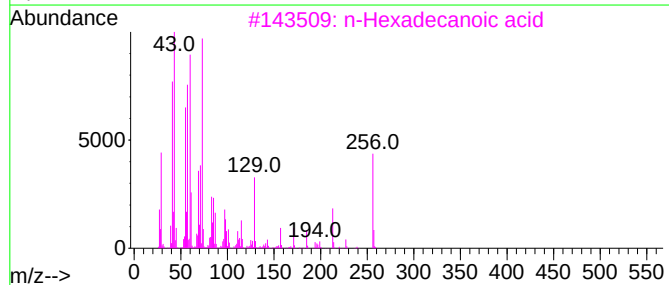
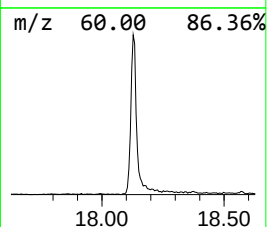
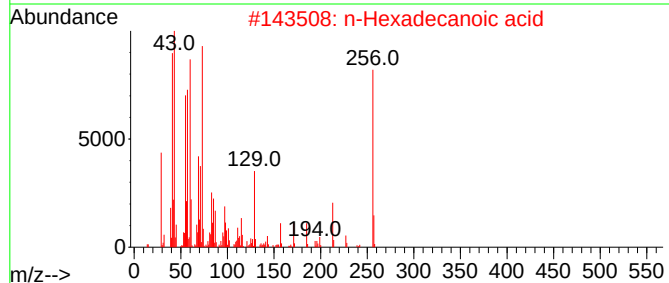
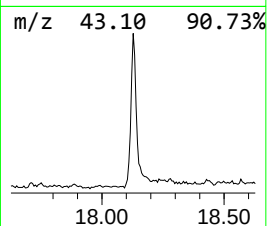
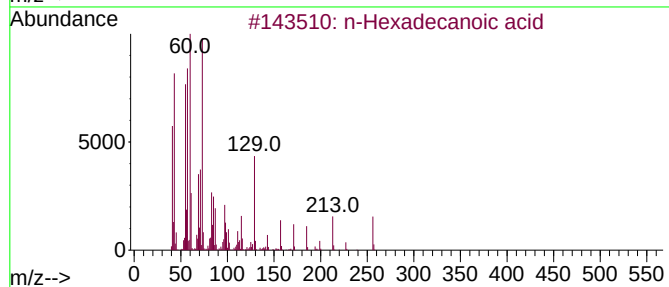
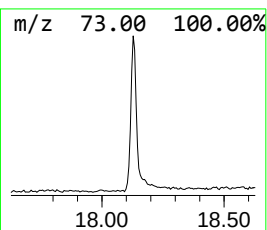
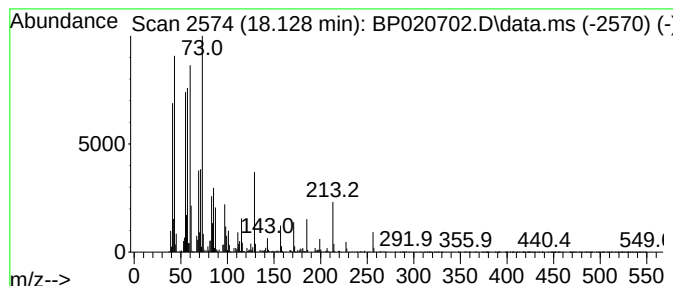
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	6.11 ng/ul	656991	Phenanthrene-d10	17.186

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	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
ALS Vial : 60 Sample Multiplier: 1

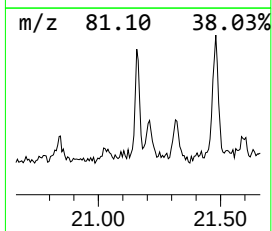
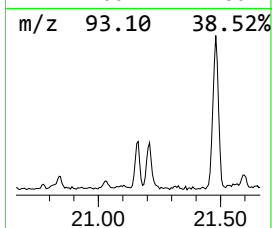
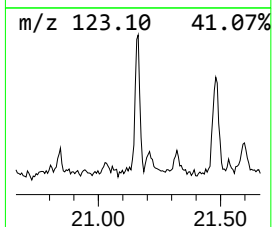
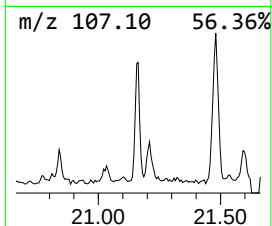
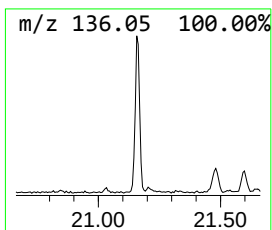
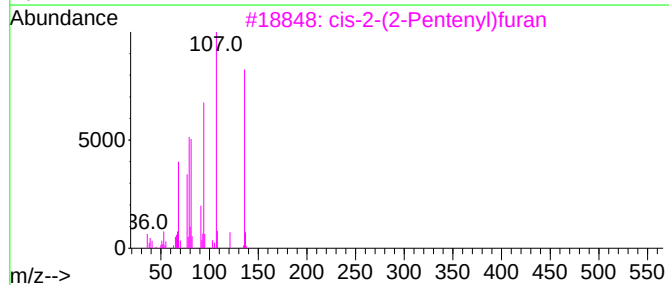
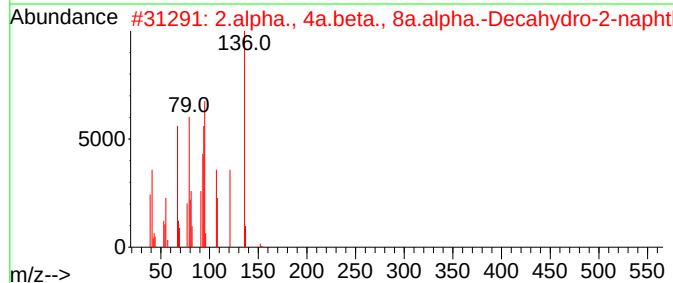
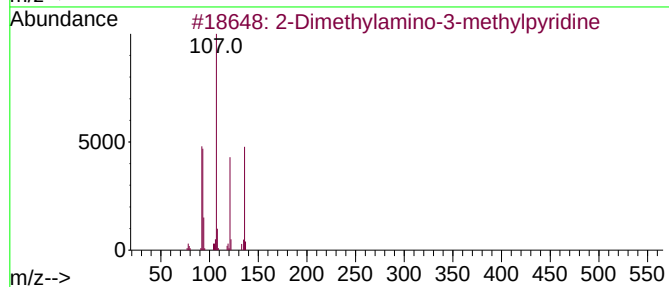
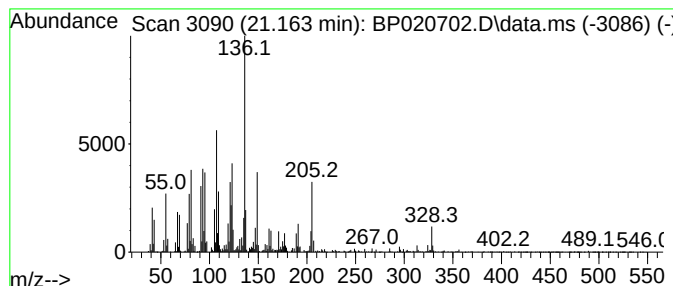
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.163	2.90 ng/ul	427800	Chrysene-d12	21.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dimethylamino-3-methylpyridine	136	C8H12N2	061713-46-0	46
2	2.alpha., 4a.beta., 8a.alpha.-De...	154	C10H18O	005779-35-1	38
3	cis-2-(2-Pentenyl)furan	136	C9H12O	070424-13-4	35
4	2-(5-Fluoropyridin-2-yl)hexanen...	192	C11H13FN2	1000482-12-7	27
5	Epiquinidine	324	C20H24N2O2	000572-59-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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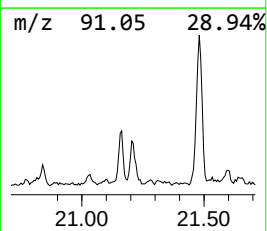
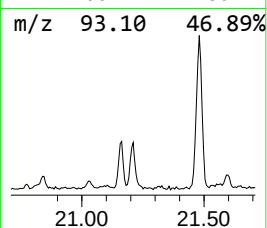
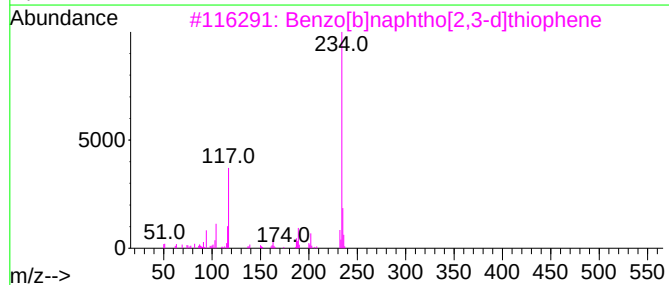
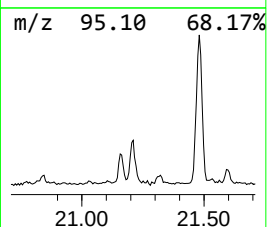
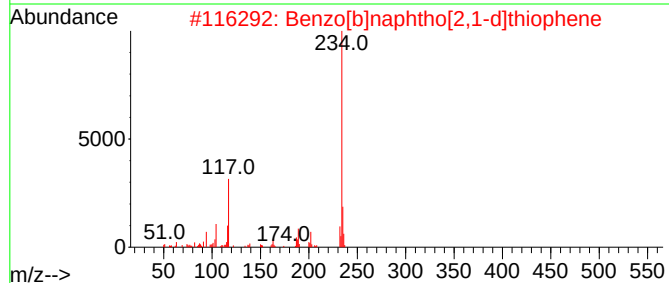
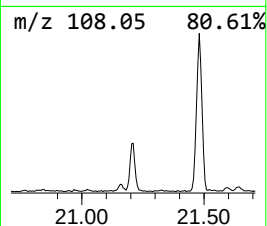
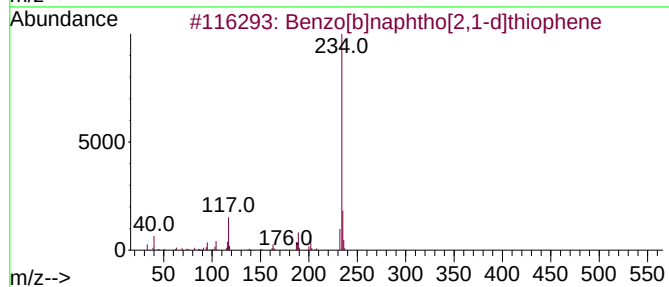
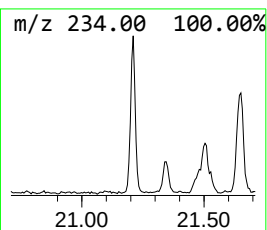
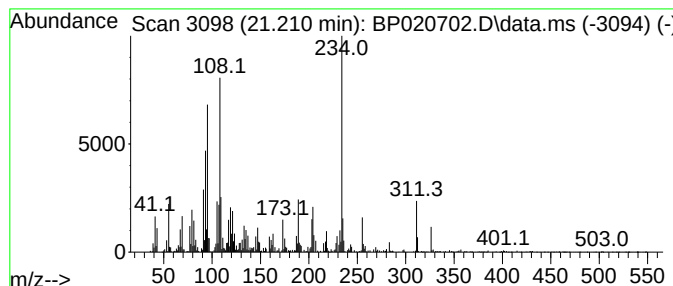
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.210	2.50 ng/ul	368550	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	38
2			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	35
3			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	30
4			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	30
5			4-Thiazoline, 4-(benzo[b]1,4-dio...	234	C11H10N2O2S	105362-06-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
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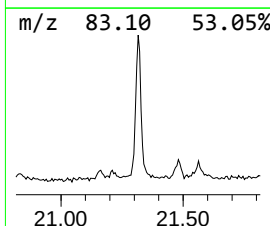
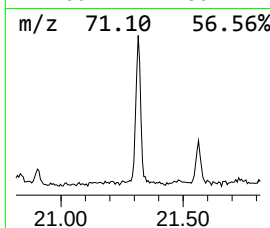
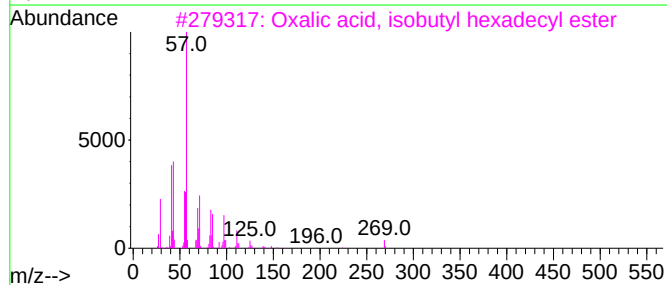
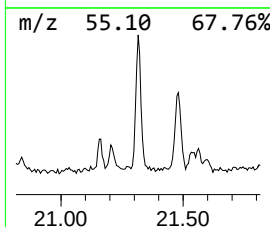
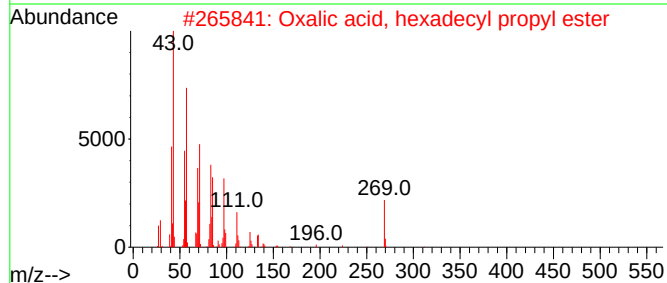
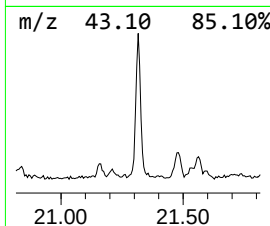
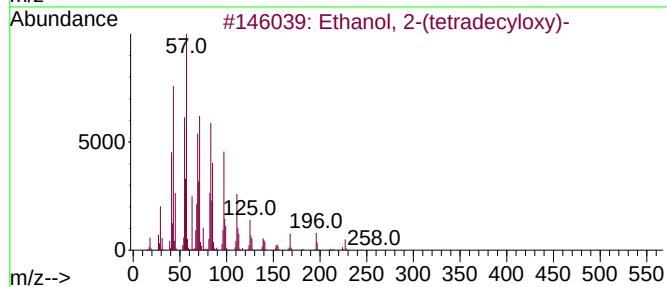
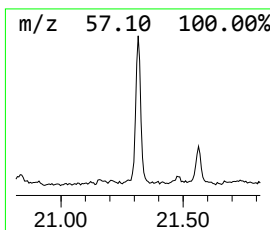
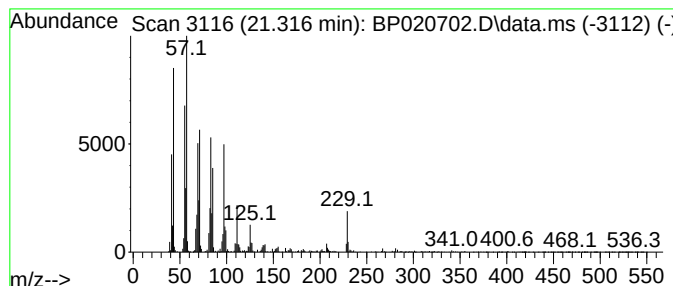
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Ethanol, 2-(tetradecyloxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.316	4.57 ng/ul	673479	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
2			Oxalic acid, hexadecyl propyl ester	356	C21H40O4	1000309-26-9	90
3			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	87
4			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	87
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
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Misc :
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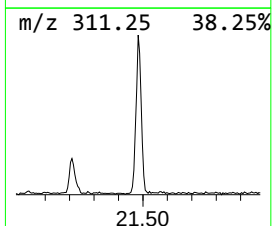
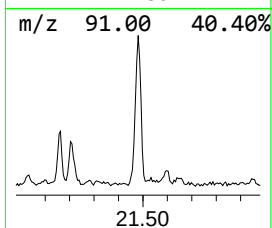
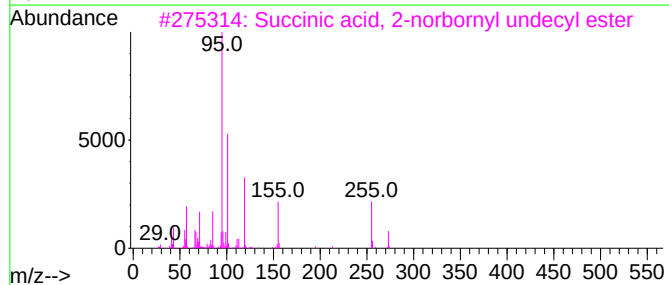
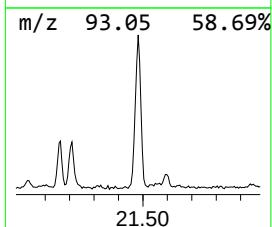
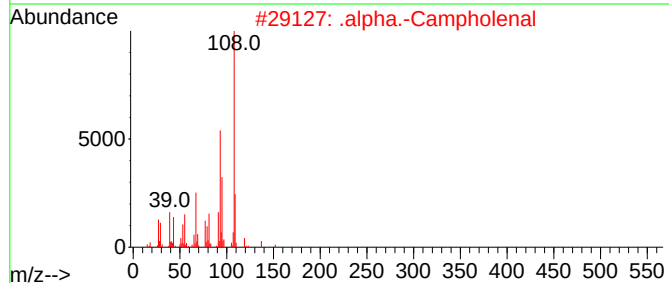
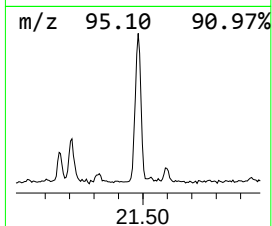
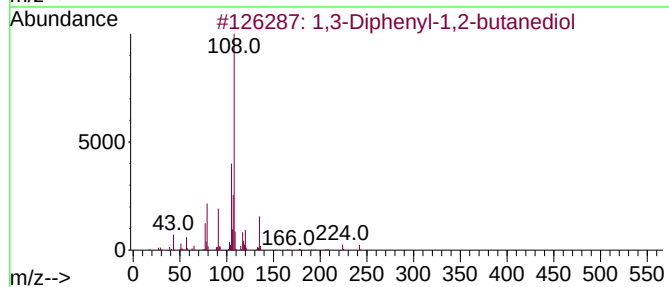
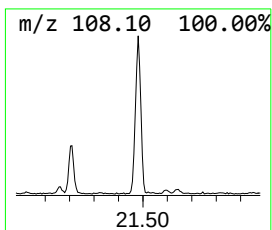
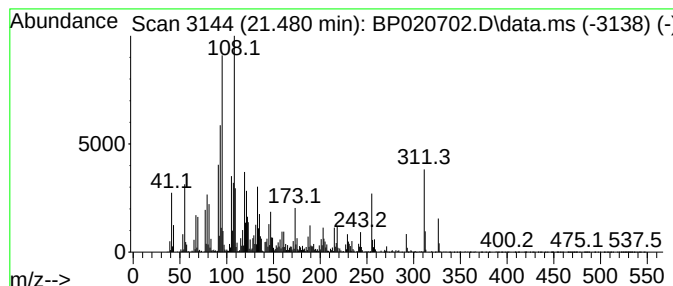
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.480	10.47 ng/ul	1542260	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Diphenyl-1,2-butanediol	242	C16H18O2	005381-88-4	25
2			.alpha.-Campholenal	152	C10H16O	004501-58-0	22
3			Succinic acid, 2-norbornyl undec...	366	C22H38O4	1000330-19-9	22
4			1,4,6-Trimethyl-1,2,3,3a,4,7,8,8...	204	C15H24	065128-08-7	20
5			1,3-Cyclohexanebis(methylamine),...	534	C16H16F14N2O2	1000454-07-2	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
ALS Vial : 60 Sample Multiplier: 1

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ClientSampleId :
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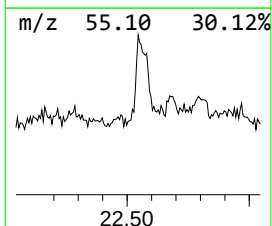
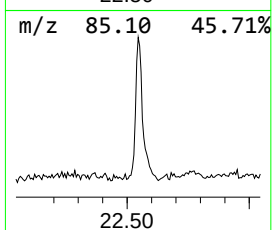
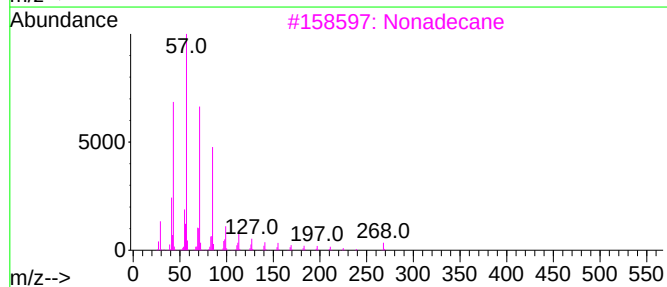
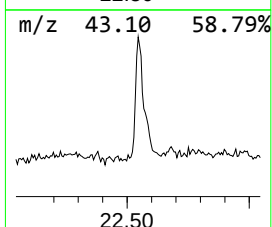
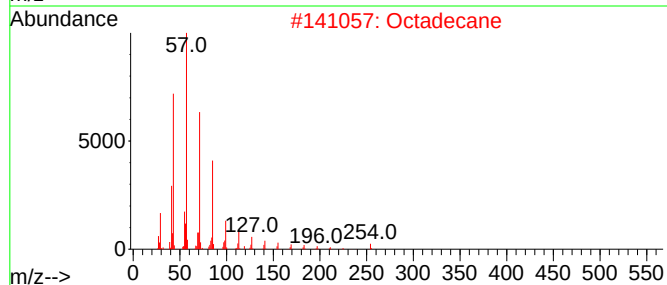
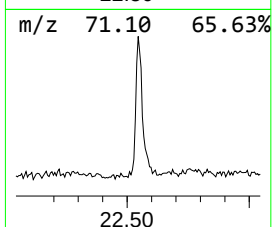
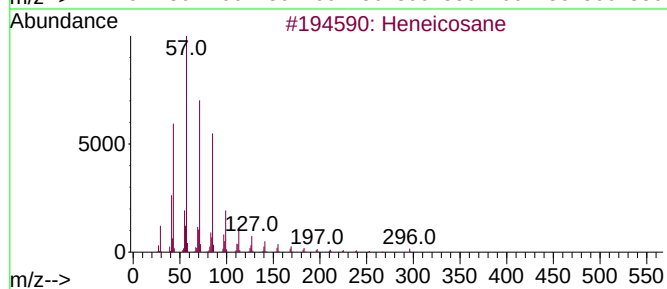
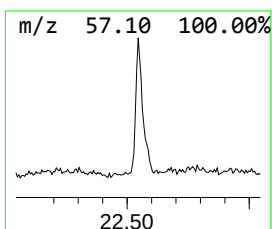
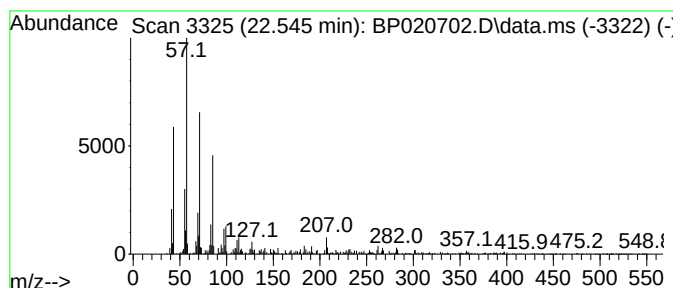
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.545	2.29 ng/ul	337056	Chrysene-d12	21.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	93
2			Octadecane	254	C18H38	000593-45-3	93
3			Nonadecane	268	C19H40	000629-92-5	91
4			Eicosane	282	C20H42	000112-95-8	91
5			Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062724\
Data File : BP020702.D
Acq On : 28 Jun 2024 21:58
Operator : MA/JU
Sample : P2934-12
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CP8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.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
Butane, 2-metho...	3.011	60.2	ng/ul	3649680	1	7.740	1212230	20.0
unknown-01	4.899	5.8	ng/ul	350590	1	7.740	1212230	20.0
Pentanedioic ac...	9.446	2.1	ng/ul	168705	2	10.516	1640940	20.0
1-Phenoxypropan...	11.257	3.3	ng/ul	270629	2	10.516	1640940	20.0
n-Hexadecanoic ...	18.128	6.1	ng/ul	656991	4	17.186	2152050	20.0
unknown-02	21.163	2.9	ng/ul	427800	5	21.651	2947450	20.0
unknown-03	21.210	2.5	ng/ul	368550	5	21.651	2947450	20.0
Ethanol, 2-(tet...	21.316	4.6	ng/ul	673479	5	21.651	2947450	20.0
unknown-04	21.480	10.5	ng/ul	1542260	5	21.651	2947450	20.0
(DEL) Alkane: S...	22.545	2.3	ng/ul	337056	5	21.651	2947450	20.0