

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
 Data File : BN030642.D
 Acq On : 04 Apr 2024 00:42
 Operator : MA/JU
 Sample : P1849-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EOLY8

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_N\Methods\SFAM-EPA-BN032824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN030642.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.199	32	36	44	rVB2	38558	66278	1.63%	0.106%
2	3.605	102	105	118	rBV2	149790	259655	6.38%	0.415%
3	3.711	118	123	130	rVV	102333	137026	3.37%	0.219%
4	3.781	132	135	141	rVB3	14452	22623	0.56%	0.036%
5	4.469	249	252	258	rVV2	17505	26009	0.64%	0.042%
6	4.828	308	313	319	rVB2	16222	35883	0.88%	0.057%
7	4.917	324	328	334	rVV	24405	35854	0.88%	0.057%
8	5.058	347	352	362	rVB	21672	44185	1.09%	0.071%
9	5.475	418	423	430	rVV	94710	152905	3.76%	0.244%
10	5.699	456	461	468	rBV4	17947	35518	0.87%	0.057%
11	6.146	532	537	543	rBV3	27397	43725	1.08%	0.070%
12	6.587	601	612	618	rBV3	9647	25482	0.63%	0.041%
13	6.687	620	629	636	rBV5	22834	44381	1.09%	0.071%
14	6.863	649	659	675	rVB	747607	1233986	30.34%	1.971%
15	7.034	682	688	700	rVB	634396	972323	23.91%	1.553%
16	7.228	712	721	730	rBV	773140	1231334	30.28%	1.967%
17	7.311	730	735	739	rBV5	13090	22771	0.56%	0.036%
18	7.569	775	779	788	rBV3	11797	32633	0.80%	0.052%
19	7.693	792	800	807	rVV	903175	1423778	35.01%	2.275%
20	7.758	807	811	825	rVB3	59835	125068	3.08%	0.200%
21	7.934	834	841	848	rBV5	22567	44895	1.10%	0.072%
22	8.399	912	920	932	rBV	751800	1241321	30.52%	1.983%
23	8.516	935	940	945	rVB	51685	86059	2.12%	0.137%
24	8.846	990	996	1004	rVV	730786	1211392	29.79%	1.935%
25	8.910	1004	1007	1012	rVV5	12841	23753	0.58%	0.038%
26	8.963	1012	1016	1020	rVV5	11886	26024	0.64%	0.042%
27	9.010	1020	1024	1032	rVB6	17624	33449	0.82%	0.053%
28	9.175	1045	1052	1059	rVV	856159	1316794	32.38%	2.104%
29	9.257	1059	1066	1072	rVB3	23951	47860	1.18%	0.076%
30	9.416	1086	1093	1102	rVB	27522	48702	1.20%	0.078%
31	9.563	1111	1118	1125	rBV	600772	996671	24.51%	1.592%
32	9.793	1153	1157	1167	rVB3	15855	30764	0.76%	0.049%
33	10.028	1191	1197	1203	rBV	57207	110827	2.73%	0.177%
34	10.104	1203	1210	1219	rVV	976202	1640279	40.33%	2.620%
35	10.193	1219	1225	1232	rVV	95634	184733	4.54%	0.295%
36	10.257	1232	1236	1244	rVB4	12681	25475	0.63%	0.041%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
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37	10.481	1266	1274	1279	rBV	1232673	2077386	51.08%	3.319%
38	10.528	1279	1282	1287	rVV	65299	103608	2.55%	0.166%
39	10.581	1287	1291	1295	rVV	33141	57206	1.41%	0.091%
40	10.616	1295	1297	1300	rVV4	18499	28293	0.70%	0.045%
41	10.669	1300	1306	1314	rVB	107298	187394	4.61%	0.299%
42	11.022	1358	1366	1373	rBV	15135	30804	0.76%	0.049%
43	11.128	1377	1384	1393	rVB2	22742	44271	1.09%	0.071%
44	11.222	1393	1400	1404	rBV	19221	35005	0.86%	0.056%
45	11.263	1404	1407	1412	rVB4	17134	24853	0.61%	0.040%
46	11.381	1420	1427	1431	rBV	34371	64743	1.59%	0.103%
47	12.069	1536	1544	1550	rVV2	22594	46617	1.15%	0.074%
48	12.140	1550	1556	1561	rVV2	18218	34495	0.85%	0.055%
49	12.540	1619	1624	1629	rVV5	14279	26154	0.64%	0.042%
50	13.163	1725	1730	1735	rVV2	32493	51539	1.27%	0.082%
51	13.234	1735	1742	1750	rVV4	26130	61723	1.52%	0.099%
52	13.304	1750	1754	1760	rVV9	16775	29068	0.71%	0.046%
53	13.751	1824	1830	1835	rVV	1697443	2342595	57.60%	3.742%
54	13.798	1835	1838	1847	rVV	103265	149497	3.68%	0.239%
55	13.893	1847	1854	1860	rVB4	20409	42610	1.05%	0.068%
56	14.028	1865	1877	1887	rBV	1970017	3028237	74.46%	4.838%
57	14.128	1889	1894	1900	rVB10	12138	26082	0.64%	0.042%
58	14.240	1909	1913	1919	rVB2	15542	23221	0.57%	0.037%
59	14.340	1919	1930	1936	rBV	1858333	2821371	69.37%	4.507%
60	14.398	1936	1940	1948	rVB	93542	146843	3.61%	0.235%
61	14.540	1958	1964	1980	rBV	572942	985218	24.22%	1.574%
62	14.740	1993	1998	2006	rVB	54232	118404	2.91%	0.189%
63	15.145	2061	2067	2073	rBV	88266	141204	3.47%	0.226%
64	15.198	2073	2076	2079	rVV	35528	41087	1.01%	0.066%
65	15.334	2093	2099	2105	rBV	2598223	3633800	89.35%	5.805%
66	15.387	2105	2108	2113	rVB	95369	131766	3.24%	0.211%
67	15.451	2113	2119	2127	rBV	390715	556372	13.68%	0.889%
68	16.063	2219	2223	2229	rBV4	38496	77520	1.91%	0.124%
69	16.116	2229	2232	2237	rVB2	27002	38736	0.95%	0.062%
70	16.269	2249	2258	2260	rBV4	45724	94918	2.33%	0.152%
71	16.298	2260	2263	2273	rVB2	142186	238278	5.86%	0.381%
72	16.492	2293	2296	2304	rVB4	57574	93526	2.30%	0.149%
73	16.598	2311	2314	2317	rVB	40055	45545	1.12%	0.073%
74	16.739	2335	2338	2344	rVB	32978	44418	1.09%	0.071%
75	16.804	2346	2349	2353	rBV4	71962	110977	2.73%	0.177%
76	16.857	2353	2358	2363	rVV	1134848	1427467	35.10%	2.281%

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77	16.904	2363	2366	2370	rVB	62913	84152	2.07%	0.134%
78	16.963	2371	2376	2380	rBV	294330	399292	9.82%	0.638%
79	17.086	2388	2397	2401	rBV	2165008	3149519	77.44%	5.032%
80	17.128	2401	2404	2409	rVB	906157	1139316	28.01%	1.820%
81	17.186	2409	2414	2419	rBV2	2445735	3523322	86.63%	5.629%
82	17.492	2462	2466	2470	rVB	110215	143887	3.54%	0.230%
83	17.798	2514	2518	2524	rVB	324801	424601	10.44%	0.678%
84	17.869	2526	2530	2534	rBV4	87707	136433	3.35%	0.218%
85	17.992	2546	2551	2560	rBV2	974767	1362097	33.49%	2.176%
86	18.157	2575	2579	2583	rVB2	101630	152270	3.74%	0.243%
87	18.286	2598	2601	2604	rBV	102544	124420	3.06%	0.199%
88	18.475	2630	2633	2636	rBV3	67846	85269	2.10%	0.136%
89	18.504	2636	2638	2643	rVB2	79003	103698	2.55%	0.166%
90	19.151	2743	2748	2757	rVB	1350309	1868910	45.95%	2.986%
91	19.280	2766	2770	2779	rBV	602996	877915	21.59%	1.403%
92	19.486	2800	2805	2808	rBV	2850089	4067044	100.00%	6.497%
93	19.516	2808	2810	2819	rVB	1077152	1360012	33.44%	2.173%
94	20.427	2962	2965	2970	rVB	410015	477999	11.75%	0.764%
95	20.845	3033	3036	3042	rBV	820953	990065	24.34%	1.582%
96	21.016	3062	3065	3073	rVB3	546585	749332	18.42%	1.197%
97	21.239	3100	3103	3108	rBV	1156301	1415890	34.81%	2.262%
98	21.327	3112	3118	3122	rVV2	1446460	2427208	59.68%	3.878%
99	24.074	3579	3585	3592	rVB2	927305	2077935	51.09%	3.320%
100	24.274	3614	3619	3640	rVB	992497	3150550	77.47%	5.033%

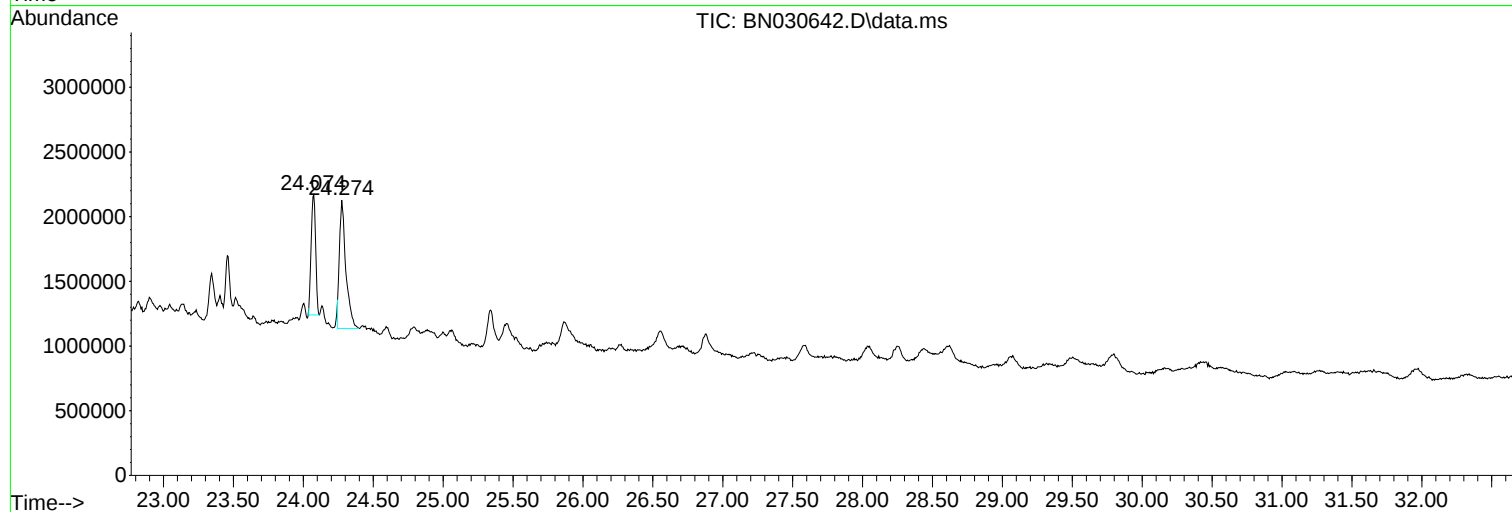
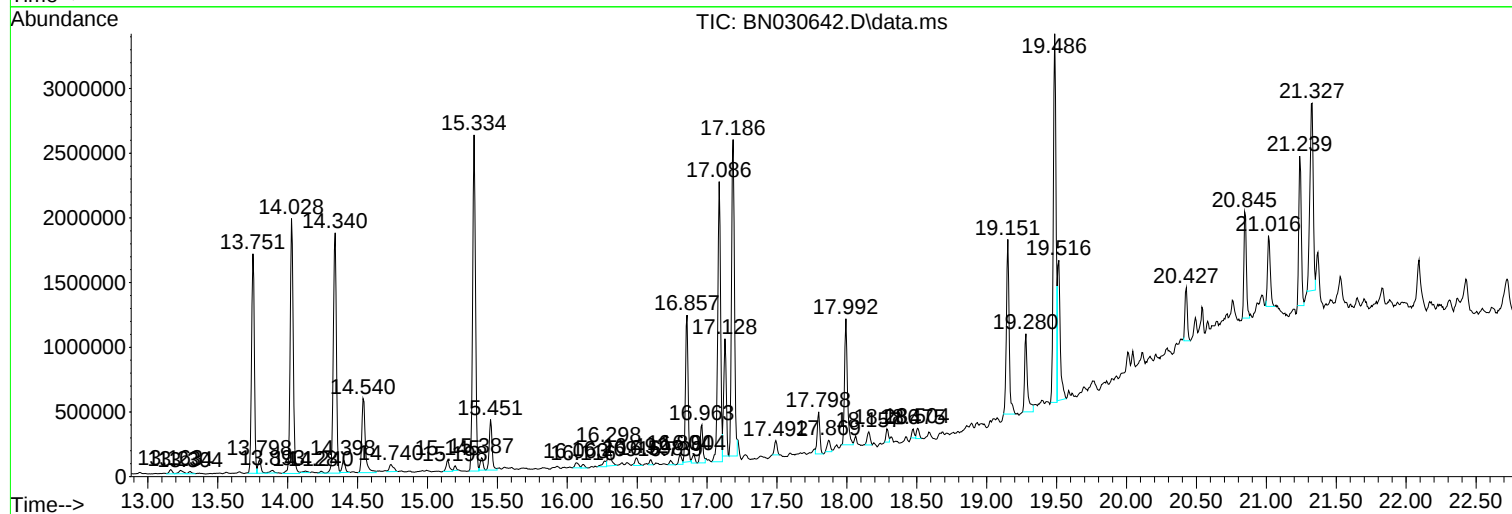
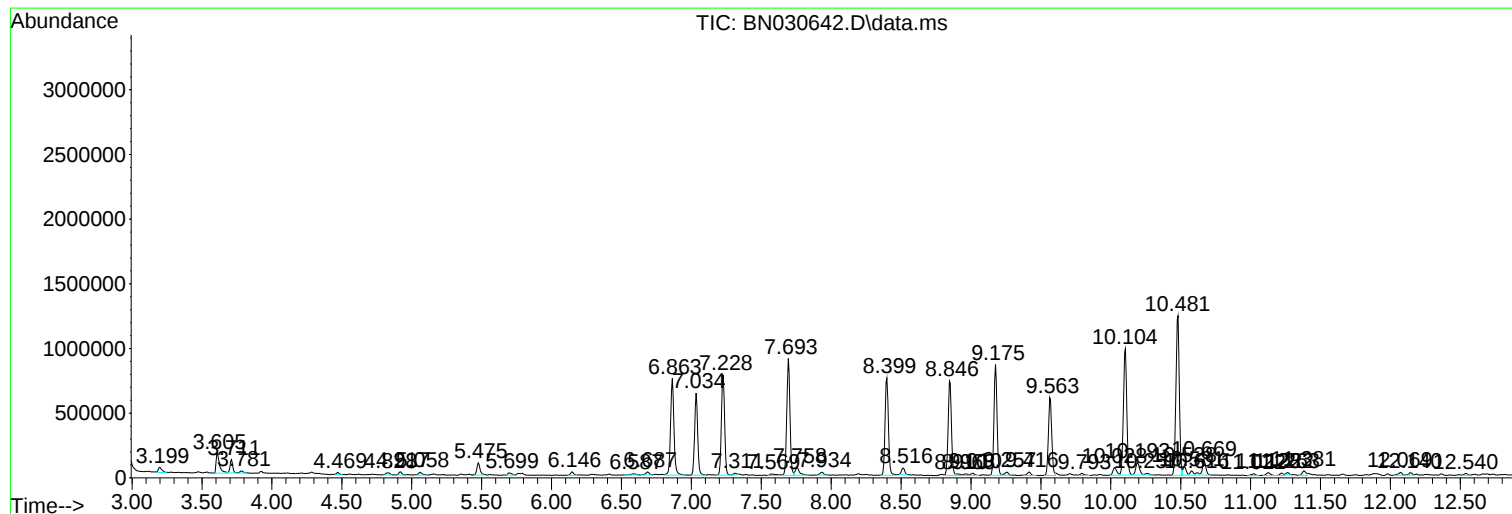
Sum of corrected areas: 62594402

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION

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



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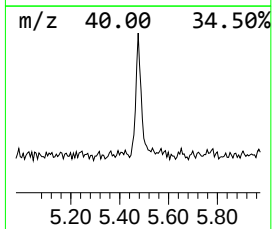
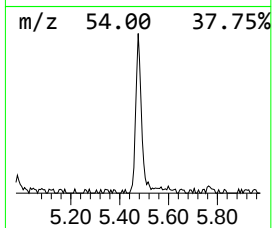
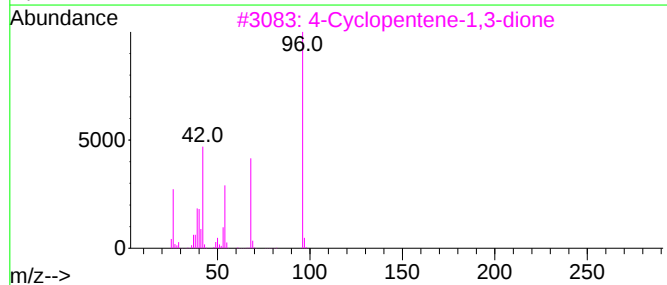
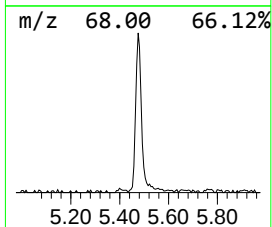
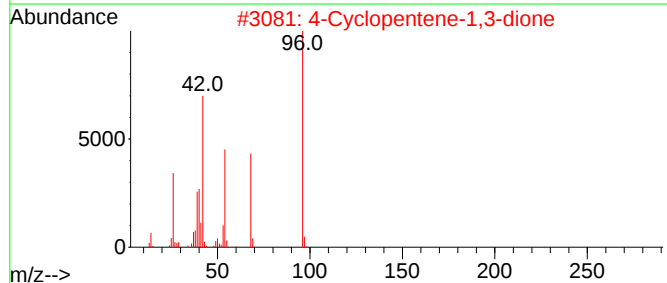
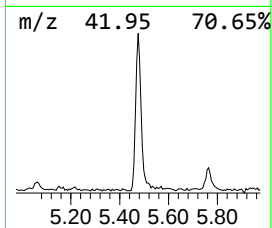
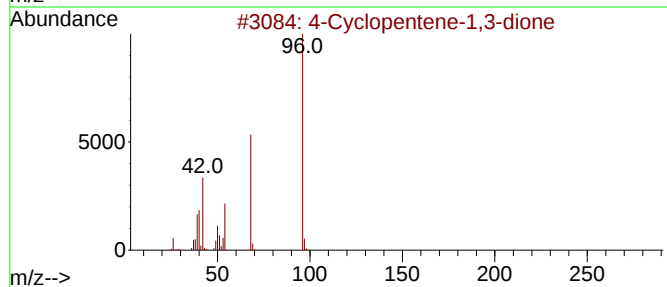
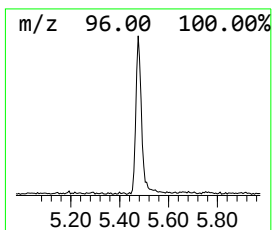
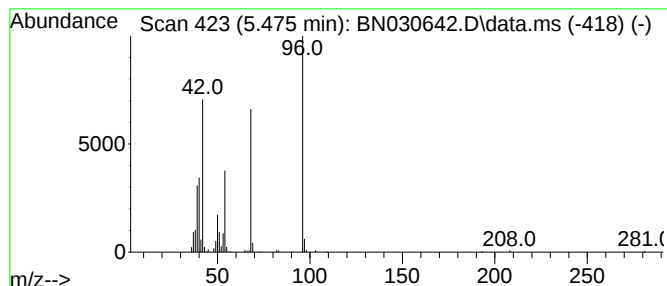
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 4-Cyclopentene-1,3-dione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.475	2.15 ng/ul	152905	1,4-Dichlorobenzene-d4	7.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Cyclopentene-1,3-dione	96	C5H4O2	000930-60-9	91
2			4-Cyclopentene-1,3-dione	96	C5H4O2	000930-60-9	80
3			4-Cyclopentene-1,3-dione	96	C5H4O2	000930-60-9	80
4			Pyrimidine, 4-hydroxy-	96	C4H4N2O	051953-18-5	64
5			4(1H)-Pyrimidinone	96	C4H4N2O	004562-27-0	53



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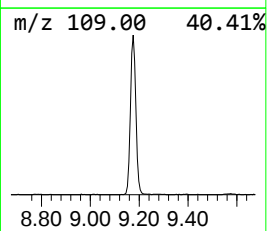
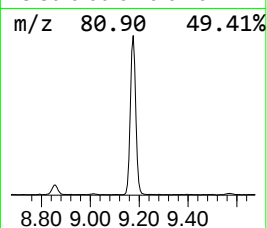
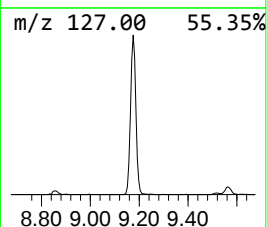
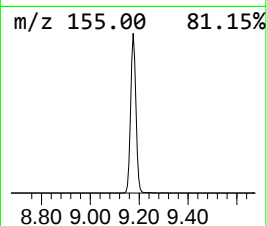
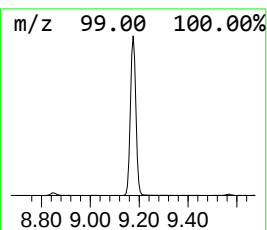
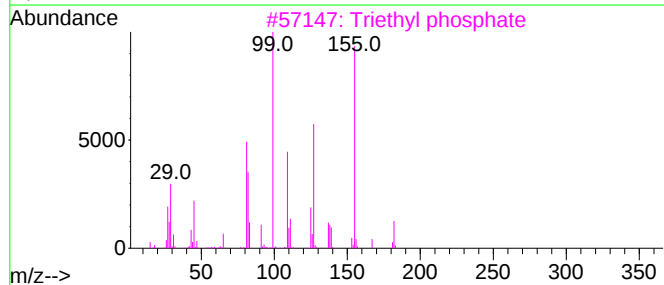
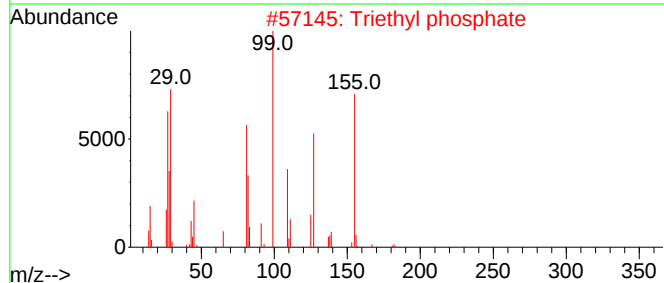
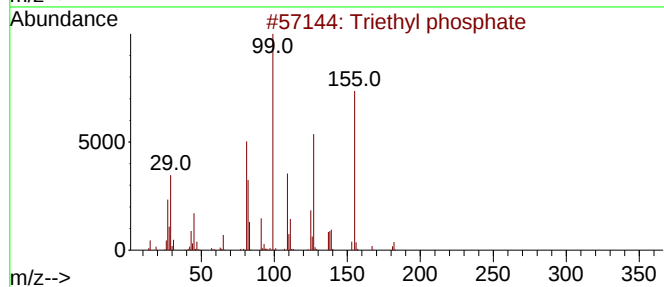
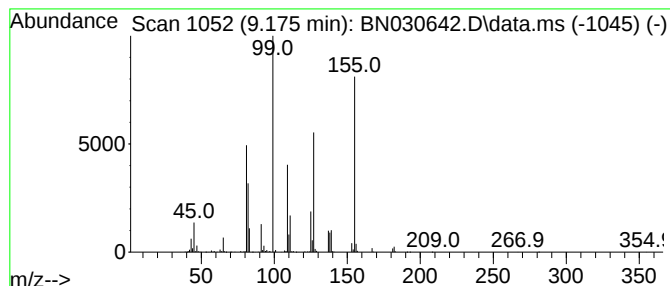
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Triethyl phosphate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.175	12.68 ng/ul	1316790	Naphthalene-d8	10.481

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethyl phosphate	182	C6H15O4P	000078-40-0	95
2			Triethyl phosphate	182	C6H15O4P	000078-40-0	94
3			Triethyl phosphate	182	C6H15O4P	000078-40-0	91
4			Triethyl phosphate	182	C6H15O4P	000078-40-0	91
5			Butyl diethyl phosphate	210	C8H19O4P	1000298-58-0	50



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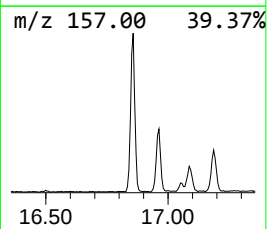
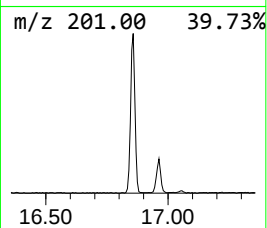
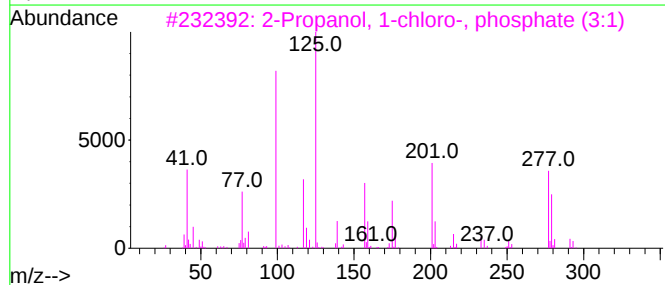
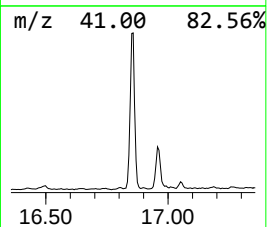
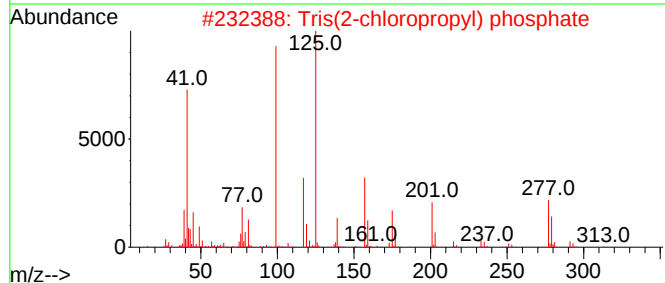
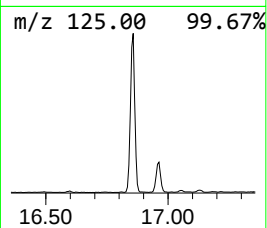
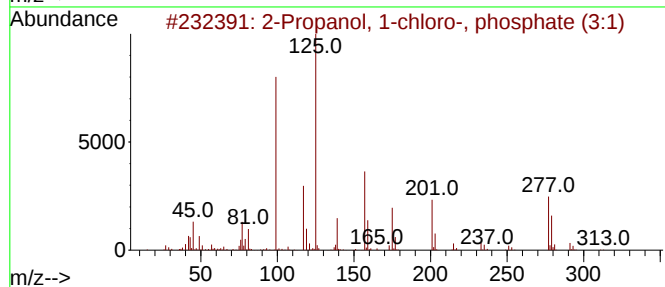
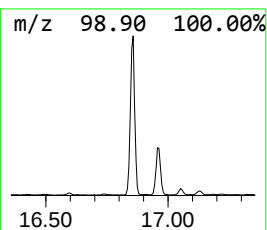
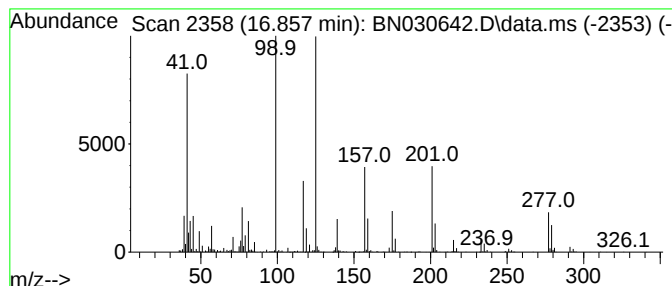
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1-chloro-, phos... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.857	9.06 ng/ul	1427470	Phenanthrene-d10	17.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	90		
2	Tris(2-chloropropyl) phosphate	326	C9H18Cl3O4P	006145-73-9	90		
3	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	90		
4	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	58		
5	Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
Acq On : 04 Apr 2024 00:42
Operator : MA/JU
Sample : P1849-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EOLY8

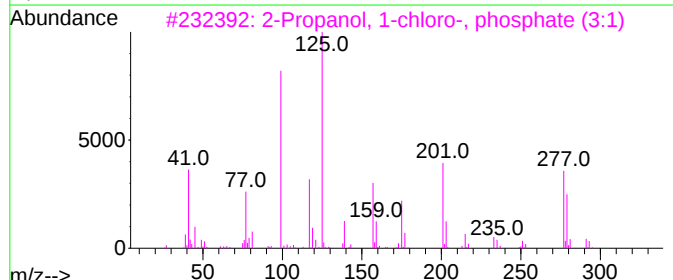
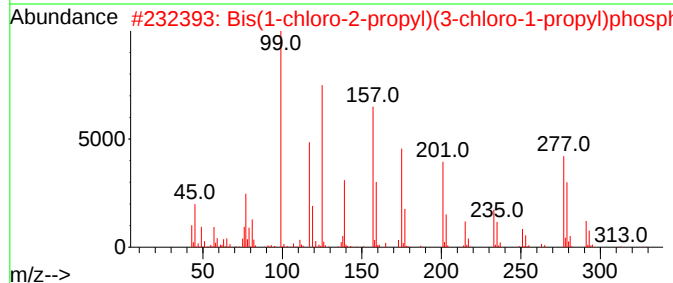
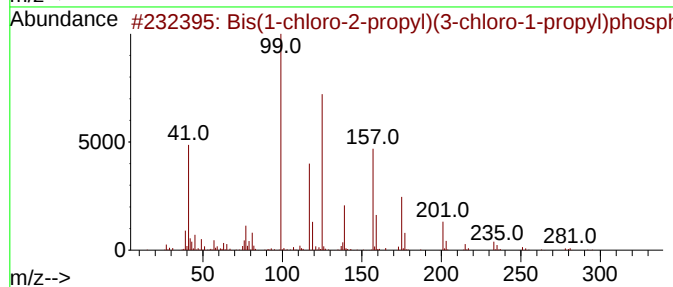
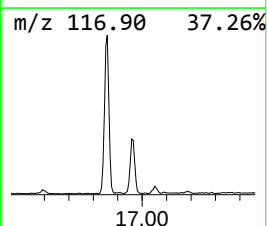
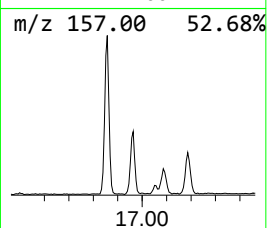
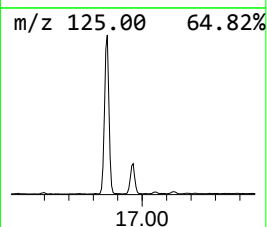
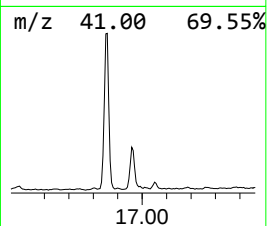
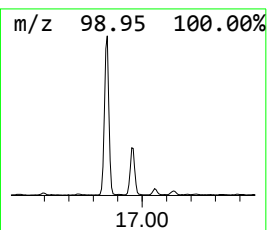
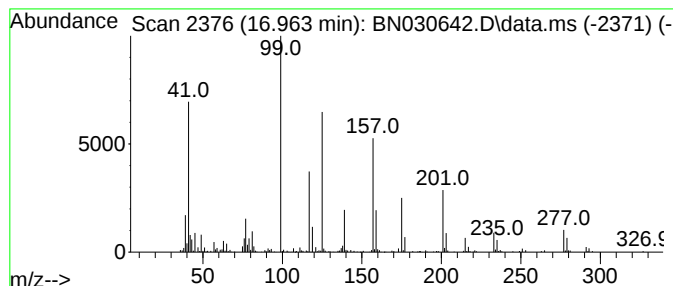
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bis(1-chloro-2-propyl)(3-ch... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.963	2.54 ng/ul	399292	Phenanthrene-d10	17.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	90
2			Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	80
3			2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	50
4			2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	50
5			Bis(3-chloro-1-propyl)(1-chloro-...	326	C9H18Cl3O4P	137888-35-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
Acq On : 04 Apr 2024 00:42
Operator : MA/JU
Sample : P1849-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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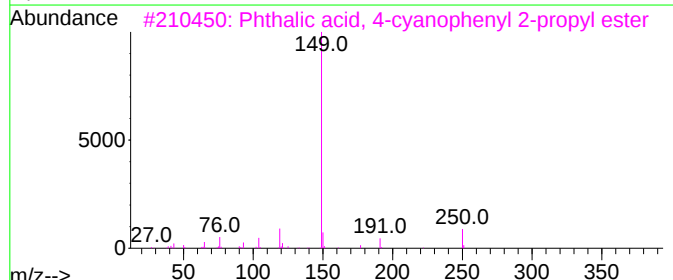
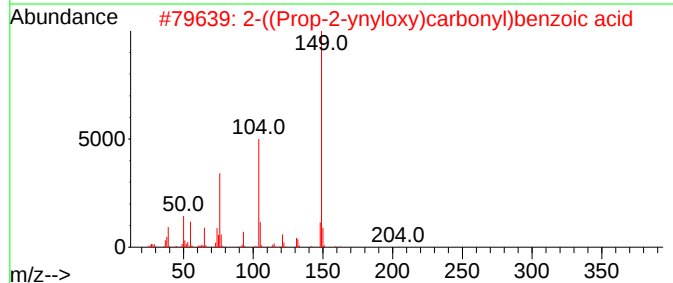
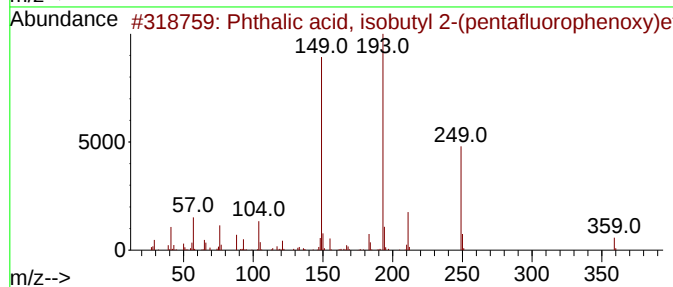
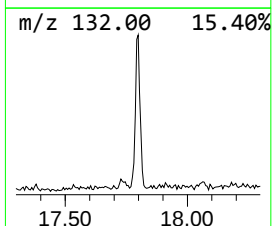
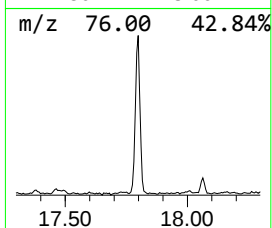
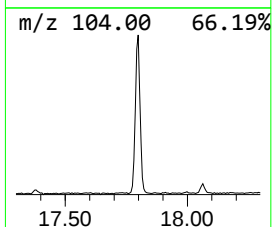
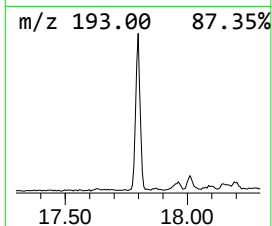
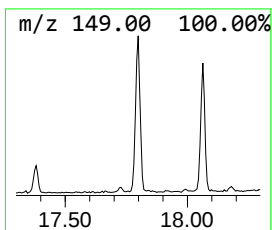
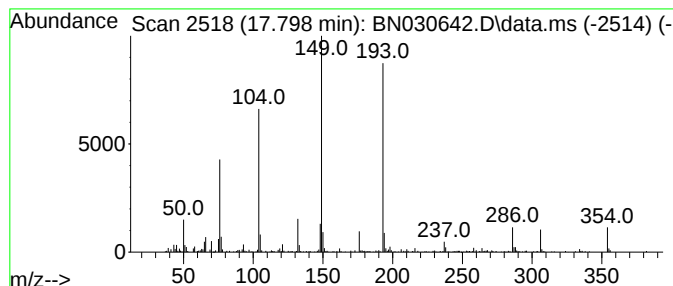
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.798	2.70 ng/ul	424601	Phenanthrene-d10	17.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, isobutyl 2-(penta...	432	C20H17F5O5	1000378-00-4	47
2			2-((Prop-2-ynyloxy)carbonyl)benz...	204	C11H8O4	006139-61-3	46
3			Phthalic acid, 4-cyanophenyl 2-p...	309	C18H15NO4	1000315-57-1	35
4			Phthalic acid, ethyl isopropyl e...	236	C13H16O4	1000314-99-6	18
5			Benzo[1,3]dioxole-5-carboxylic a...	354	C19H18N2O5	1000310-68-7	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
Acq On : 04 Apr 2024 00:42
Operator : MA/JU
Sample : P1849-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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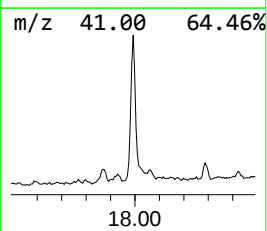
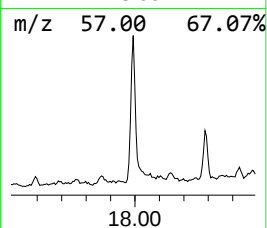
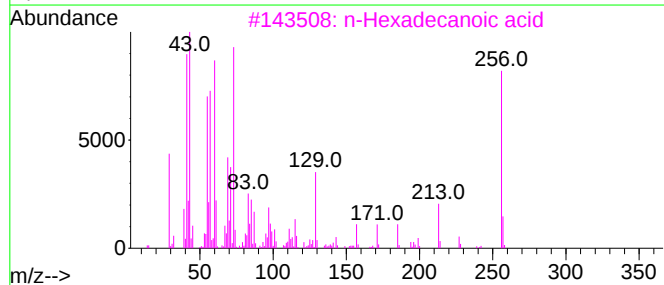
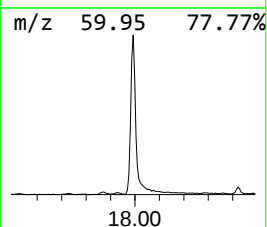
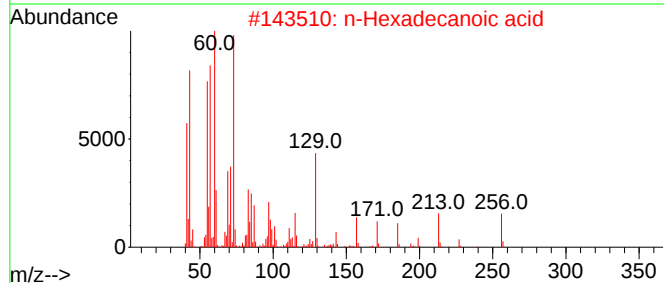
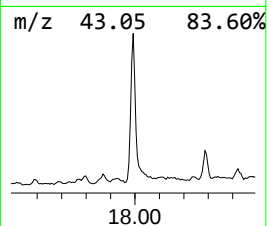
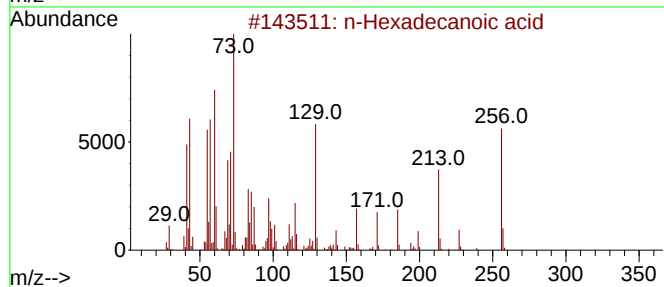
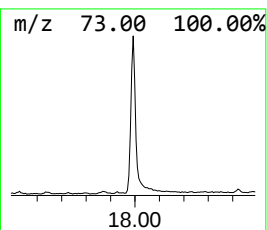
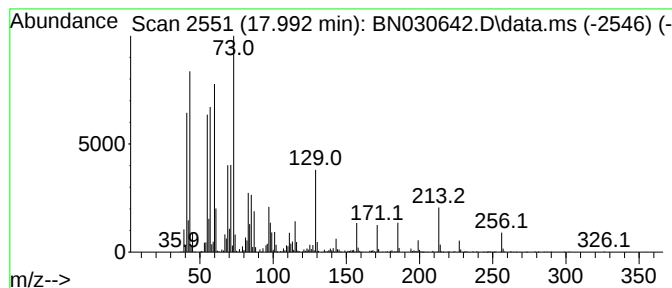
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.992	8.65 ng/ul	1362100	Phenanthrene-d10	17.087

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
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Operator : MA/JU
Sample : P1849-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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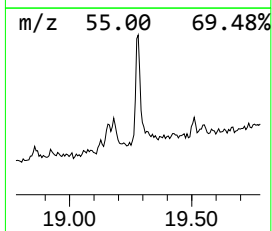
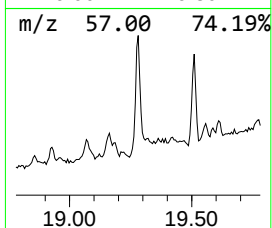
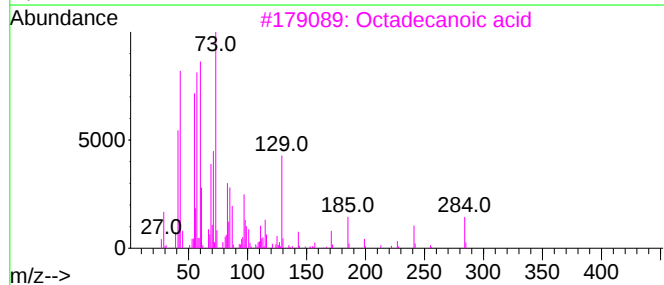
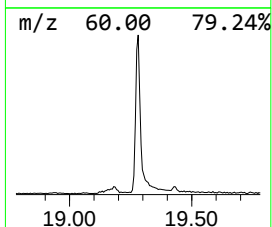
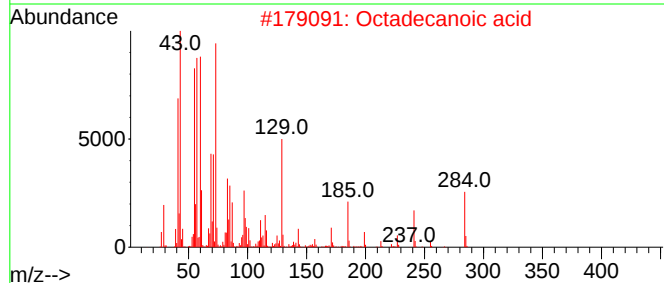
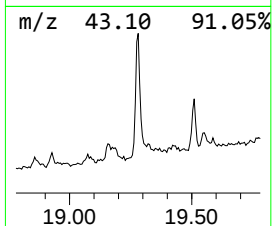
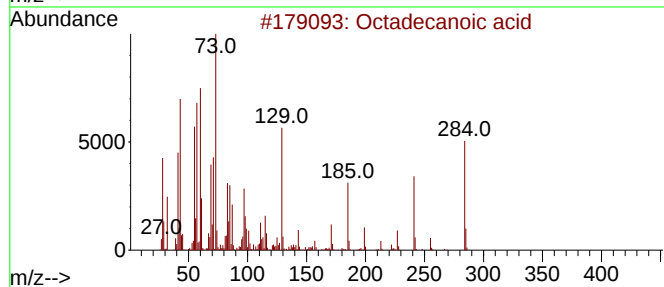
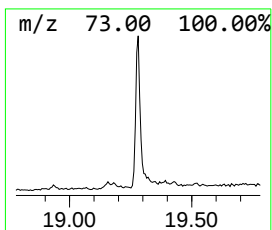
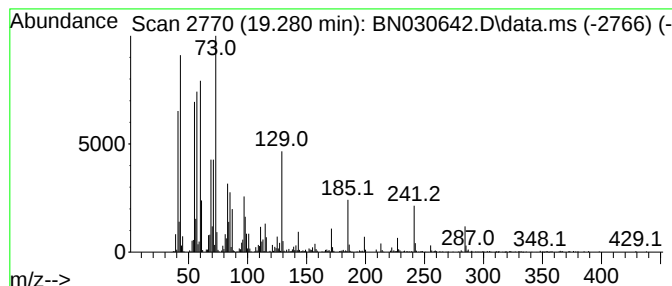
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.281	7.23 ng/ul	877915	Chrysene-d12	21.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
Acq On : 04 Apr 2024 00:42
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Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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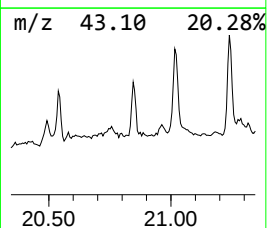
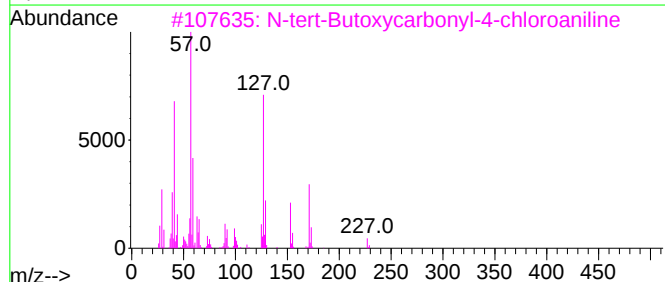
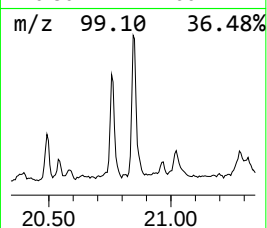
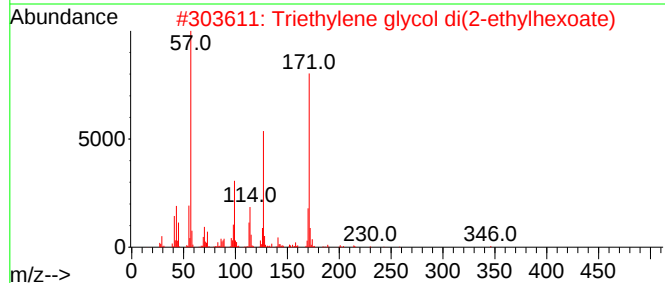
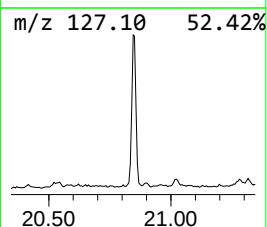
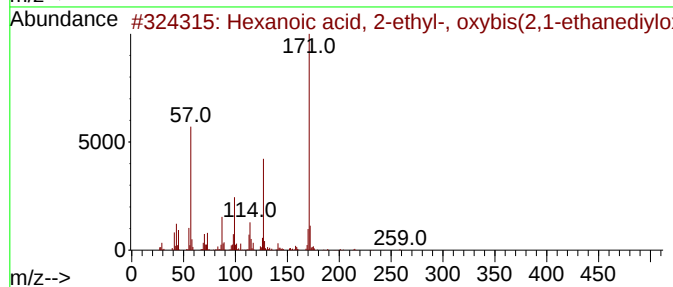
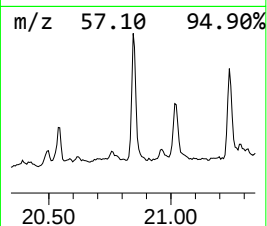
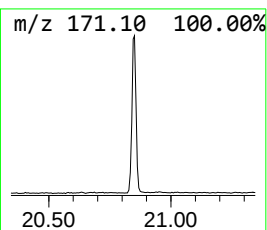
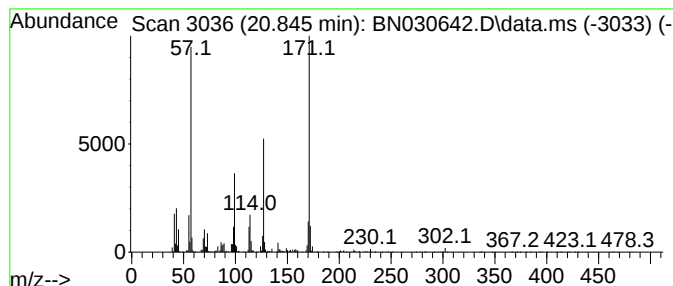
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexanoic acid, 2-ethyl-, ox... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.845	8.16 ng/ul	990065	Chrysene-d12	21.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-, oxybis(...	446	C24H46O7	018268-70-7	91
2			Triethylene glycol di(2-ethylhex...	402	C22H42O6	000094-28-0	58
3			N-tert-Butoxycarbonyl-4-chloroan...	227	C11H14ClNO2	018437-66-6	53
4			3-Chlorothiobenzamide	171	C7H6ClNS	002548-79-0	43
5			1-(1-Butoxypropan-2-yloxy)propan...	302	C17H34O4	1000367-12-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
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Sample : P1849-04
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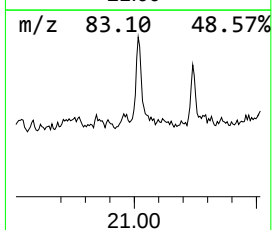
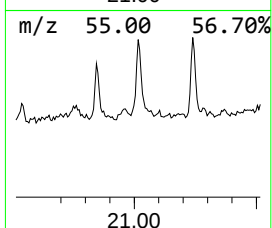
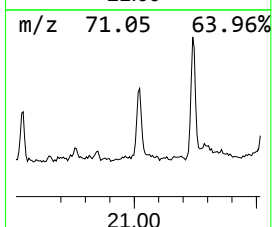
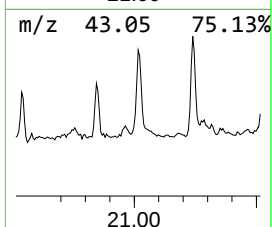
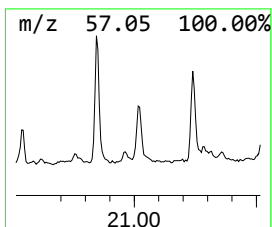
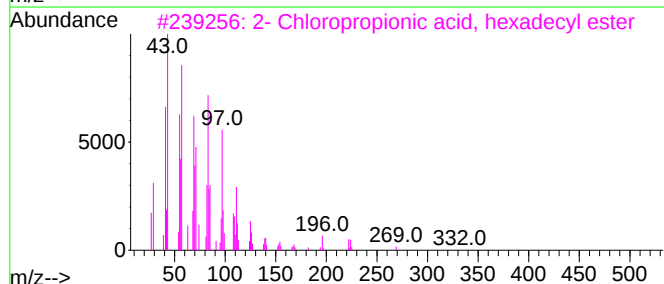
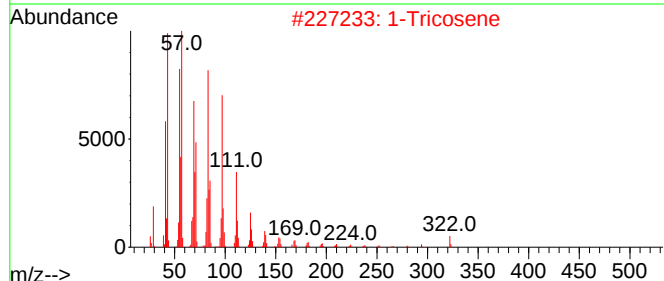
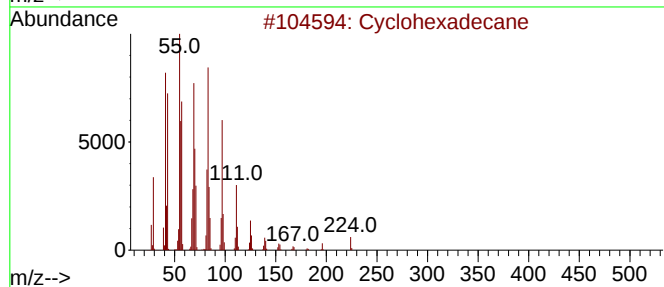
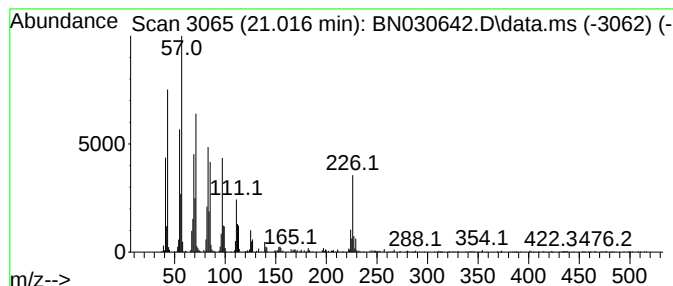
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic21.016 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.016	6.17 ng/ul	749332	Chrysene-d12	21.327

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	95
2			1-Tricosene	322	C23H46	018835-32-0	90
3			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	90
4			Cyclooctacosane	392	C28H56	000297-24-5	89
5			7-Hexadecene, (Z)-	224	C16H32	035507-09-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040424\
Data File : BN030642.D
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Sample : P1849-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.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-Cyclopentene-...	5.475	2.1	ng/ul	152905	1	7.693	1423780	20.0
Triethyl phosphate	9.175	12.7	ng/ul	1316790	2	10.481	2077390	20.0
2-Propanol, 1-c...	16.857	9.1	ng/ul	1427470	4	17.087	3149520	20.0
Bis(1-chloro-2-...	16.963	2.5	ng/ul	399292	4	17.087	3149520	20.0
unknown-01	17.798	2.7	ng/ul	424601	4	17.087	3149520	20.0
n-Hexadecanoic ...	17.992	8.7	ng/ul	1362100	4	17.087	3149520	20.0
Octadecanoic acid	19.281	7.2	ng/ul	877915	5	21.327	2427210	20.0
Hexanoic acid, ...	20.845	8.2	ng/ul	990065	5	21.327	2427210	20.0
(DEL) Alkane: C...	21.016	6.2	ng/ul	749332	5	21.327	2427210	20.0