

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
 Data File : BN032240.D
 Acq On : 25 Jun 2024 13:21
 Operator : MA/JU
 Sample : P2844-23
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4C48

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-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN032240.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.170	27	31	38	rVB	44144	61821	1.32%	0.092%
2	3.434	73	76	87	rBV	84997	141080	3.02%	0.210%
3	3.587	97	102	111	rBV3	59309	98193	2.10%	0.146%
4	3.675	113	117	125	rVB	58923	89627	1.92%	0.134%
5	3.875	148	151	156	rVB	21002	29085	0.62%	0.043%
6	4.375	231	236	240	rBV	17600	27892	0.60%	0.042%
7	4.422	240	244	251	rVB	38894	61338	1.31%	0.091%
8	4.769	297	303	311	rBV	126305	199238	4.26%	0.297%
9	4.875	317	321	327	rVB4	10019	17211	0.37%	0.026%
10	5.011	339	344	357	rBV	150379	254899	5.45%	0.380%
11	5.734	460	467	476	rVB	24814	44810	0.96%	0.067%
12	6.311	561	565	572	rVB2	20960	32472	0.69%	0.048%
13	6.611	611	616	621	rBV2	17969	30808	0.66%	0.046%
14	6.669	621	626	629	rVV3	8505	15177	0.32%	0.023%
15	6.816	642	651	669	rVV	745520	1305941	27.94%	1.948%
16	6.975	672	678	687	rVV	644795	1016941	21.75%	1.517%
17	7.052	687	691	697	rVB4	16243	26524	0.57%	0.040%
18	7.169	703	711	719	rBV	839303	1386408	29.66%	2.068%
19	7.240	719	723	733	rVB	52277	102934	2.20%	0.154%
20	7.546	771	775	779	rVB6	10793	16315	0.35%	0.024%
21	7.628	782	789	797	rBV	1123969	1818141	38.89%	2.712%
22	7.705	797	802	807	rVB3	13660	27603	0.59%	0.041%
23	7.834	820	824	828	rBV5	9231	17322	0.37%	0.026%
24	8.346	903	911	924	rBV	712069	1243355	26.60%	1.855%
25	8.452	924	929	937	rVB	55264	102704	2.20%	0.153%
26	8.787	978	986	996	rBV	744261	1282818	27.44%	1.914%
27	8.910	1004	1007	1010	rVV4	10149	16030	0.34%	0.024%
28	8.940	1010	1012	1017	rVB4	11910	17230	0.37%	0.026%
29	9.157	1041	1049	1054	rBV5	8975	19435	0.42%	0.029%
30	9.252	1058	1065	1070	rVV6	10116	27038	0.58%	0.040%
31	9.352	1074	1082	1094	rVB2	75452	147045	3.15%	0.219%
32	9.505	1101	1108	1121	rBV	614795	1086228	23.24%	1.620%
33	9.699	1135	1141	1145	rBV6	9620	21301	0.46%	0.032%
34	9.752	1147	1150	1157	rVV7	9269	16896	0.36%	0.025%
35	9.828	1157	1163	1173	rVB	87385	155634	3.33%	0.232%
36	10.040	1192	1199	1220	rBV	917292	1674856	35.83%	2.498%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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37	10.410	1253	1262	1269	rBV	1436613	2535248	54.23%	3.782%
38	10.563	1281	1288	1301	rBV	297667	636930	13.63%	0.950%
39	10.957	1344	1355	1361	rBV5	38594	75222	1.61%	0.112%
40	11.040	1364	1369	1374	rVB4	11008	19202	0.41%	0.029%

41	11.157	1381	1389	1396	rBV	176788	342735	7.33%	0.511%
42	11.775	1488	1494	1500	rBV8	9909	19119	0.41%	0.029%
43	11.881	1504	1512	1524	rBV	2860486	4573912	97.84%	6.823%
44	11.999	1529	1532	1540	rVB2	24257	47438	1.01%	0.071%
45	12.075	1540	1545	1550	rBV3	10809	20544	0.44%	0.031%

46	12.440	1603	1607	1613	rBV6	15295	26247	0.56%	0.039%
47	12.493	1613	1616	1621	rVB6	12836	20341	0.44%	0.030%
48	12.651	1640	1643	1648	rVB5	12132	15904	0.34%	0.024%
49	12.881	1675	1682	1687	rBV9	7010	17725	0.38%	0.026%
50	13.128	1721	1724	1728	rVB3	23616	34215	0.73%	0.051%

51	13.234	1738	1742	1749	rVB3	11015	20292	0.43%	0.030%
52	13.457	1776	1780	1784	rBV4	12020	16941	0.36%	0.025%
53	13.522	1789	1791	1797	rVB6	9107	14958	0.32%	0.022%
54	13.581	1797	1801	1806	rBV6	13496	27645	0.59%	0.041%
55	13.693	1813	1820	1826	rVV	1649703	2403849	51.42%	3.586%

56	13.751	1826	1830	1835	rVV4	28095	49675	1.06%	0.074%
57	13.793	1835	1837	1841	rVV5	10077	15204	0.33%	0.023%
58	13.969	1855	1867	1879	rBV2	2033320	3208503	68.64%	4.786%
59	14.169	1894	1901	1906	rBV3	19067	32850	0.70%	0.049%
60	14.275	1908	1919	1926	rBV	2168190	3334900	71.34%	4.975%

61	14.334	1926	1929	1938	rVB3	26469	62032	1.33%	0.093%
62	14.522	1950	1961	1976	rBV	601402	1257045	26.89%	1.875%
63	14.657	1976	1984	1990	rBV4	125941	293241	6.27%	0.437%
64	14.951	2030	2034	2038	rVB5	16957	23370	0.50%	0.035%
65	15.069	2049	2054	2058	rBV5	14266	28603	0.61%	0.043%

66	15.122	2062	2063	2067	rVV3	14510	17537	0.38%	0.026%
67	15.157	2067	2069	2074	rVB5	11255	15136	0.32%	0.023%
68	15.275	2082	2089	2096	rBV	2560846	3734847	79.90%	5.571%
69	15.328	2096	2098	2105	rVV3	27804	39474	0.84%	0.059%
70	15.416	2107	2113	2122	rBV	319683	483851	10.35%	0.722%

71	15.610	2141	2146	2153	rBV3	130394	193054	4.13%	0.288%
72	15.834	2180	2184	2189	rBV7	20890	40395	0.86%	0.060%
73	15.934	2198	2201	2205	rBV6	12128	16255	0.35%	0.024%
74	16.010	2207	2214	2219	rVV4	81611	136485	2.92%	0.204%
75	16.439	2284	2287	2292	rBV6	17132	27513	0.59%	0.041%

76	16.569	2305	2309	2313	rBV2	33447	46813	1.00%	0.070%
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77	16.692	2326	2330	2335	rVB2	22052	32813	0.70%	0.049%
78	16.934	2363	2371	2377	rBV2	25241	71951	1.54%	0.107%
79	17.028	2381	2387	2392	rVV	2571286	3780692	80.88%	5.640%
80	17.069	2392	2394	2398	rVV	283928	355091	7.60%	0.530%
81	17.128	2398	2404	2414	rVB	2612471	3907497	83.59%	5.829%
82	17.263	2424	2427	2432	rVB	46287	60240	1.29%	0.090%
83	17.504	2464	2468	2474	rVB2	96480	137893	2.95%	0.206%
84	17.592	2479	2483	2491	rBV9	34630	67161	1.44%	0.100%
85	17.698	2498	2501	2506	rVB5	43398	56364	1.21%	0.084%
86	17.898	2531	2535	2536	rBV	97324	121044	2.59%	0.181%
87	17.928	2536	2540	2549	rVV2	785823	1281089	27.40%	1.911%
88	17.998	2550	2552	2555	rVV	64057	70483	1.51%	0.105%
89	18.039	2556	2559	2563	rVB2	61837	79612	1.70%	0.119%
90	18.098	2565	2569	2573	rBV2	59869	108215	2.31%	0.161%
91	18.222	2586	2590	2593	rBV4	39572	58186	1.24%	0.087%
92	18.463	2628	2631	2637	rVB2	40169	50246	1.07%	0.075%
93	19.092	2731	2738	2746	rBV	708946	1162502	24.87%	1.734%
94	19.216	2755	2759	2771	rBV2	448479	793409	16.97%	1.184%
95	19.433	2790	2796	2800	rBV	3029329	4582977	98.04%	6.836%
96	20.951	3050	3054	3062	rVB2	834592	1202200	25.72%	1.793%
97	21.269	3102	3108	3113	rBV2	2830368	4454614	95.29%	6.645%
98	21.998	3229	3232	3235	rBV2	258825	372970	7.98%	0.556%
99	23.986	3564	3570	3578	rVB2	1327346	2999109	64.16%	4.474%
100	24.186	3597	3604	3614	rVB	1807461	4674678	100.00%	6.973%

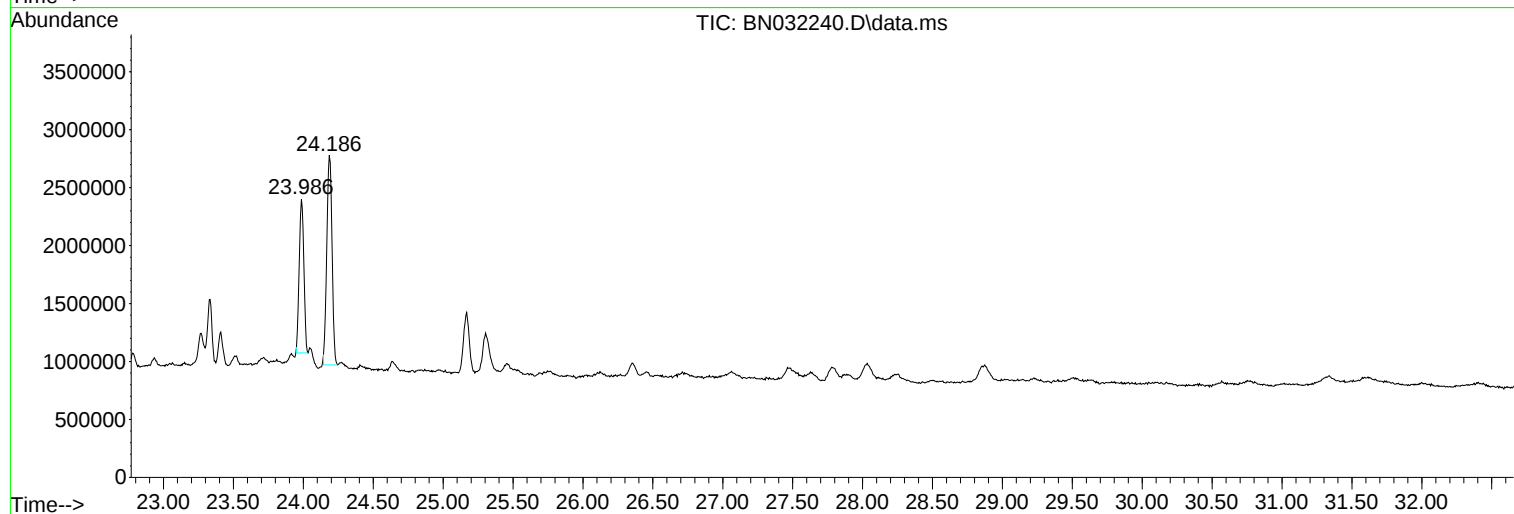
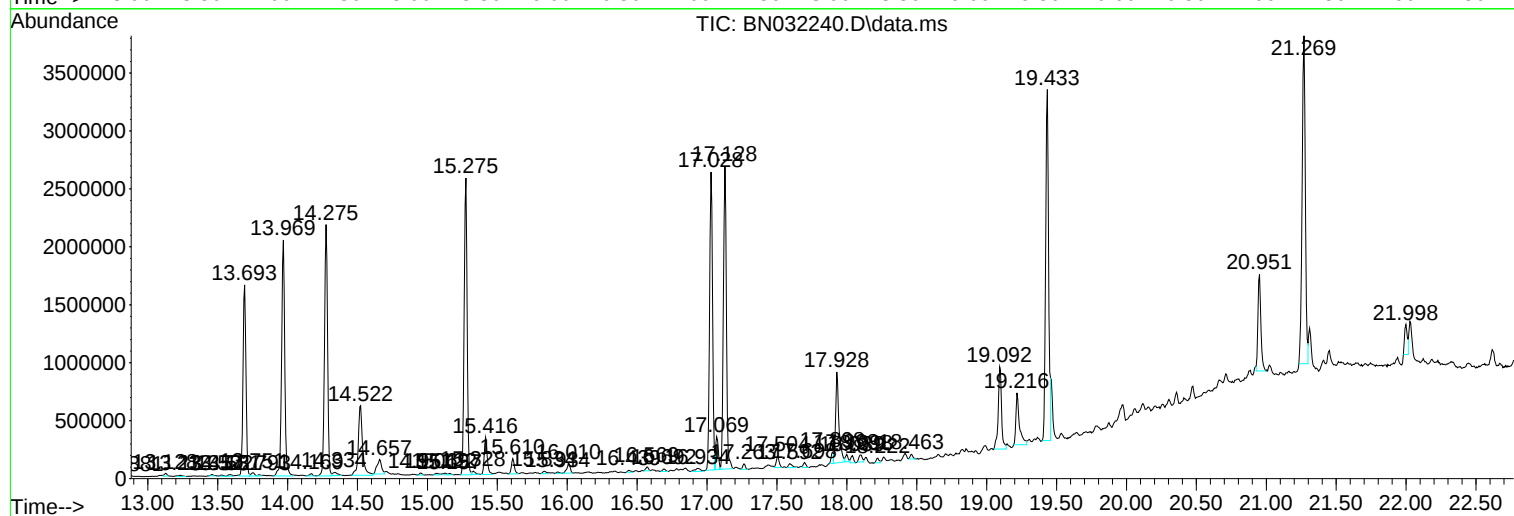
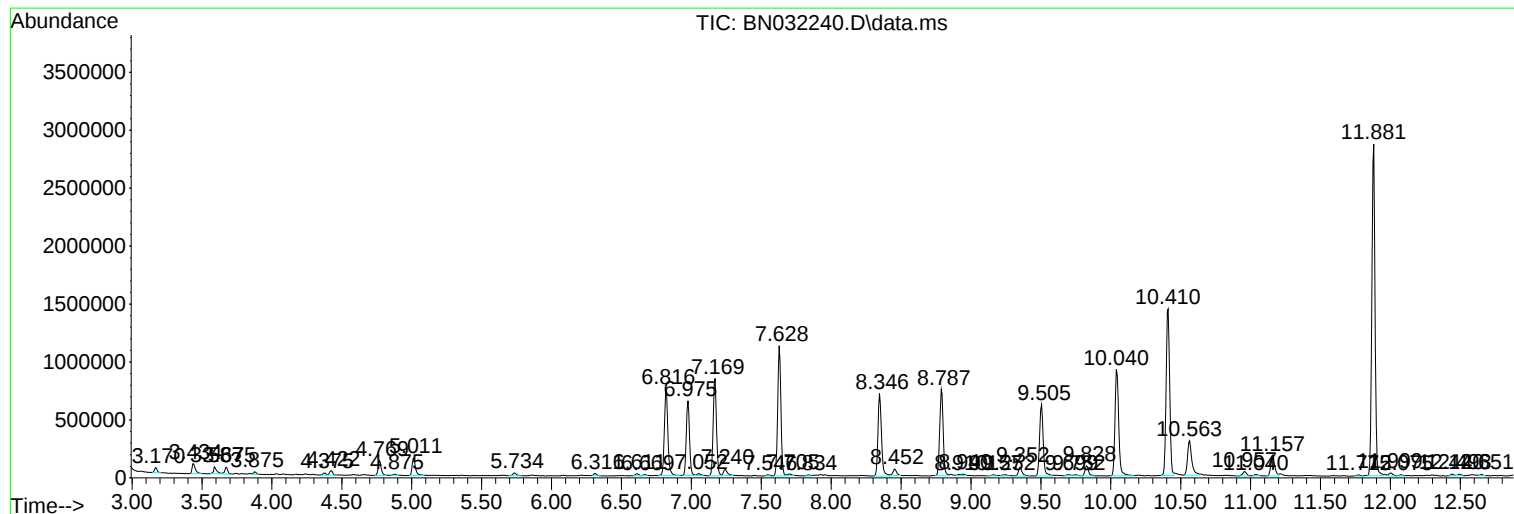
Sum of corrected areas: 67038636

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Instrument :
BNA_N
ClientSampleId :
A4C48

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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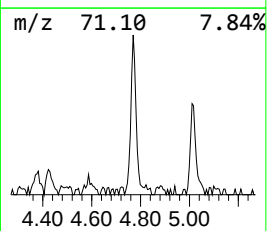
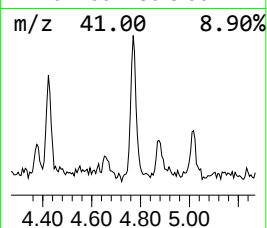
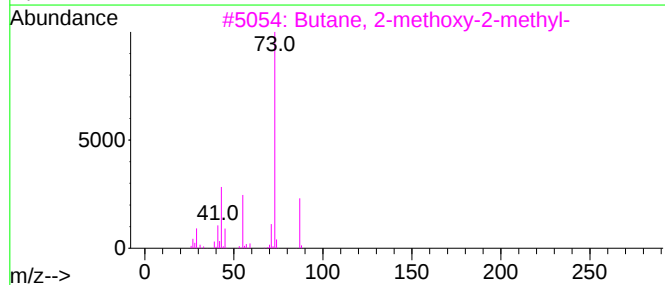
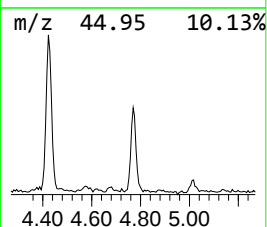
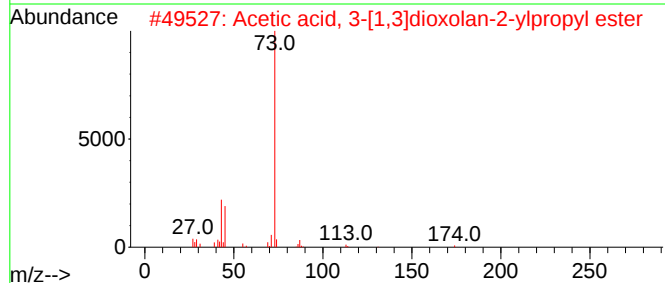
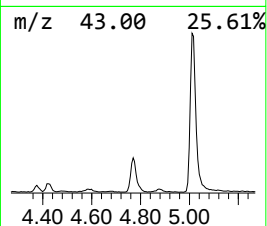
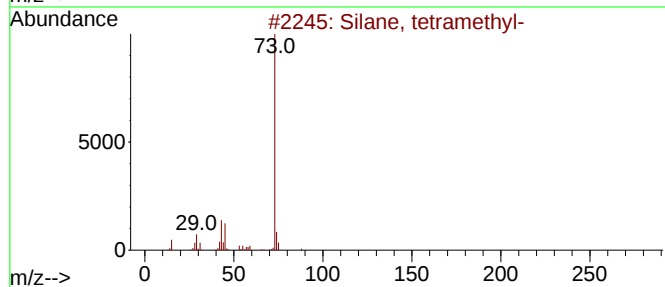
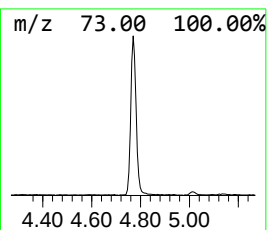
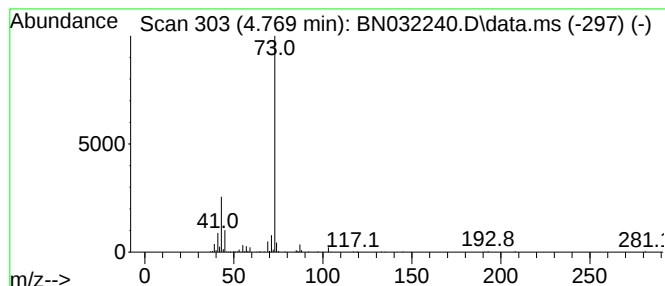
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.769	2.19 ng/ul	199238	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
2			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	33
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	28
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	28
5			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	25



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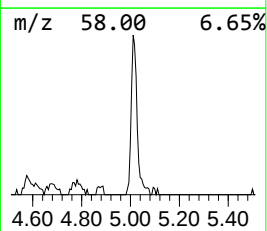
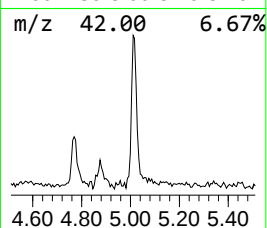
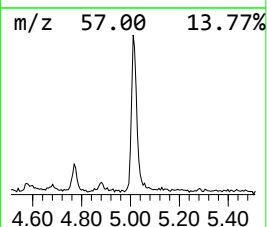
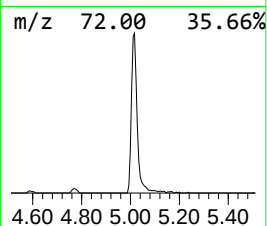
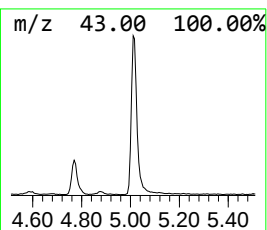
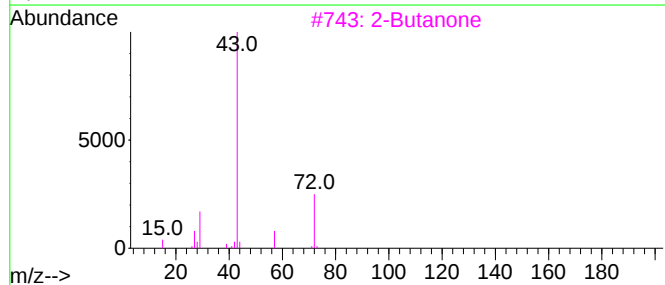
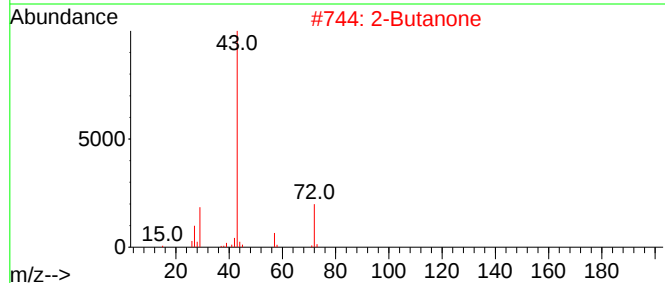
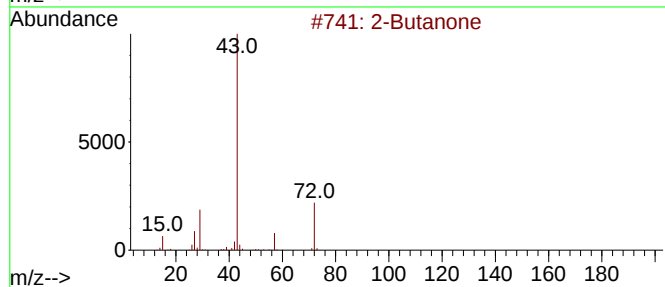
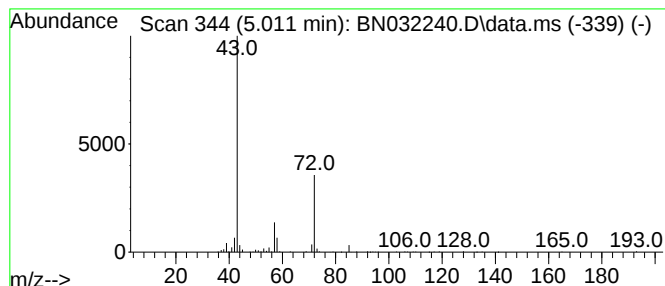
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Butanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.011	2.80 ng/ul	254899	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone			72	C4H8O	000078-93-3	72
2	2-Butanone			72	C4H8O	000078-93-3	64
3	2-Butanone			72	C4H8O	000078-93-3	59
4	2-Butanone			72	C4H8O	000078-93-3	59
5	2-Butanone			72	C4H8O	000078-93-3	59



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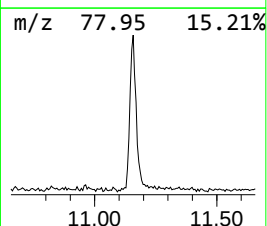
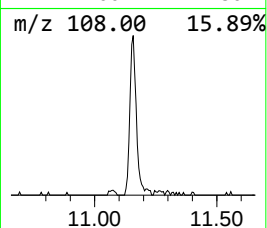
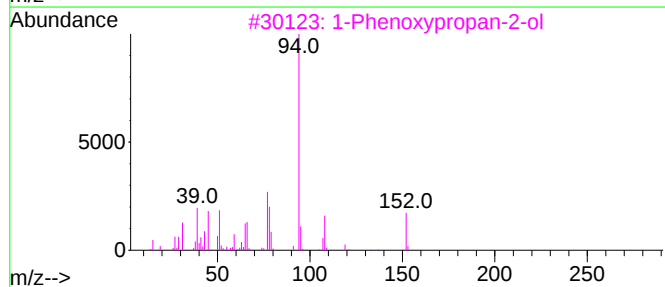
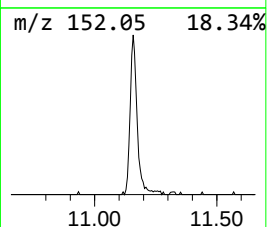
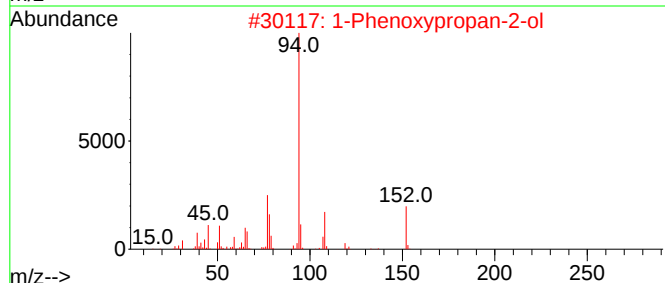
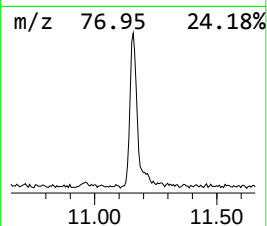
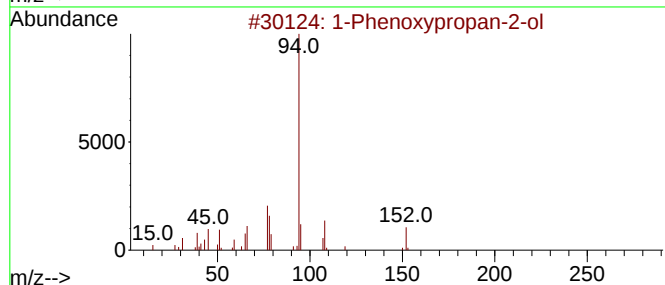
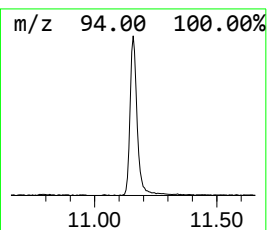
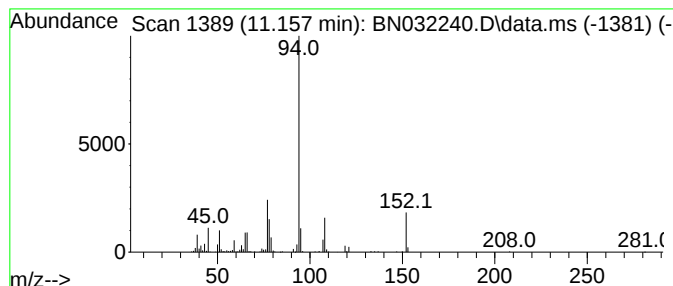
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.157	2.70 ng/ul	342735	Naphthalene-d8	10.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
Data File : BN032240.D
Acq On : 25 Jun 2024 13:21
Operator : MA/JU
Sample : P2844-23
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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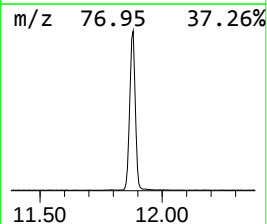
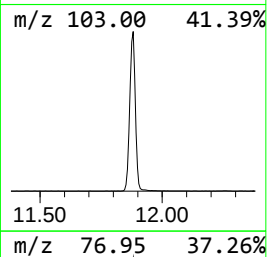
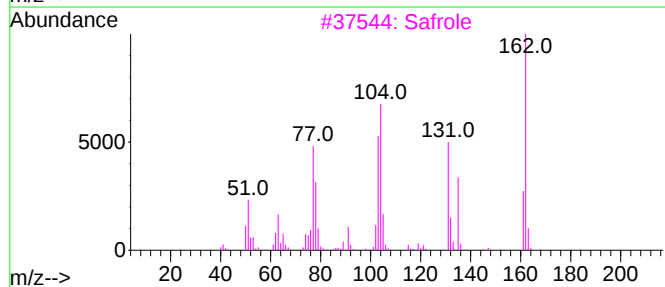
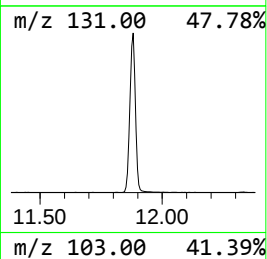
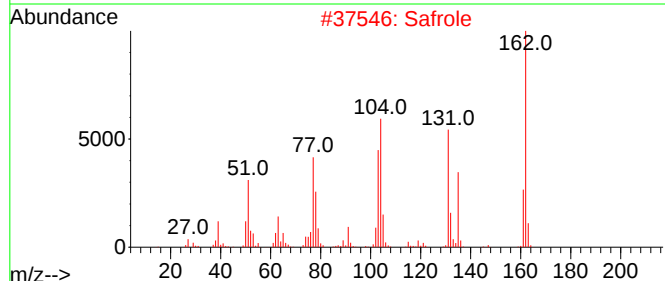
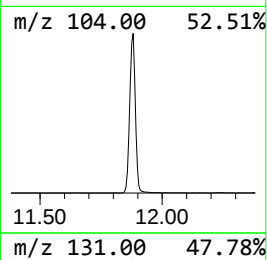
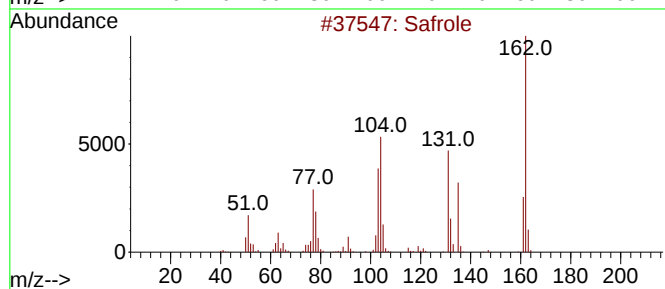
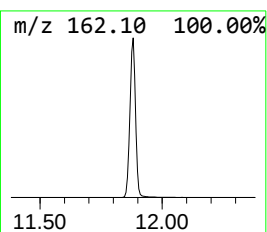
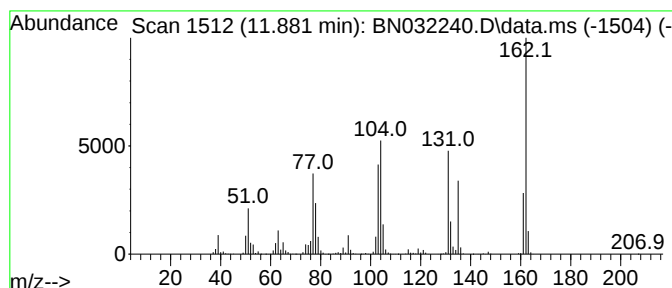
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Safrole Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.881	36.08 ng/ul	4573910	Naphthalene-d8	10.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Safrole	162	C10H10O2	000094-59-7	98
2			Safrole	162	C10H10O2	000094-59-7	98
3			Safrole	162	C10H10O2	000094-59-7	97
4			Safrole	162	C10H10O2	000094-59-7	96
5			1,3-Benzodioxole, 5-(1-propenyl)-	162	C10H10O2	000120-58-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
Data File : BN032240.D
Acq On : 25 Jun 2024 13:21
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Sample : P2844-23
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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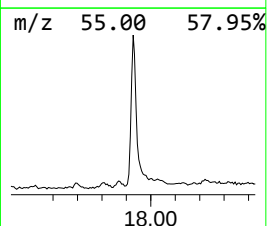
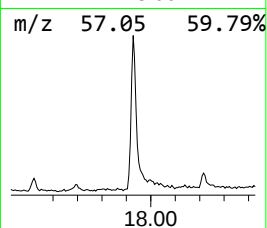
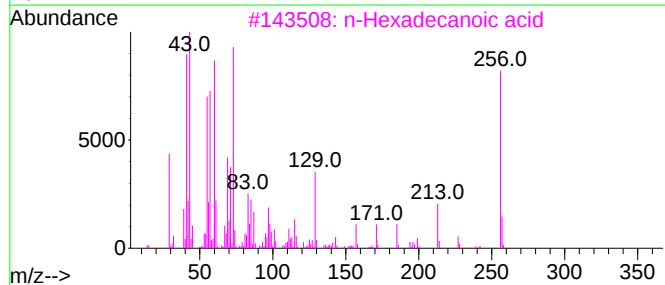
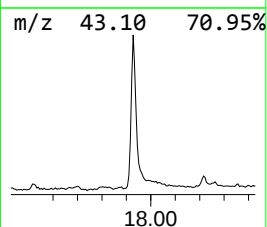
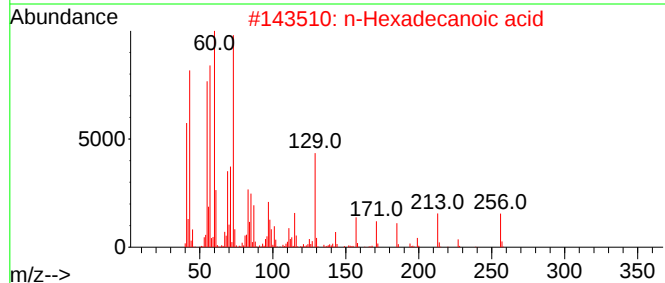
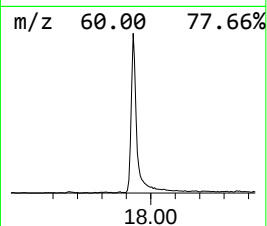
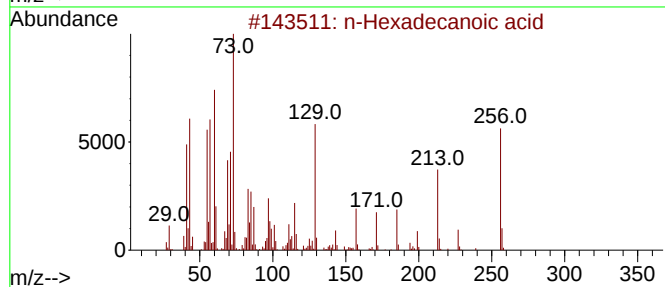
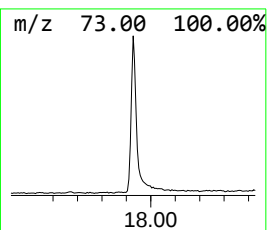
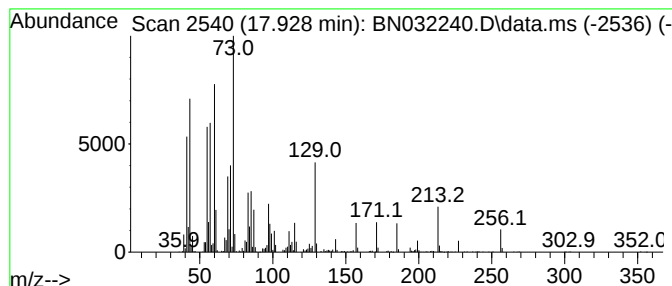
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.928	6.78 ng/ul	1281090	Phenanthrene-d10	17.028

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			Tridecanoic acid	214	C13H26O2	000638-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
Data File : BN032240.D
Acq On : 25 Jun 2024 13:21
Operator : MA/JU
Sample : P2844-23
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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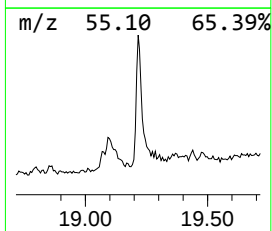
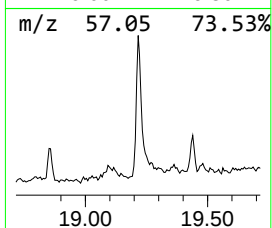
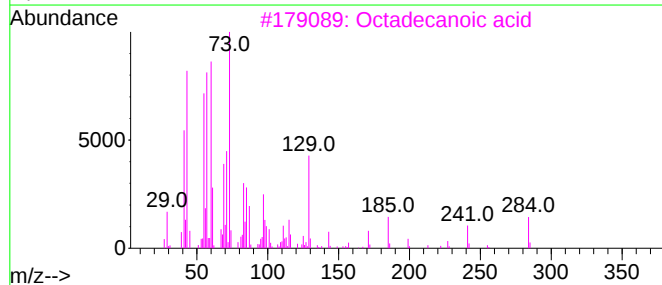
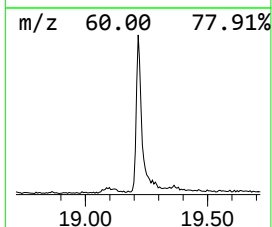
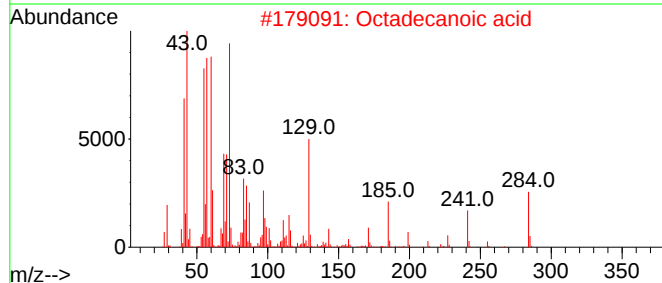
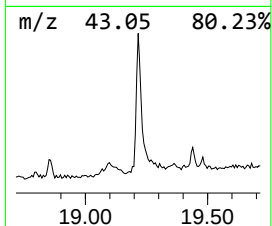
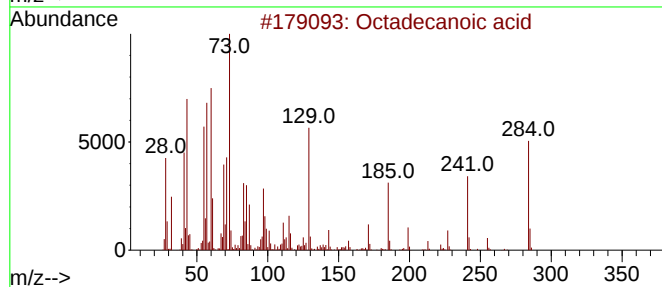
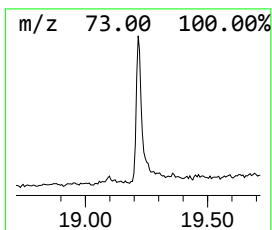
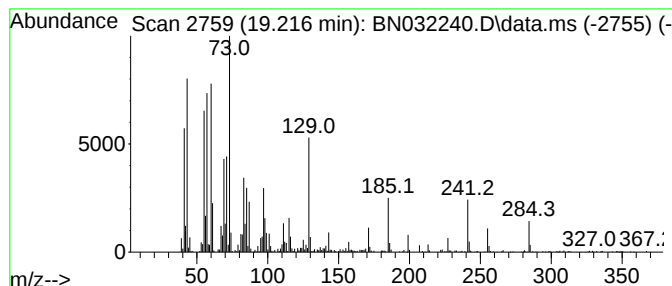
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.216	3.56 ng/ul	793409	Chrysene-d12	21.269

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	99
5			Octadecanoic acid	284	C18H36O2	000057-11-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
Data File : BN032240.D
Acq On : 25 Jun 2024 13:21
Operator : MA/JU
Sample : P2844-23
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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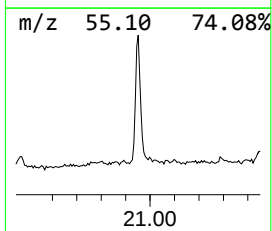
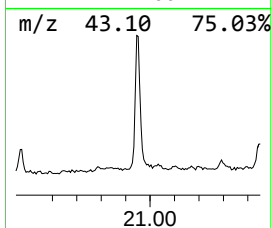
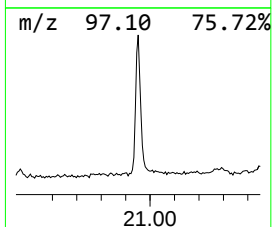
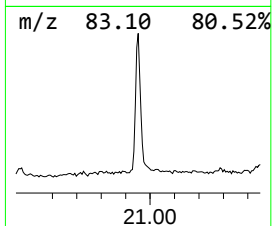
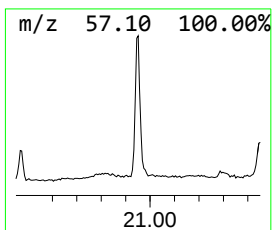
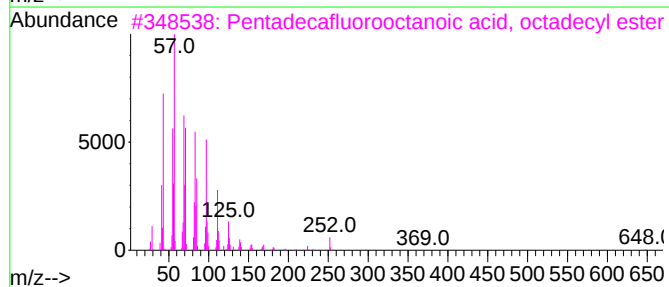
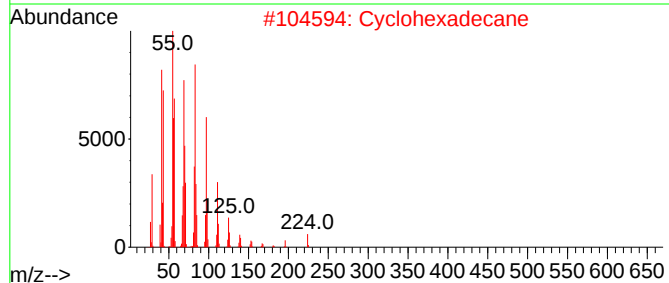
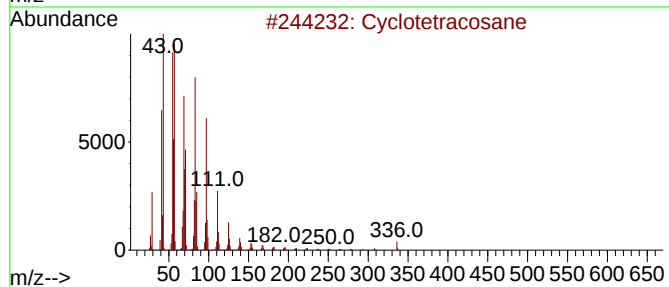
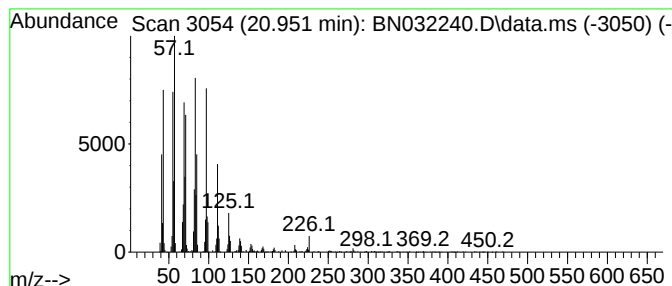
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic20.951 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.951	5.40 ng/ul	1202200	Chrysene-d12	21.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	95
2			Cyclohexadecane	224	C16H32	000295-65-8	94
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
4			1-Heneicosanol	312	C21H44O	015594-90-8	93
5			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062424\
Data File : BN032240.D
Acq On : 25 Jun 2024 13:21
Operator : MA/JU
Sample : P2844-23
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Butanone	5.011	2.8	ng/ul	254899	1	7.628	1818140	20.0
1-Phenoxypropan...	11.157	2.7	ng/ul	342735	2	10.410	2535250	20.0
Safrrole	11.881	36.1	ng/ul	4573910	2	10.410	2535250	20.0
n-Hexadecanoic ...	17.928	6.8	ng/ul	1281090	4	17.028	3780690	20.0
Octadecanoic acid	19.216	3.6	ng/ul	793409	5	21.269	4454610	20.0
(DEL) Alkane: C...	20.951	5.4	ng/ul	1202200	5	21.269	4454610	20.0