

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
 Data File : BN015102.D
 Acq On : 26 Jun 2021 06:15
 Operator : CG/JU
 Sample : M2704-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD86

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-BN061821MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015102.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	30	34	41	rVB	27913	39946	0.84%	0.095%
2	3.681	97	101	111	rBV	149424	237149	4.96%	0.565%
3	3.781	116	118	125	rVB3	14841	18475	0.39%	0.044%
4	3.834	125	127	134	rVB5	5991	11246	0.24%	0.027%
5	3.999	146	155	162	rVB	59417	95921	2.01%	0.229%
6	4.358	211	216	220	rBV2	21570	32744	0.69%	0.078%
7	4.481	234	237	241	rBV	16435	19877	0.42%	0.047%
8	4.534	241	246	257	rVB	56241	77665	1.63%	0.185%
9	5.281	368	373	381	rVB	42115	59523	1.25%	0.142%
10	5.411	388	395	406	rBV	123574	250154	5.24%	0.596%
11	5.775	451	457	465	rVB	59501	91899	1.92%	0.219%
12	5.840	465	468	474	rBV3	7522	12171	0.25%	0.029%
13	6.746	617	622	626	rVB3	6475	10501	0.22%	0.025%
14	6.893	640	647	657	rBV	448651	704746	14.75%	1.680%
15	7.063	669	676	685	rBV	364698	561708	11.76%	1.339%
16	7.163	689	693	698	rVB4	7508	10588	0.22%	0.025%
17	7.258	702	709	719	rVB	450020	729913	15.28%	1.740%
18	7.363	719	727	732	rBV4	6132	11595	0.24%	0.028%
19	7.728	782	789	795	rVV	626849	987145	20.66%	2.353%
20	7.787	795	799	805	rVB2	9835	14962	0.31%	0.036%
21	7.963	825	829	837	rVB3	13493	24873	0.52%	0.059%
22	8.275	875	882	886	rBV	9462	17102	0.36%	0.041%
23	8.416	899	906	914	rBV	524525	816220	17.08%	1.946%
24	8.534	921	926	933	rVB	19046	30259	0.63%	0.072%
25	8.810	967	973	977	rBV2	13271	20901	0.44%	0.050%
26	8.869	977	983	992	rVV	410183	658559	13.78%	1.570%
27	9.310	1052	1058	1068	rVB3	7082	12406	0.26%	0.030%
28	9.581	1096	1104	1111	rBV	285056	478088	10.01%	1.140%
29	10.116	1188	1195	1206	rBV	519139	837827	17.53%	1.997%
30	10.275	1217	1222	1234	rBV	88788	147209	3.08%	0.351%
31	10.499	1252	1260	1266	rBV	836665	1395768	29.21%	3.328%
32	10.546	1266	1268	1274	rVV	8440	10771	0.23%	0.026%
33	10.628	1274	1282	1295	rVV	476776	805714	16.86%	1.921%
34	11.540	1430	1437	1441	rBV6	4714	9764	0.20%	0.023%
35	11.934	1495	1504	1511	rBV4	12888	22914	0.48%	0.055%
36	12.087	1523	1530	1535	rVB2	9813	17278	0.36%	0.041%

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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-BN061821MA.M
 Title : SVOA CALIBRATION

37	12.157	1538	1542	1550	rVB4	7334	12427	0.26%	0.030%
38	12.898	1664	1668	1672	rBV2	10119	15975	0.33%	0.038%
39	13.187	1709	1717	1725	rBV	40093	67086	1.40%	0.160%
40	13.675	1795	1800	1802	rBV4	7985	10658	0.22%	0.025%
41	13.763	1809	1815	1821	rVB	865527	1222127	25.58%	2.914%
42	13.851	1826	1830	1834	rVB3	19738	26776	0.56%	0.064%
43	13.898	1834	1838	1843	rVB5	8990	16229	0.34%	0.039%
44	14.040	1853	1862	1873	rBV	1044357	1614717	33.79%	3.850%
45	14.269	1895	1901	1905	rBV	57233	78679	1.65%	0.188%
46	14.351	1908	1915	1924	rVV	1238292	1797981	37.63%	4.286%
47	14.539	1939	1947	1963	rVV	288457	429226	8.98%	1.023%
48	14.704	1968	1975	1979	rBV7	7946	20844	0.44%	0.050%
49	14.769	1979	1986	1992	rVV3	41089	84847	1.78%	0.202%
50	14.828	1992	1996	1999	rVV5	7591	11096	0.23%	0.026%
51	14.887	2003	2006	2012	rVB4	15930	24627	0.52%	0.059%
52	14.957	2012	2018	2024	rBV7	9615	19431	0.41%	0.046%
53	15.222	2059	2063	2068	rVB	56027	70049	1.47%	0.167%
54	15.345	2077	2084	2090	rBV	1401536	1953393	40.88%	4.657%
55	15.404	2090	2094	2098	rVV	21449	30249	0.63%	0.072%
56	15.457	2098	2103	2109	rVV	190135	266603	5.58%	0.636%
57	15.510	2109	2112	2118	rVB2	30502	43336	0.91%	0.103%
58	15.586	2121	2125	2128	rVV	15492	21500	0.45%	0.051%
59	15.628	2128	2132	2138	rVB	30914	46418	0.97%	0.111%
60	15.698	2141	2144	2152	rVV4	8529	19115	0.40%	0.046%
61	15.775	2152	2157	2163	rVV	222328	285813	5.98%	0.681%
62	15.839	2164	2168	2176	rVB4	18163	34709	0.73%	0.083%
63	16.086	2205	2210	2212	rBV	61107	88531	1.85%	0.211%
64	16.110	2212	2214	2224	rVB3	45248	66746	1.40%	0.159%
65	16.422	2258	2267	2271	rBV9	13566	35987	0.75%	0.086%
66	16.551	2285	2289	2292	rBV5	9138	13756	0.29%	0.033%
67	16.645	2301	2305	2306	rBV4	9829	13810	0.29%	0.033%
68	16.828	2334	2336	2338	rBV3	11780	12617	0.26%	0.030%
69	16.875	2341	2344	2348	rVV2	90607	113931	2.38%	0.272%
70	16.928	2348	2353	2364	rVB2	49212	98630	2.06%	0.235%
71	17.098	2376	2382	2386	rBV	1472112	2082032	43.57%	4.964%
72	17.139	2386	2389	2393	rVV	198017	267939	5.61%	0.639%
73	17.198	2393	2399	2411	rVV2	1435114	2208550	46.22%	5.265%
74	17.292	2411	2415	2420	rVB8	22476	46937	0.98%	0.112%
75	17.475	2443	2446	2448	rBV4	17131	21998	0.46%	0.052%
76	17.616	2466	2470	2475	rBV	104075	124684	2.61%	0.297%

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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-BN061821MA.M
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77	17.786	2496	2499	2504	rVB3	29893	36055	0.75%	0.086%
78	17.975	2528	2531	2533	rBV3	25112	29101	0.61%	0.069%
79	18.016	2533	2538	2543	rVB3	53316	109296	2.29%	0.261%
80	18.169	2560	2564	2568	rVB3	36786	51037	1.07%	0.122%
81	18.245	2575	2577	2584	rVB7	24137	38504	0.81%	0.092%
82	18.310	2584	2588	2594	rBV	115577	174494	3.65%	0.416%
83	18.428	2603	2608	2612	rBV	218160	292402	6.12%	0.697%
84	18.775	2664	2667	2670	rBV2	35374	44292	0.93%	0.106%
85	18.945	2692	2696	2705	rBV	126127	221404	4.63%	0.528%
86	19.163	2728	2733	2739	rVB	300290	421385	8.82%	1.005%
87	19.492	2784	2789	2793	rBV2	1773732	2669519	55.87%	6.364%
88	20.063	2883	2886	2890	rBV	81455	85217	1.78%	0.203%
89	20.516	2958	2963	2968	rBV	4181346	4778138	100.00%	11.391%
90	21.021	3046	3049	3063	rVB4	198538	433765	9.08%	1.034%
91	21.221	3079	3083	3088	rVB	314871	360639	7.55%	0.860%
92	21.292	3089	3095	3099	rBV	1846485	2488721	52.09%	5.933%
93	22.157	3238	3242	3248	rVB	205837	281653	5.89%	0.671%
94	22.874	3359	3364	3375	rVB2	253098	547426	11.46%	1.305%
95	23.398	3446	3453	3458	rBV2	1233604	2278471	47.69%	5.432%
96	23.445	3458	3461	3469	rVB2	309551	500445	10.47%	1.193%
97	23.539	3470	3477	3484	rBV2	1179021	2273534	47.58%	5.420%
98	24.098	3567	3572	3579	rVB	236335	458292	9.59%	1.093%
99	24.845	3695	3699	3709	rVB	213274	464034	9.71%	1.106%
100	25.727	3845	3849	3857	rVB3	100397	203823	4.27%	0.486%

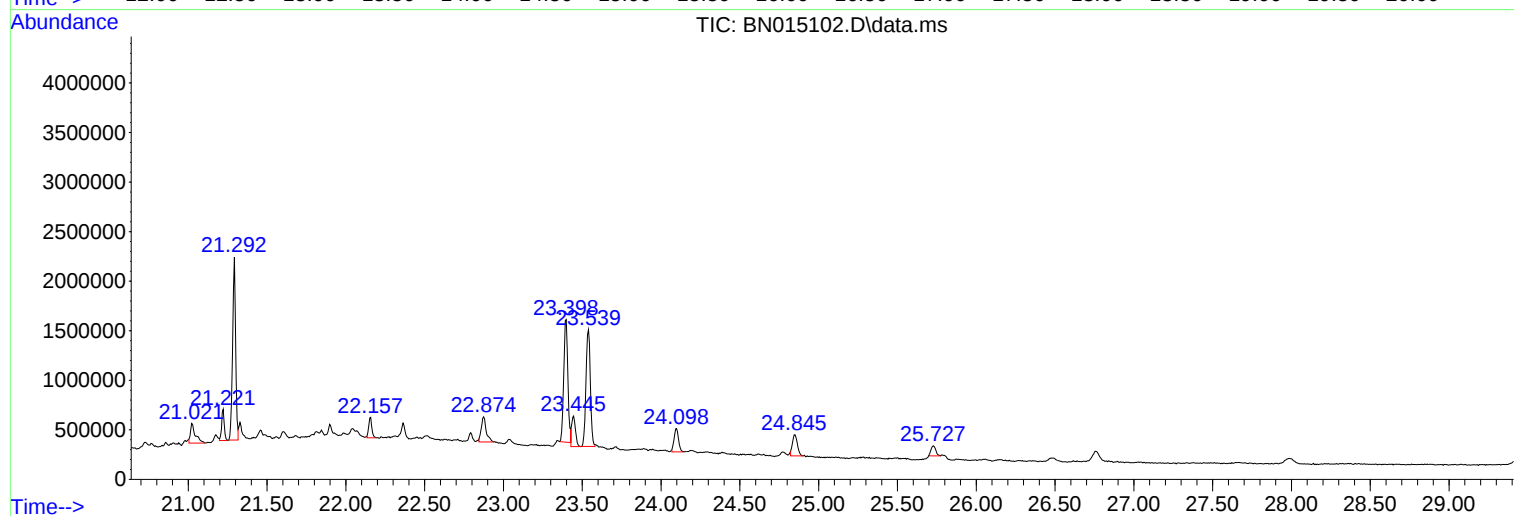
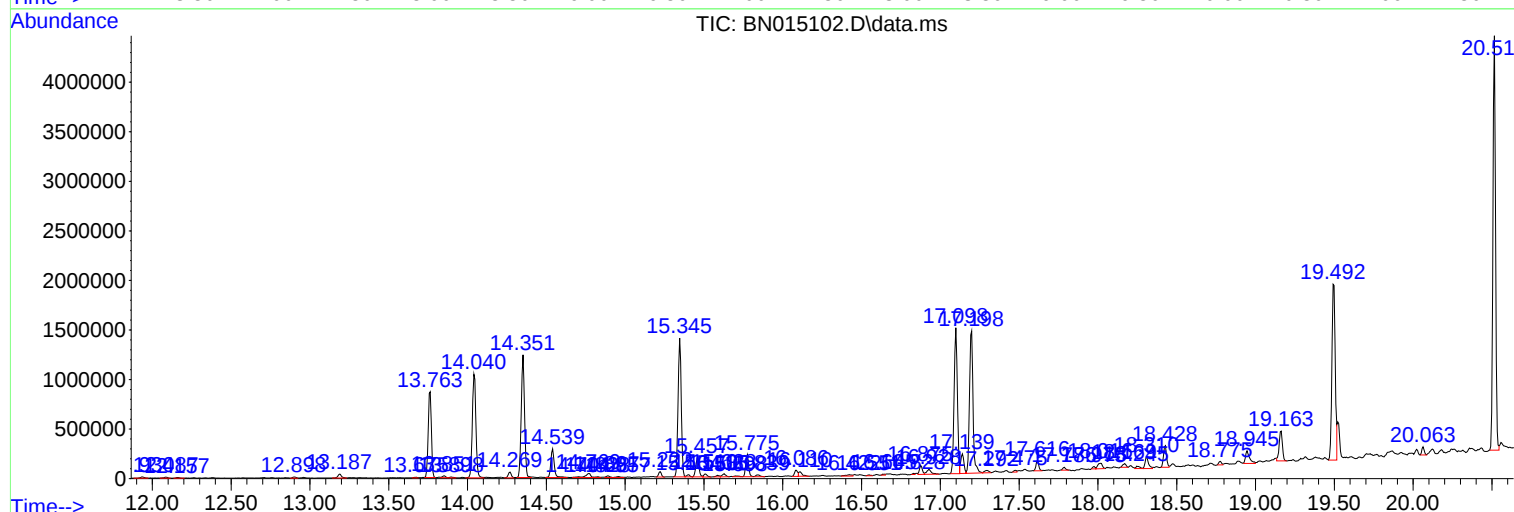
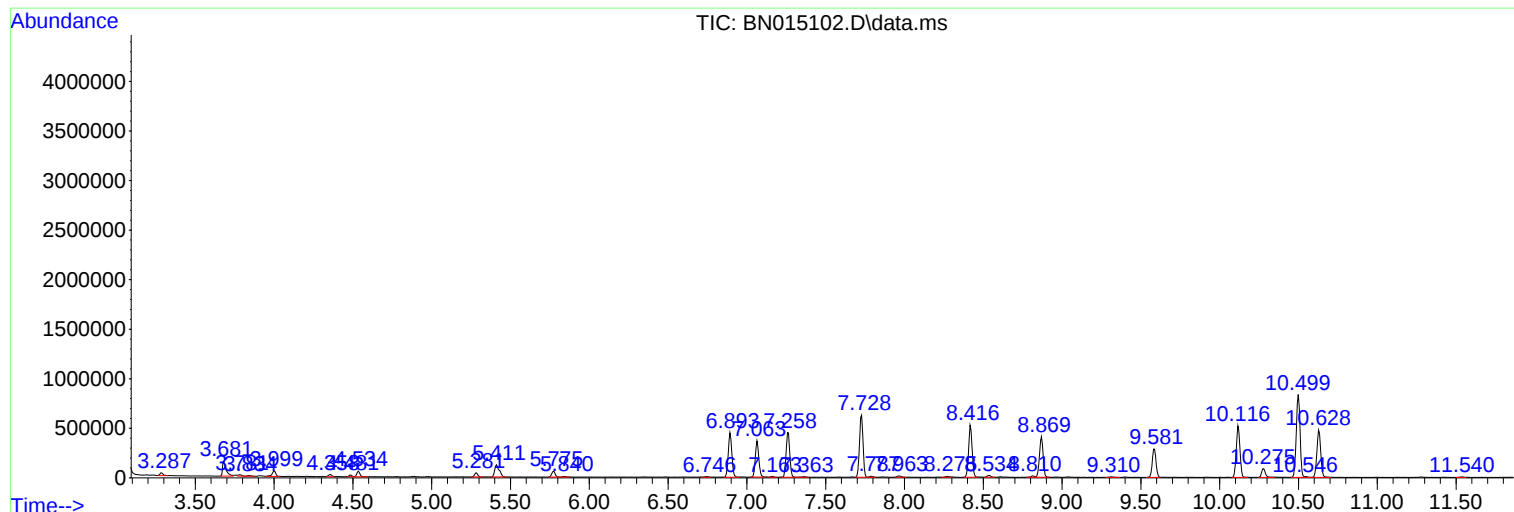
Sum of corrected areas: 41945287

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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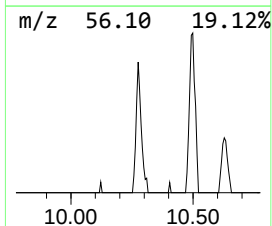
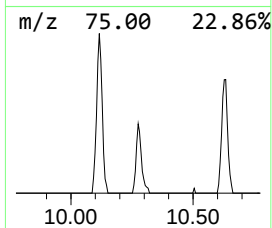
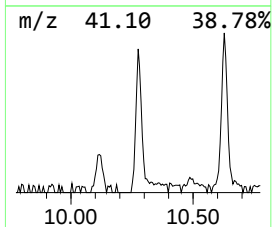
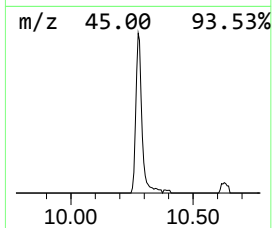
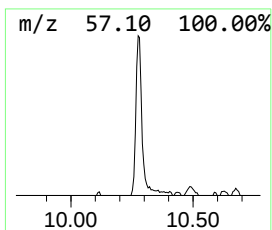
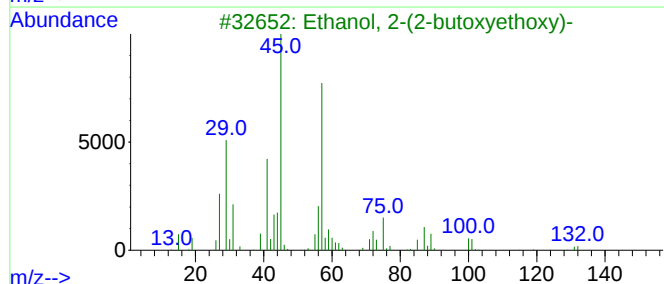
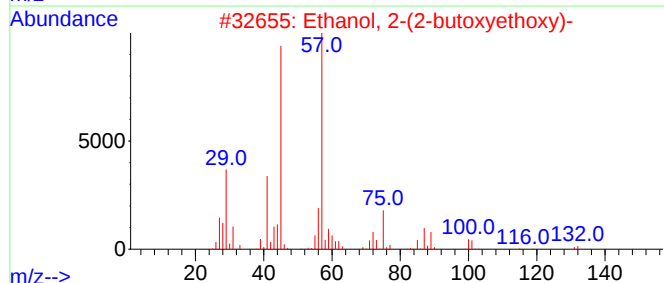
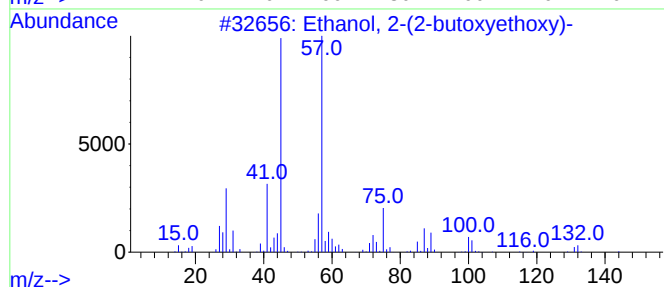
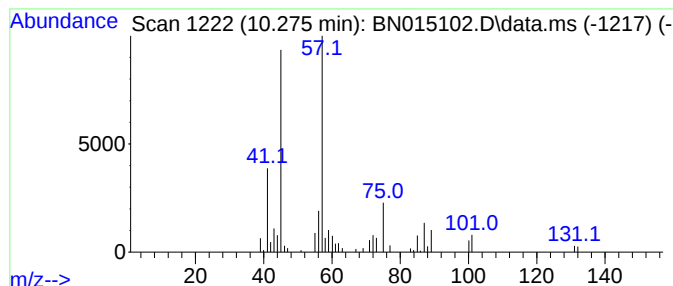
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.275	2.11 ng/ul	147209	Naphthalene-d8	10.499

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	78
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64



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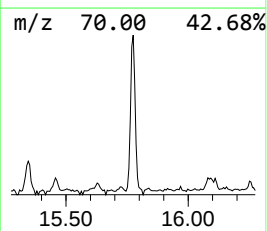
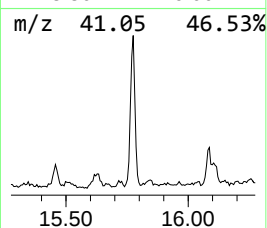
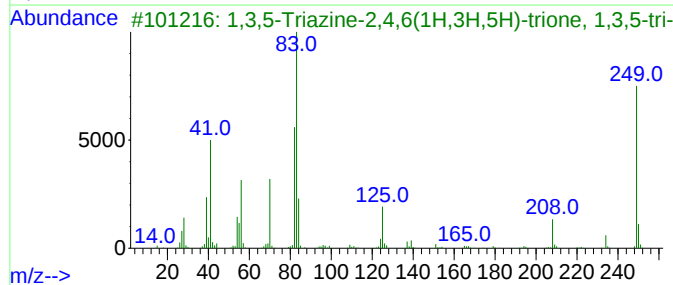
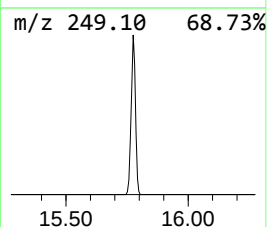
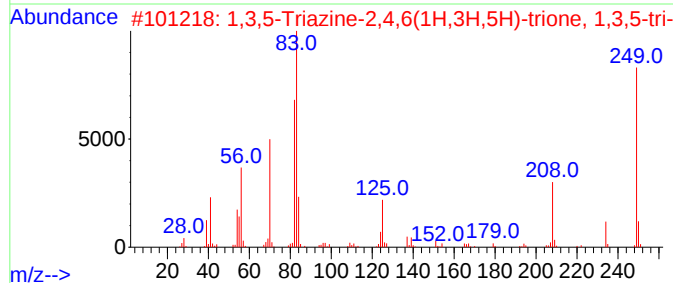
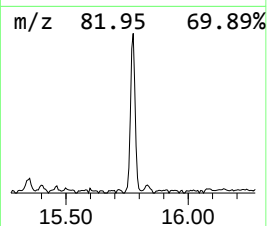
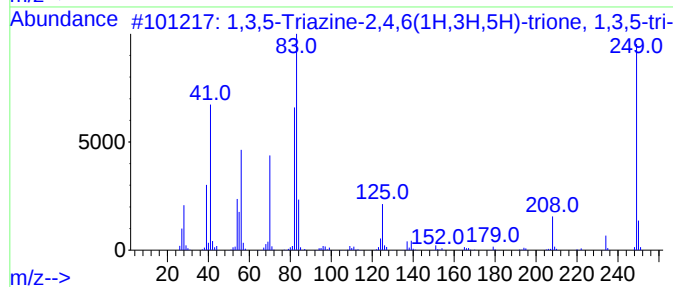
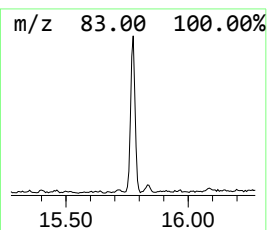
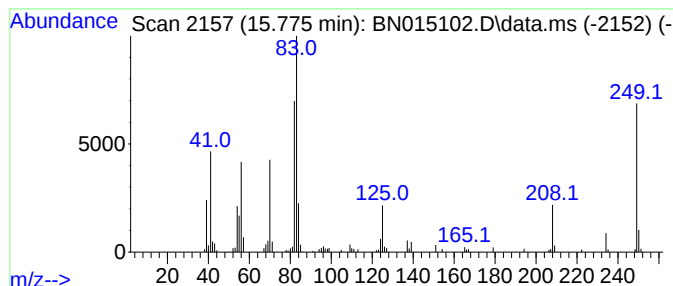
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,3,5-Triazine-2,4,6(1H,3H,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	2.75 ng/ul	285813	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	99
2			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	99
3			1,3,5-Triazine-2,4,6(1H,3H,5H)-t...	249	C12H15N3O3	001025-15-6	98
4			N-Carboxy-1-[4-chlorophenyl]-2-[...	295	C15H18ClNO3	058158-38-6	37
5			n-Pentadecylcyclohexane	294	C21H42	006006-95-7	16



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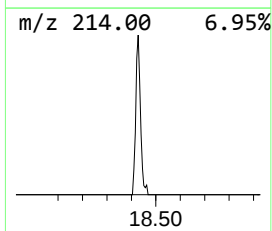
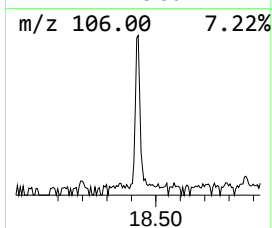
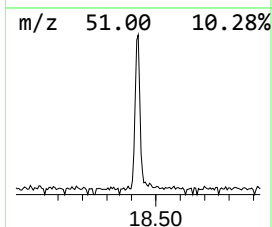
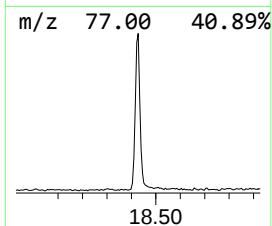
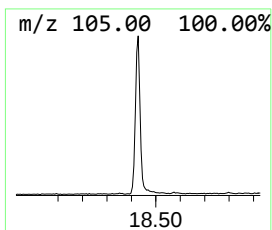
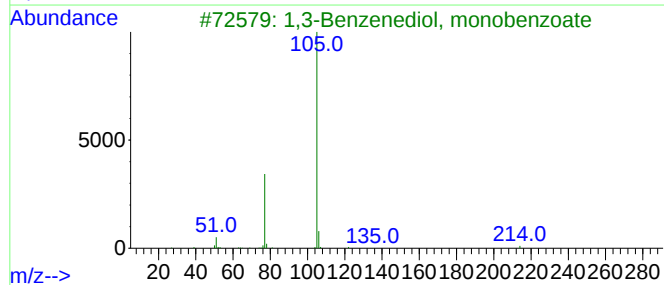
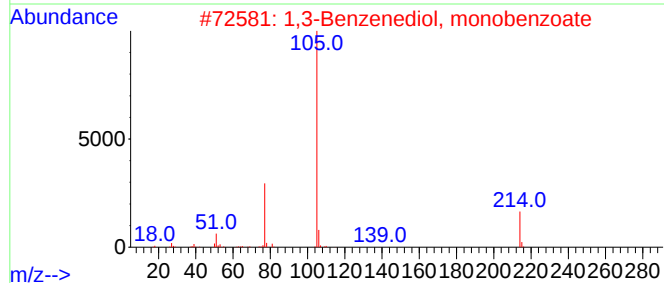
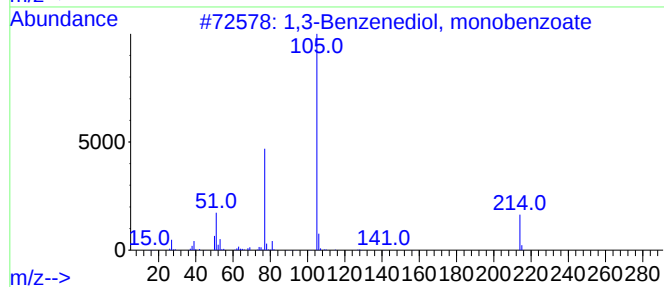
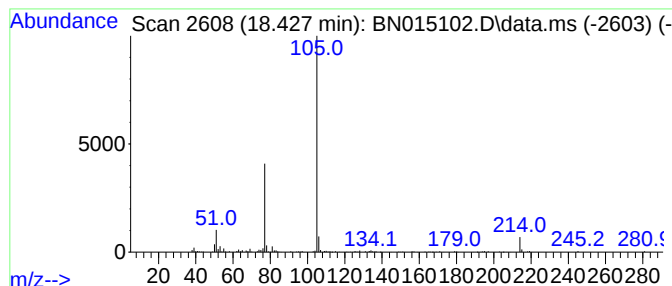
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1,3-Benzenediol, monobenzoate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.427	2.81 ng/ul	292402	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	90
2			1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	87
3			1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	87
4			1,4-Benzenediol, monobenzoate	214	C13H10O3	002444-19-1	80
5			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015102.D
Acq On : 26 Jun 2021 06:15
Operator : CG/JU
Sample : M2704-05
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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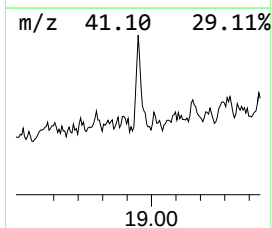
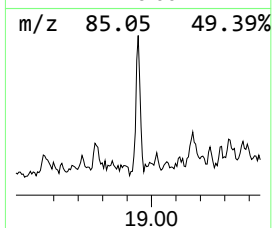
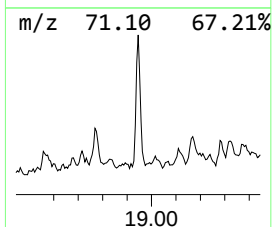
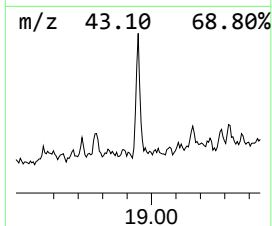
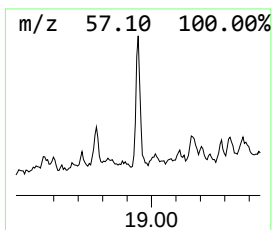
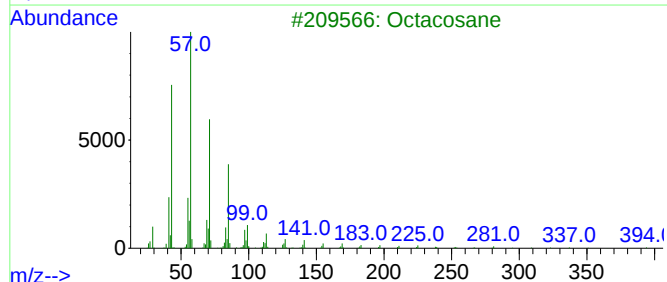
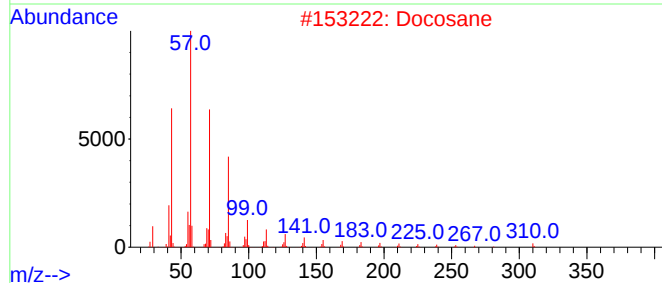
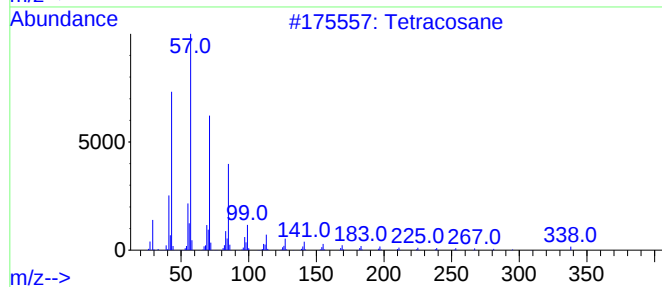
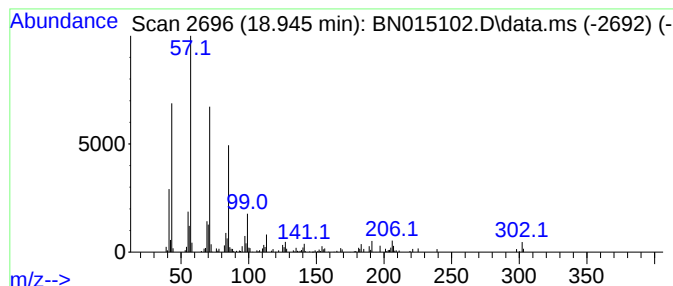
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.945	2.13 ng/ul	221404	Phenanthrene-d10	17.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	83
2		Docosane	310	C22H46	000629-97-0	80
3		Octacosane	394	C28H58	000630-02-4	74
4		Pentadecane	212	C15H32	000629-62-9	74
5		Heneicosane	296	C21H44	000629-94-7	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015102.D
Acq On : 26 Jun 2021 06:15
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Sample : M2704-05
Misc :
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Instrument :
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ClientSampleId :
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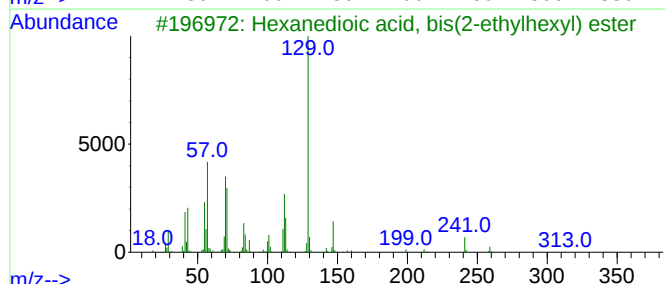
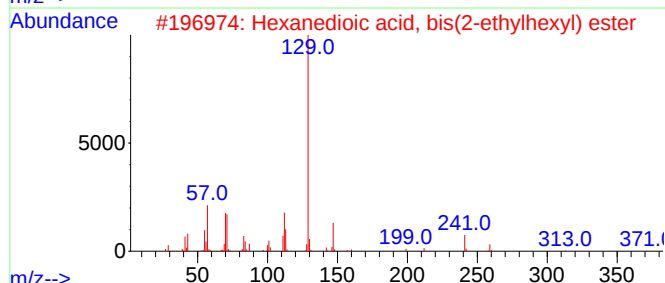
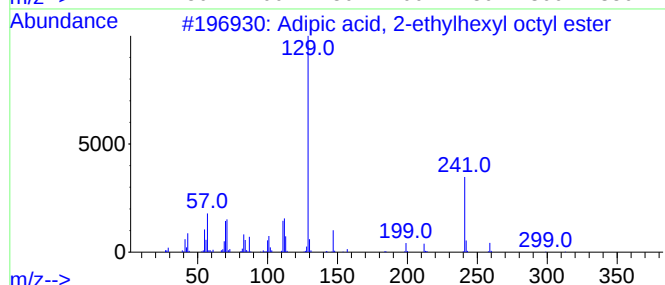
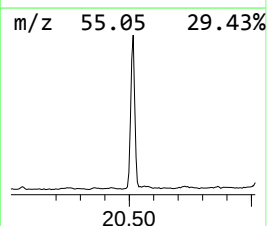
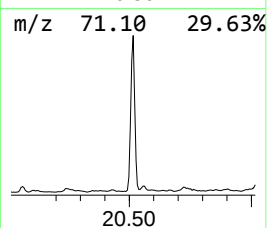
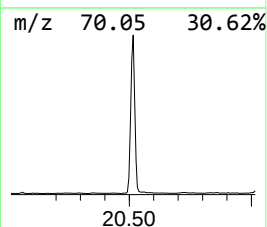
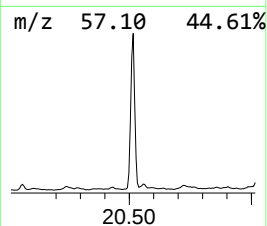
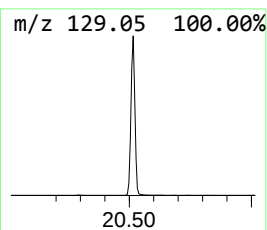
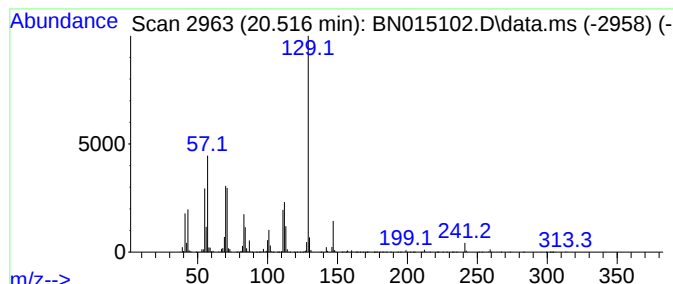
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Adipic acid, 2-ethylhexyl o... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	38.40 ng/ul	4778140	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	90
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
4			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	87
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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Acq On : 26 Jun 2021 06:15
Operator : CG/JU
Sample : M2704-05
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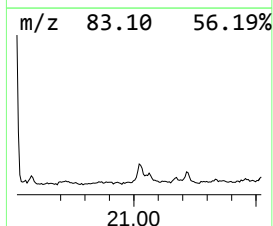
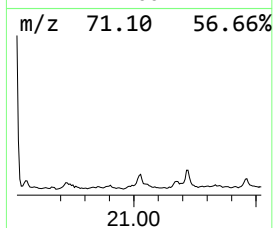
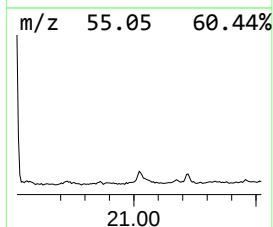
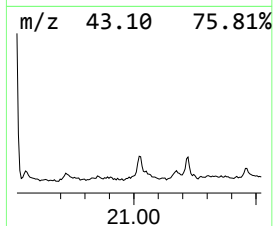
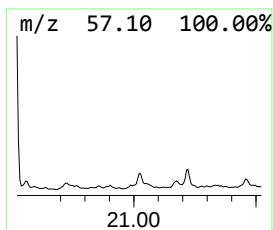
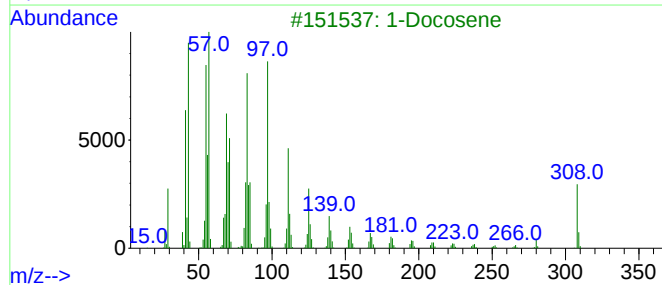
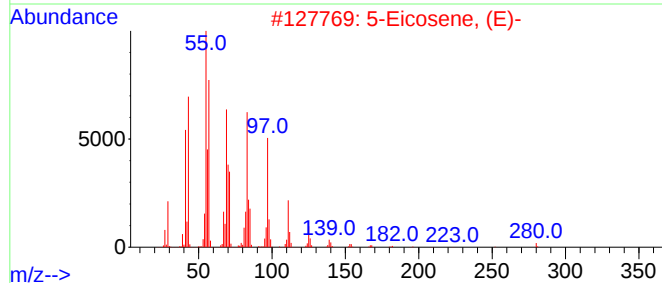
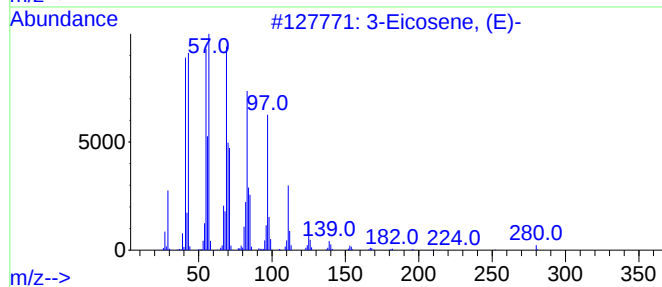
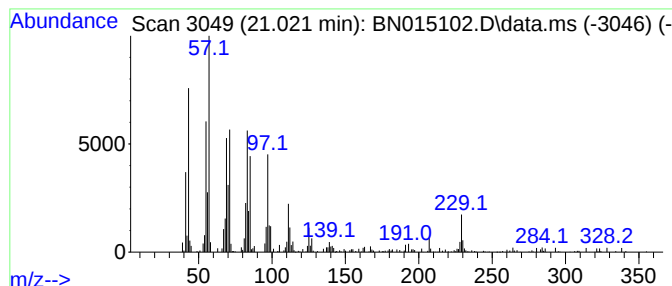
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.022	3.49 ng/ul	433765	Chrysene-d12	21.292	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	98
2	5-Eicosene, (E)-	280	C20H40	074685-30-6	92
3	1-Docosene	308	C22H44	001599-67-3	91
4	9-Eicosene, (E)-	280	C20H40	074685-29-3	83
5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015102.D
Acq On : 26 Jun 2021 06:15
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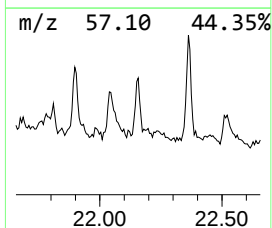
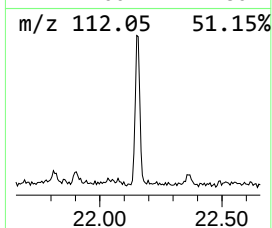
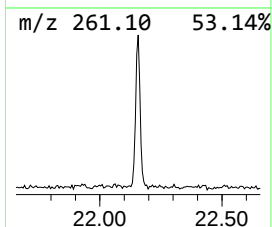
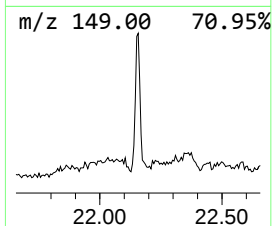
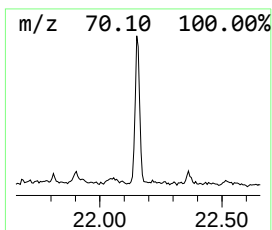
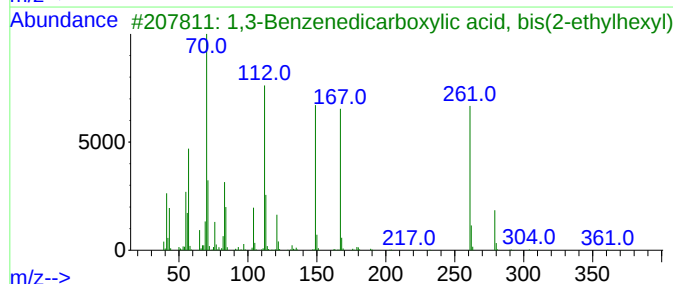
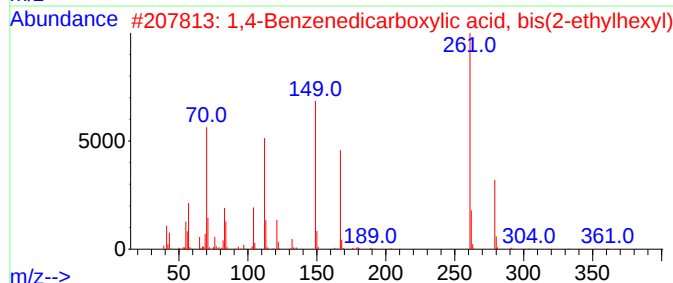
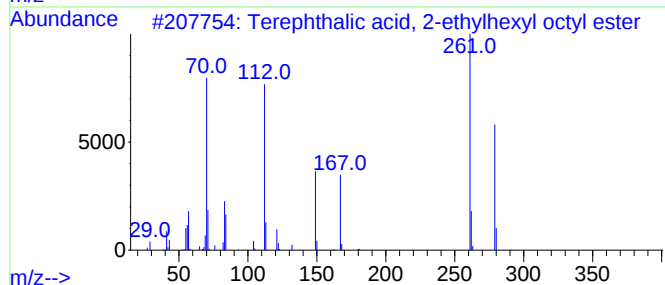
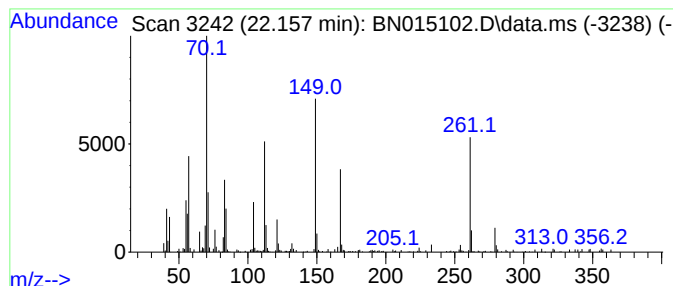
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Terephthalic acid, 2-ethylh... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.157	2.26 ng/ul	281653	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Terephthalic acid, 2-ethylhexyl ...	390	C24H38O4	1000324-00-5	58
2			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	53
3			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	50
4			Terephthalic acid, octyl 2-penty...	348	C21H32O4	1000323-55-2	49
5			Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015102.D
Acq On : 26 Jun 2021 06:15
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Sample : M2704-05
Misc :
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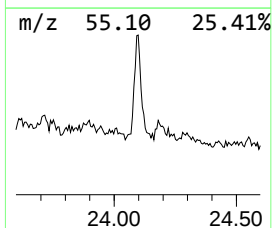
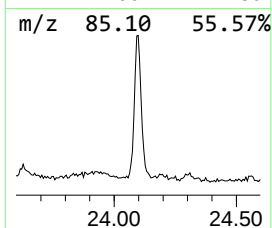
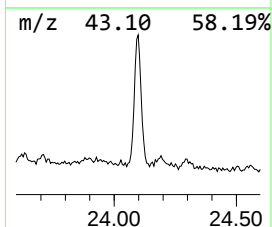
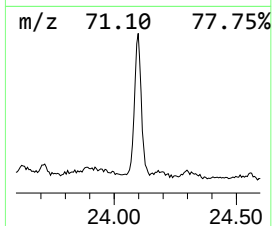
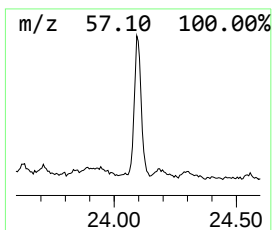
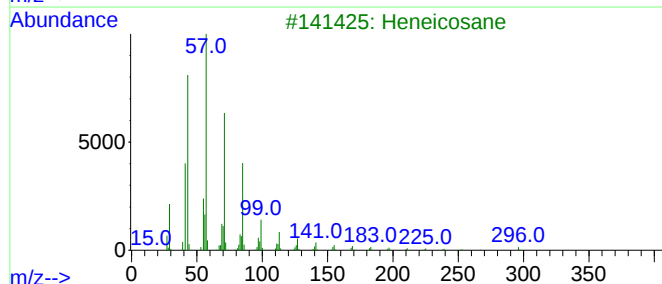
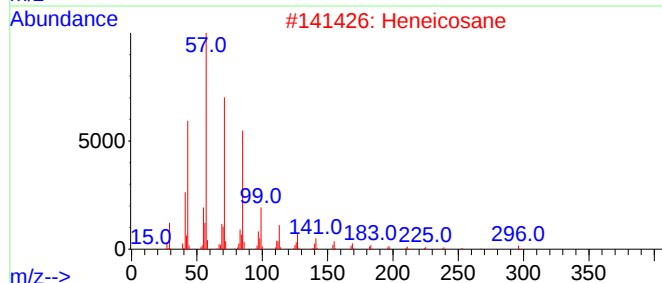
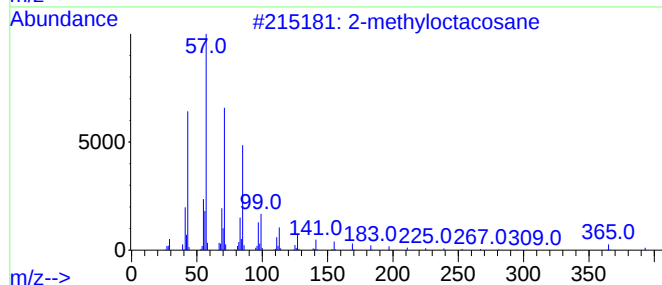
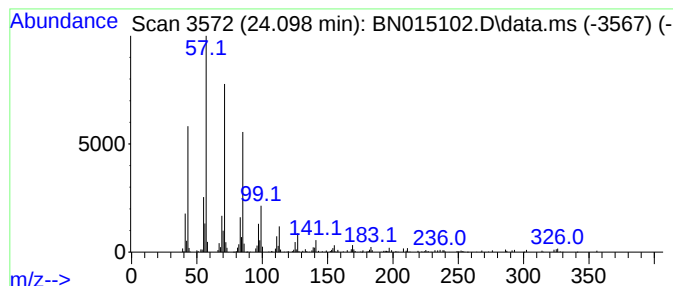
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.098	4.03 ng/ul	458292	Perylene-d12	23.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octacosane	394	C28H58	000630-02-4	90
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015102.D
Acq On : 26 Jun 2021 06:15
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Sample : M2704-05
Misc :
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Instrument :
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ClientSampleId :
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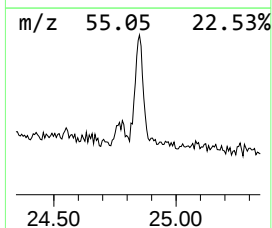
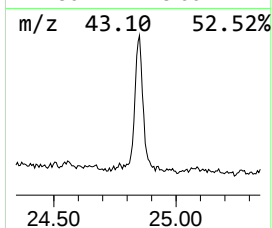
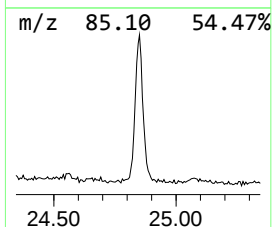
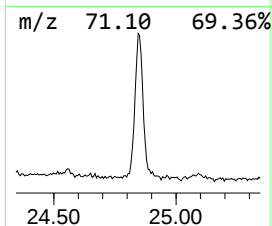
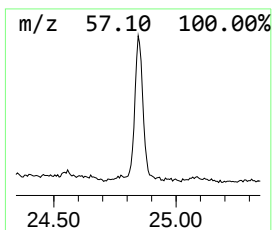
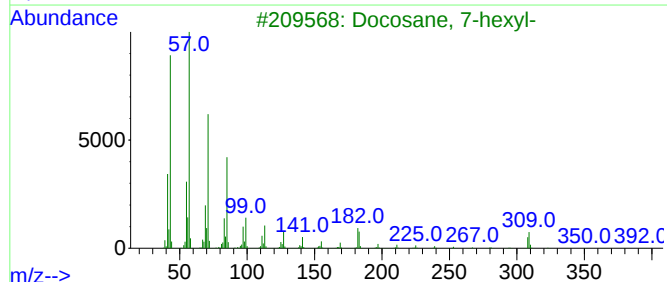
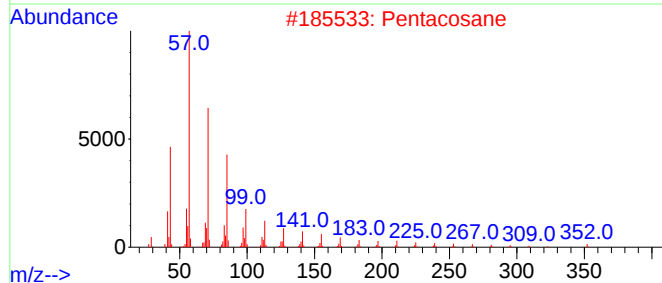
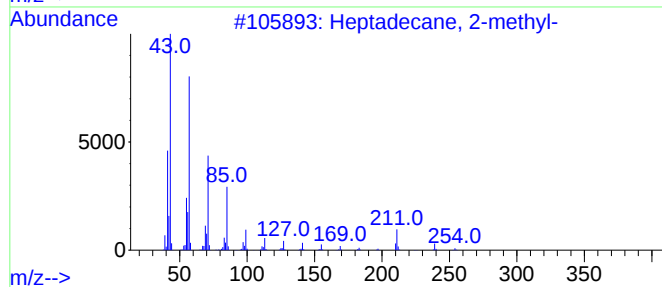
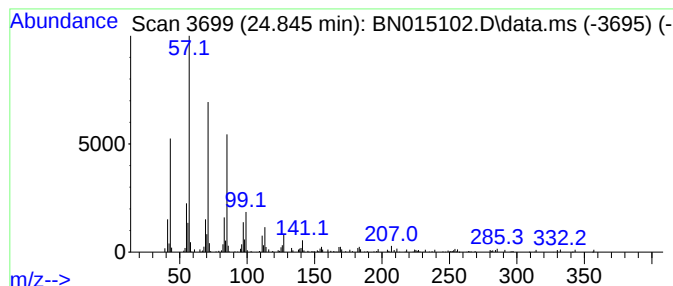
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.845	4.08 ng/ul	464034	Perylene-d12	23.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	90
2		Pentacosane	352	C25H52	000629-99-2	87
3		Docosane, 7-hexyl-	394	C28H58	055373-86-9	87
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
5		Dodecane, 2-methyl-	184	C13H28	001560-97-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
 Data File : BN015102.D
 Acq On : 26 Jun 2021 06:15
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 Sample : M2704-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.275	2.1	ng/ul	147209	2	10.499	1395770	20.0
1,3,5-Triazine-...	15.775	2.8	ng/ul	285813	4	17.098	2082030	20.0
1,3-Benzenediol...	18.427	2.8	ng/ul	292402	4	17.098	2082030	20.0
(DEL) Alkane: S...	18.945	2.1	ng/ul	221404	4	17.098	2082030	20.0
Adipic acid, 2-...	20.516	38.4	ng/ul	4778140	5	21.292	2488720	20.0
(DEL) Alkane: S...	21.022	3.5	ng/ul	433765	5	21.292	2488720	20.0
Terephthalic ac...	22.157	2.3	ng/ul	281653	5	21.292	2488720	20.0
(DEL) Alkane: S...	24.098	4.0	ng/ul	458292	6	23.539	2273530	20.0
(DEL) Alkane: S...	24.845	4.1	ng/ul	464034	6	23.539	2273530	20.0