

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
 Data File : BN010393.D
 Acq On : 02 Apr 2020 15:32
 Operator : CG/JU
 Sample : L2065-10DL2 200X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBF06DL2

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.899	15	19	26	rBV3	31372	47998	0.40%	0.074%
2	3.917	188	192	198	rVB3	16457	29437	0.24%	0.045%
3	5.187	404	408	412	rBV3	9181	13021	0.11%	0.020%
4	5.311	425	429	435	rBV4	15272	24437	0.20%	0.038%
5	6.334	598	603	609	rBV2	62327	101490	0.84%	0.156%
6	6.628	649	653	658	rVB5	16573	23773	0.20%	0.037%
7	6.728	665	670	674	rBV2	25937	40751	0.34%	0.063%
8	6.781	675	679	682	rVV4	12981	20988	0.17%	0.032%
9	6.822	682	686	688	rVV4	13981	18554	0.15%	0.029%
10	6.858	688	692	698	rVB	34417	49110	0.41%	0.075%
11	7.075	722	729	732	rVB4	24790	39831	0.33%	0.061%
12	7.275	757	763	769	rBV2	81093	130334	1.08%	0.200%
13	7.340	769	774	782	rVB3	25585	49654	0.41%	0.076%
14	7.611	813	820	828	rBV	1390187	2155611	17.83%	3.311%
15	7.758	837	845	850	rBV2	133001	212652	1.76%	0.327%
16	7.846	855	860	865	rBV4	28103	43595	0.36%	0.067%
17	7.981	878	883	890	rVB	172742	268816	2.22%	0.413%
18	8.140	902	910	917	rBV	374229	594267	4.91%	0.913%
19	8.252	924	929	936	rBV5	9891	18271	0.15%	0.028%
20	8.716	999	1008	1009	rBV6	11474	18591	0.15%	0.029%
21	8.734	1009	1011	1016	rVV5	16327	24436	0.20%	0.038%
22	8.787	1016	1020	1024	rVV2	14244	23785	0.20%	0.037%
23	8.840	1024	1029	1033	rVV2	23750	34270	0.28%	0.053%
24	8.952	1042	1048	1055	rVB3	26824	42460	0.35%	0.065%
25	9.075	1061	1069	1074	rBV3	52960	114740	0.95%	0.176%
26	9.134	1074	1079	1084	rVB3	18073	29583	0.24%	0.045%
27	9.228	1090	1095	1101	rVB8	11938	19980	0.17%	0.031%
28	9.281	1101	1104	1109	rBV6	10299	17065	0.14%	0.026%
29	9.628	1155	1163	1171	rBV5	24836	58085	0.48%	0.089%
30	9.752	1178	1184	1189	rBV	835048	1356188	11.22%	2.083%
31	9.793	1189	1191	1201	rVV5	145638	269007	2.22%	0.413%
32	9.881	1201	1206	1210	rVV4	35264	69247	0.57%	0.106%
33	9.928	1210	1214	1218	rVV6	24153	41907	0.35%	0.064%
34	10.005	1221	1227	1233	rVB2	33822	61253	0.51%	0.094%

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 ClientSampleId :
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Integration Parameters: LSCINT.P

Integrator: RTE
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 Title : SVOA CALIBRATION

35	10.246	1265	1268	1272	rBV6	8908	12193	0.10%	0.019%
36	10.369	1282	1289	1293	rBV	2011802	3384020	27.99%	5.198%
37	10.422	1293	1298	1308	rVV	7259226	12090987	100.00%	18.573%
38	10.534	1310	1317	1328	rVB	205942	369805	3.06%	0.568%
39	11.199	1425	1430	1436	rBV2	32493	58530	0.48%	0.090%
40	11.928	1549	1554	1561	rVB2	24863	46875	0.39%	0.072%
41	12.034	1566	1572	1579	rBV	1161214	1804631	14.93%	2.772%
42	12.151	1586	1592	1597	rVB2	33163	55472	0.46%	0.085%
43	12.257	1599	1610	1616	rBV	576171	914621	7.56%	1.405%
44	13.063	1742	1747	1754	rVB	184450	263516	2.18%	0.405%
45	13.263	1776	1781	1788	rBV	28626	49657	0.41%	0.076%
46	13.398	1796	1804	1805	rBV4	32487	44918	0.37%	0.069%
47	13.551	1823	1830	1835	rBV2	53236	91714	0.76%	0.141%
48	13.610	1835	1840	1843	rVV3	31520	49001	0.41%	0.075%
49	13.651	1843	1847	1852	rVB	29154	38487	0.32%	0.059%
50	13.793	1866	1871	1875	rBV3	18238	32303	0.27%	0.050%
51	13.922	1887	1893	1896	rBV2	23894	41374	0.34%	0.064%
52	13.963	1896	1900	1907	rVB6	18372	30635	0.25%	0.047%
53	14.234	1939	1946	1952	rBV	3588192	5188128	42.91%	7.970%
54	14.298	1952	1957	1963	rVV	744202	1056559	8.74%	1.623%
55	14.363	1963	1968	1974	rVB	117254	187256	1.55%	0.288%
56	14.551	1997	2000	2005	rVB4	11900	16258	0.13%	0.025%
57	14.634	2008	2014	2021	rBV2	394500	590000	4.88%	0.906%
58	14.728	2025	2030	2032	rBV5	10828	14384	0.12%	0.022%
59	15.228	2111	2115	2120	rBV	37445	54647	0.45%	0.084%
60	15.287	2120	2125	2131	rVB2	268544	387924	3.21%	0.596%
61	15.451	2149	2153	2159	rVB6	13895	21993	0.18%	0.034%
62	15.545	2164	2169	2172	rBV6	9930	16915	0.14%	0.026%
63	15.587	2172	2176	2184	rVB5	16688	27154	0.22%	0.042%
64	15.734	2195	2201	2207	rBV5	19487	36007	0.30%	0.055%
65	15.957	2234	2239	2247	rBV3	28399	42855	0.35%	0.066%
66	16.087	2258	2261	2267	rVB6	10258	14986	0.12%	0.023%
67	16.175	2272	2276	2281	rVV5	15877	23766	0.20%	0.037%
68	16.257	2284	2290	2294	rBV8	13130	27823	0.23%	0.043%
69	16.769	2373	2377	2379	rBV2	22698	30003	0.25%	0.046%
70	16.798	2379	2382	2387	rVB3	21528	31532	0.26%	0.048%
71	16.981	2405	2413	2418	rBV	4738618	6739938	55.74%	10.353%

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72	17.022	2418	2420	2425	rVV	317004	372333	3.08%	0.572%
73	17.081	2425	2430	2440	rVV3	117589	241906	2.00%	0.372%
74	17.381	2472	2481	2488	rBV	370182	557933	4.61%	0.857%
75	17.481	2495	2498	2503	rVB6	11478	17438	0.14%	0.027%
76	17.645	2523	2526	2529	rBV5	9195	13276	0.11%	0.020%
77	17.904	2565	2570	2576	rBV5	50637	84141	0.70%	0.129%
78	18.133	2607	2609	2611	rBV3	12481	12224	0.10%	0.019%
79	18.204	2617	2621	2623	rBV4	18478	20595	0.17%	0.032%
80	18.728	2706	2710	2712	rBV5	8611	12241	0.10%	0.019%
81	19.045	2759	2764	2771	rBV3	33780	67998	0.56%	0.104%
82	19.198	2787	2790	2792	rBV4	13630	17308	0.14%	0.027%
83	19.351	2813	2816	2817	rBV3	17124	19882	0.16%	0.031%
84	19.375	2817	2820	2825	rVV2	61877	86333	0.71%	0.133%
85	21.180	3122	3127	3134	rBV	7791467	9097977	75.25%	13.976%
86	22.251	3306	3309	3314	rVB2	296731	351077	2.90%	0.539%
87	22.745	3389	3393	3398	rBV	503900	651493	5.39%	1.001%
88	23.292	3482	3486	3490	rBV	694134	1001474	8.28%	1.538%
89	23.357	3490	3497	3505	rVB	5375455	9798885	81.04%	15.052%
90	23.916	3587	3592	3602	rVB	757674	1335750	11.05%	2.052%
91	24.633	3708	3714	3721	rVB	705434	1388737	11.49%	2.133%

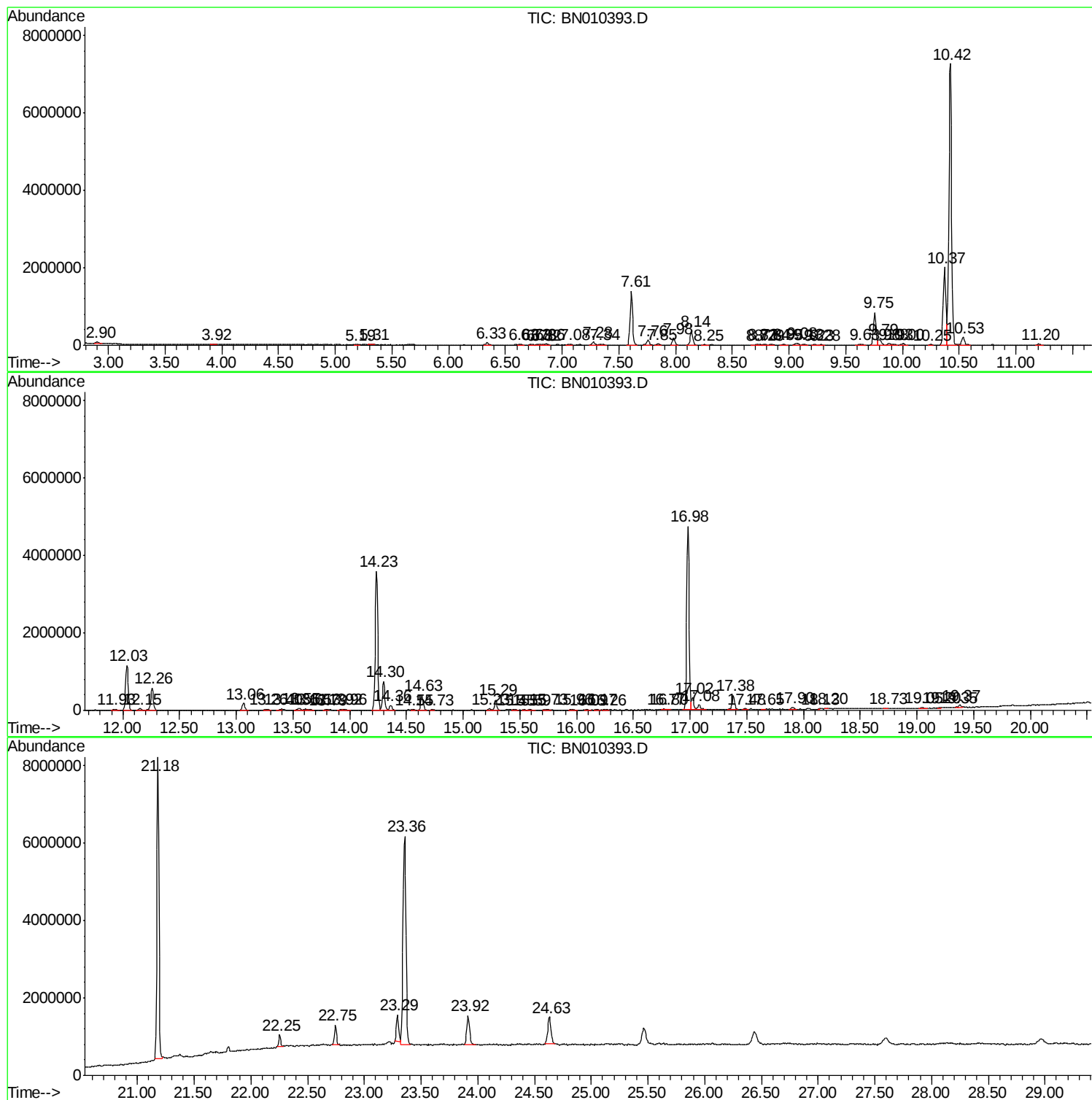
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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



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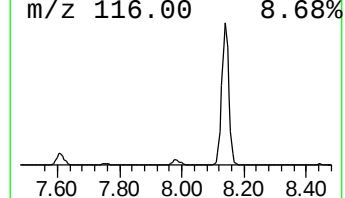
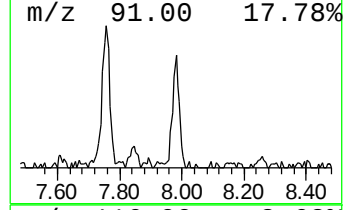
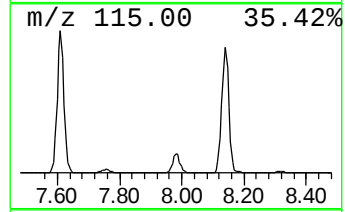
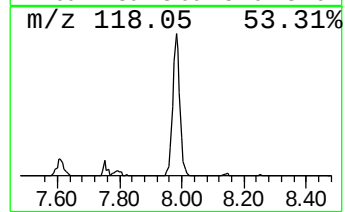
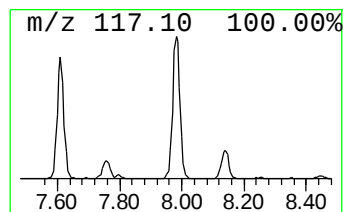
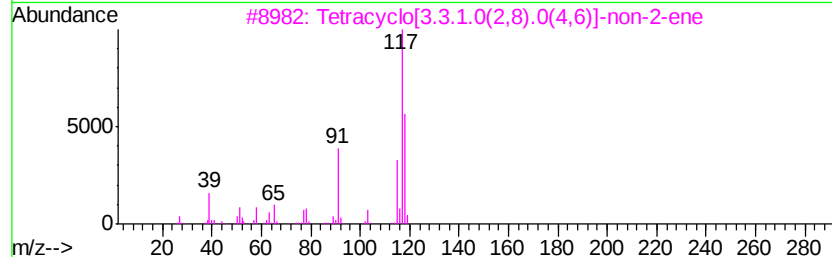
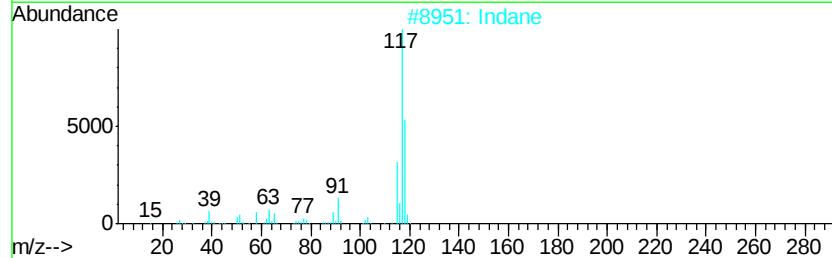
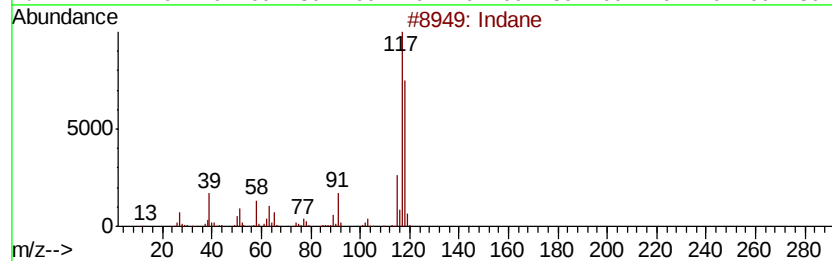
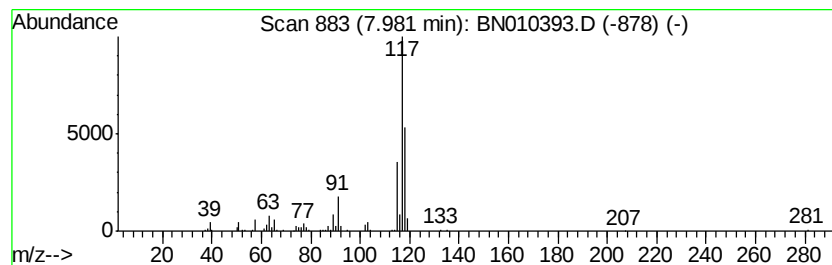
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.98	2.49 ng/ul	268816	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	87
2		Indane	118	C9H10	000496-11-7	87
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	83
4		Indane	118	C9H10	000496-11-7	83
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	81



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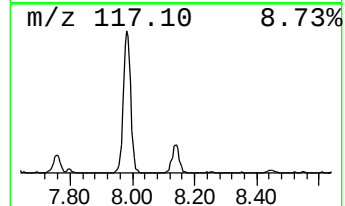
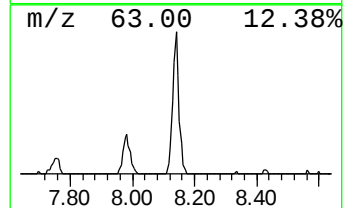
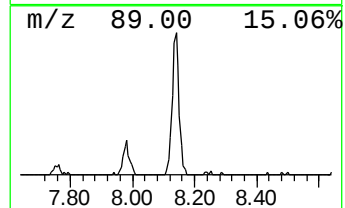
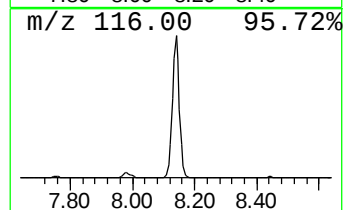
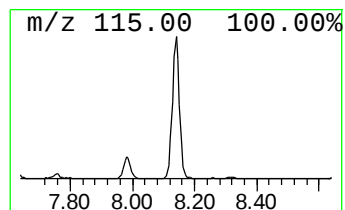
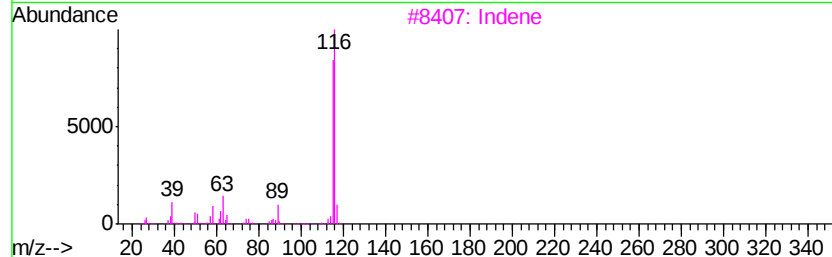
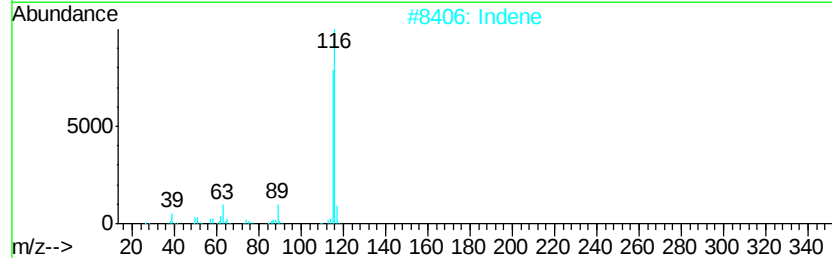
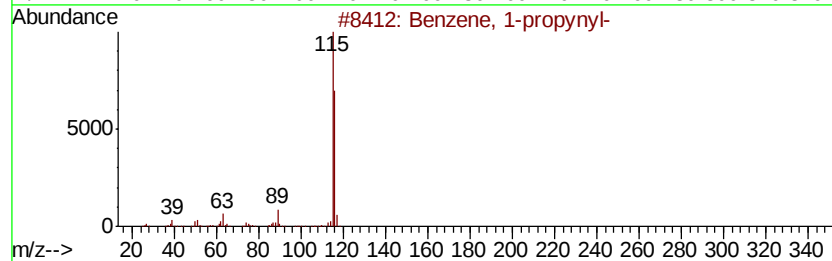
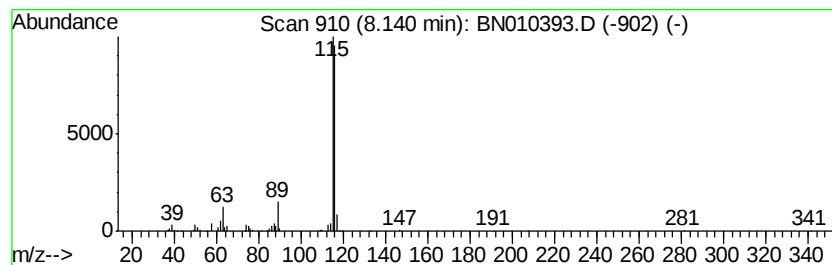
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1-propynyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.14	5.51 ng/ul	594267	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
2		Indene	116	C9H8	000095-13-6	94
3		Indene	116	C9H8	000095-13-6	91
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5		3-Methylphenylacetylene	116	C9H8	000766-82-5	91



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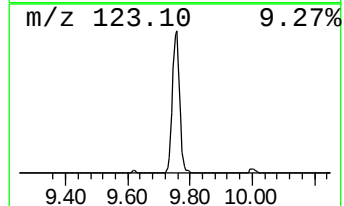
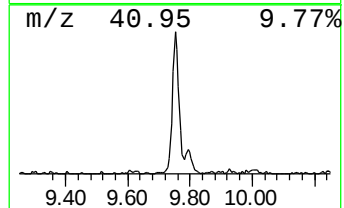
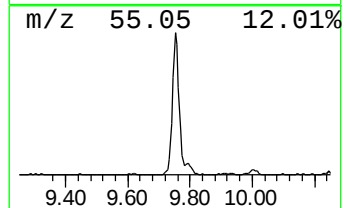
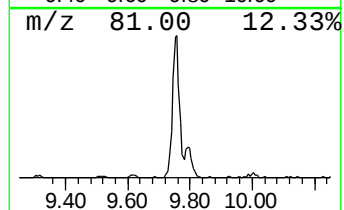
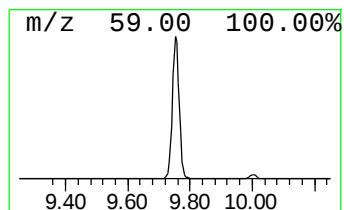
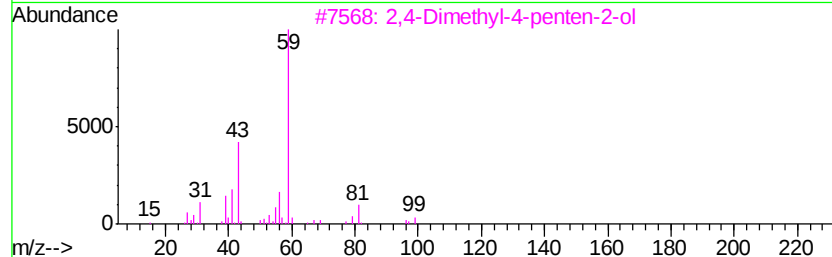
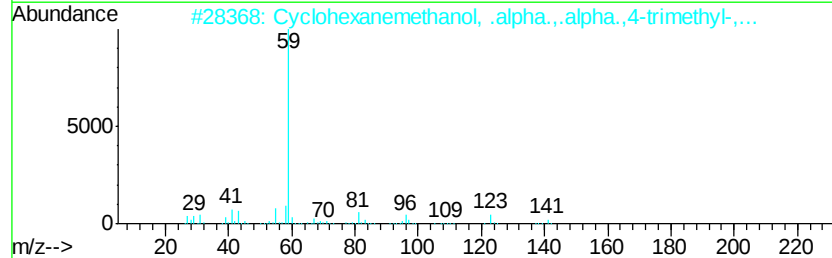
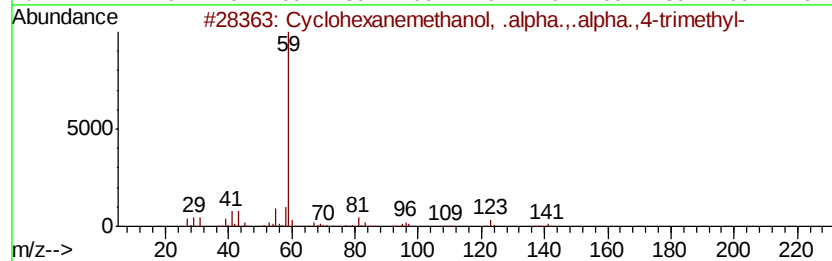
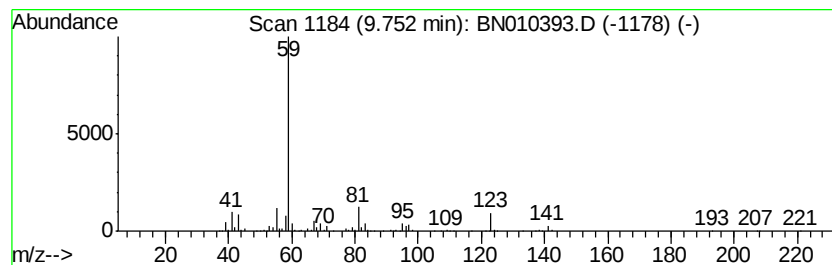
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclohexanemethanol, .alpha... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.75	8.02 ng/ul	1356190	Naphthalene-d8	10.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanemethanol, .alpha...al...	156	C10H20O	000498-81-7	74
2		Cyclohexanemethanol, .alpha...al...	156	C10H20O	007322-63-6	64
3		2,4-Dimethyl-4-penten-2-ol	114	C7H14O	019781-53-4	64
4		Cyclohexanemethanol, .alpha...al...	156	C10H20O	005114-00-1	64
5		1-Cyclohexylethanol, methyl ether	142	C9H18O	1000365-12-8	53



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DBF06DL2

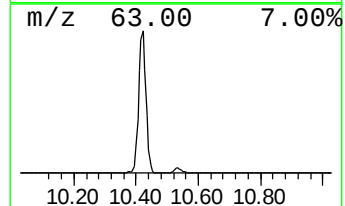
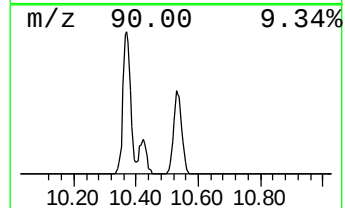
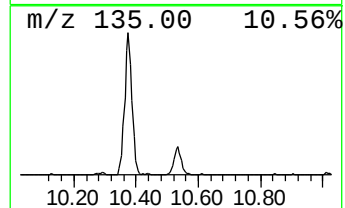
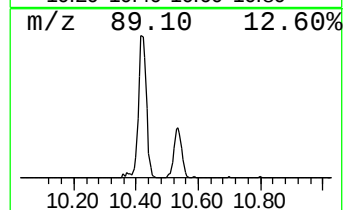
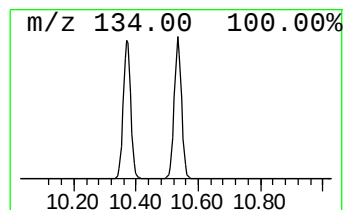
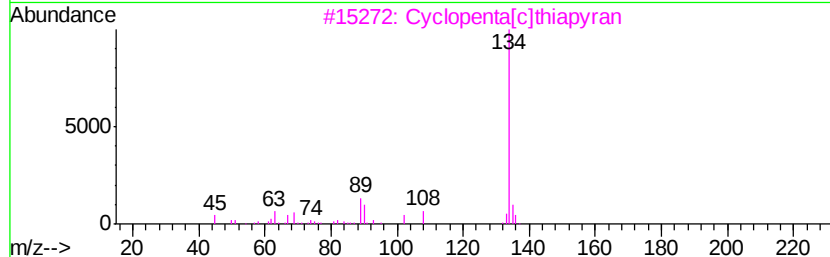
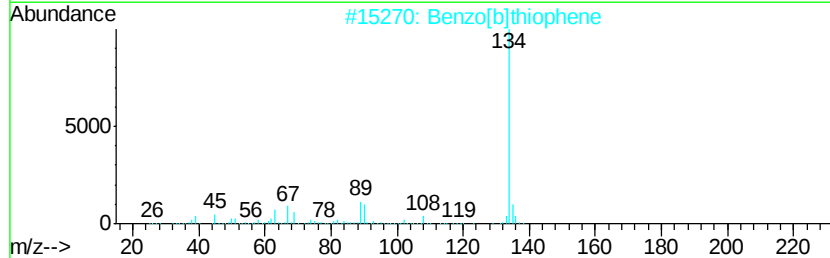
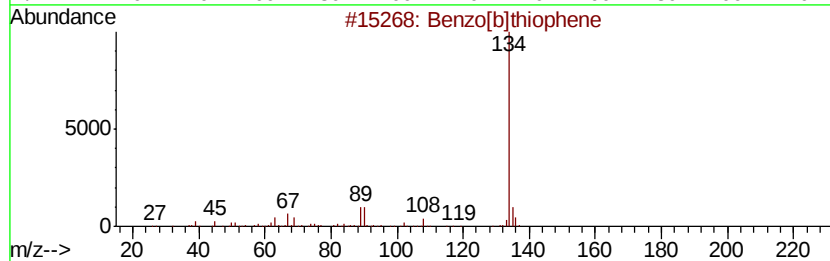
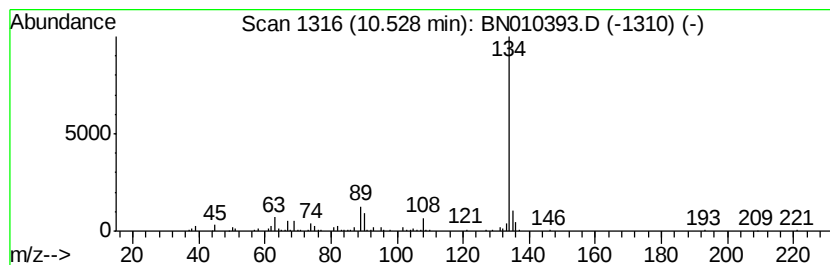
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzo[b]thiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.53	2.19 ng/ul	369805	Naphthalene-d8	10.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene	134	C8H6S	000095-15-8	91
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	91
3		Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	90
4		Benzo[c]thiophene	134	C8H6S	000270-82-6	50
5		Benzo[b]thiophene	134	C8H6S	000095-15-8	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010393.D
Acq On : 02 Apr 2020 15:32
Operator : CG/JU
Sample : L2065-10DL2 200X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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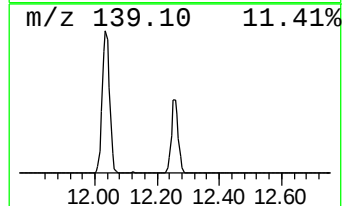
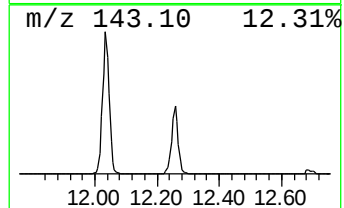
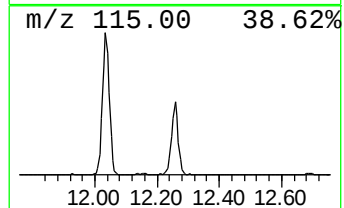
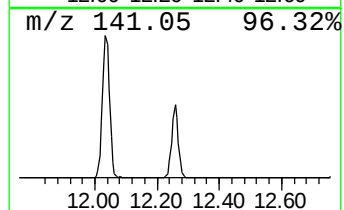
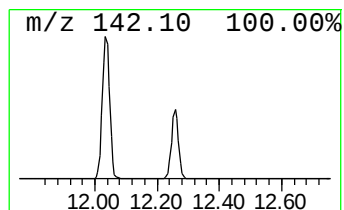
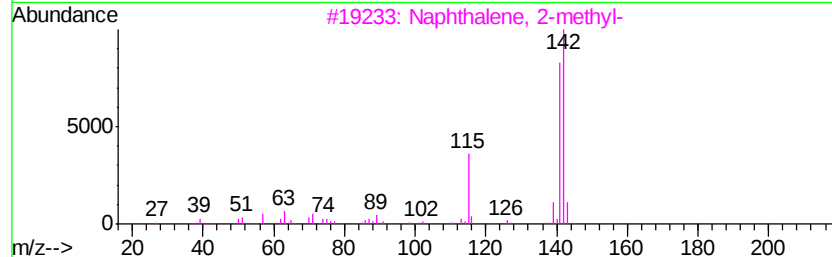
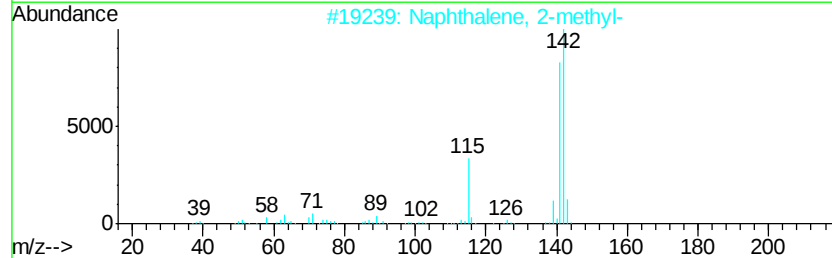
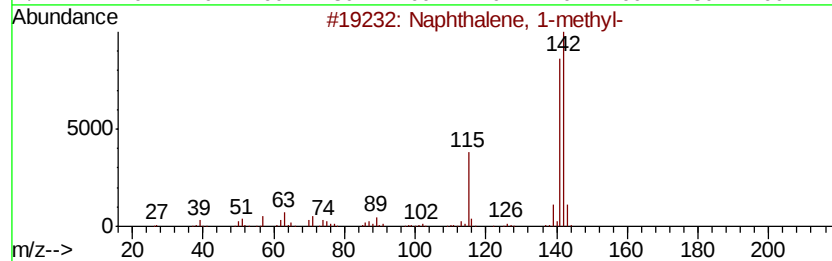
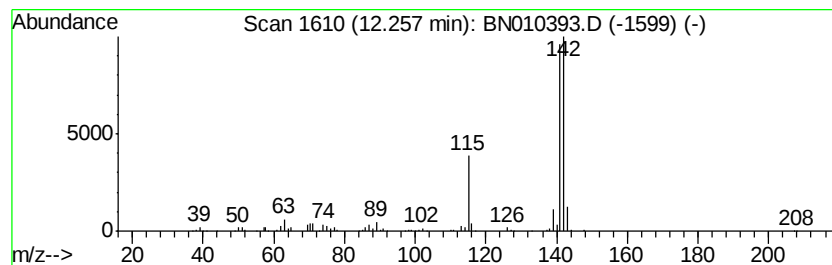
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	5.41 ng/ul	914621	Naphthalene-d8	10.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010393.D
Acq On : 02 Apr 2020 15:32
Operator : CG/JU
Sample : L2065-10DL2 200X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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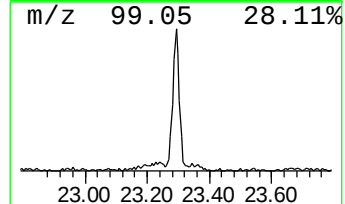
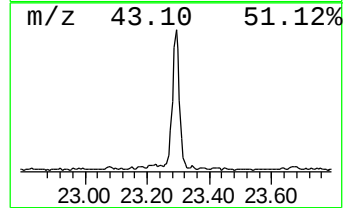
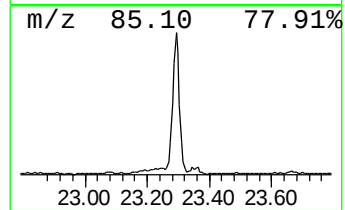
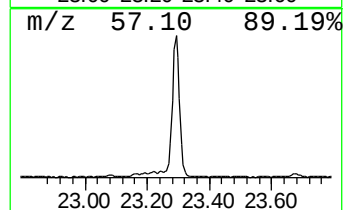
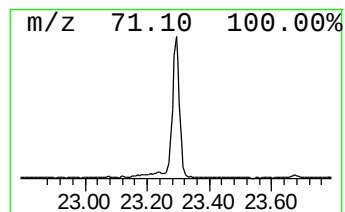
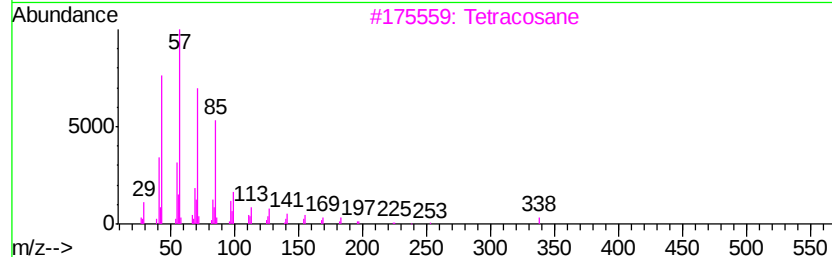
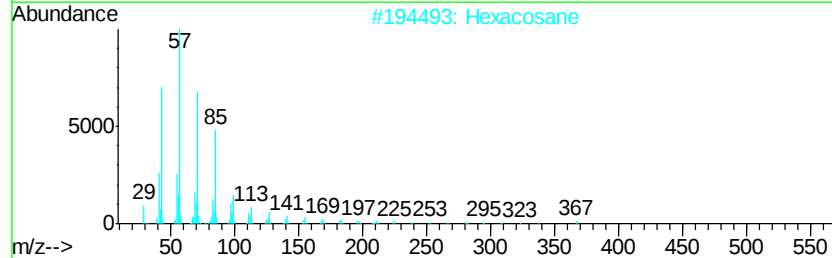
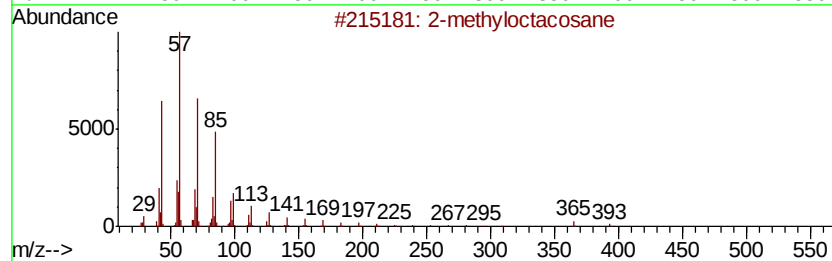
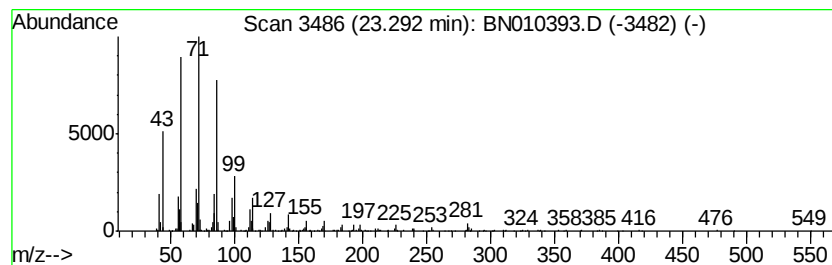
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.29	2.04 ng/ul	1001470	Perylene-d12	23.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Tetracosane	338	C24H50	000646-31-1	87
4		Octacosane	394	C28H58	000630-02-4	87
5		triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010393.D
Acq On : 02 Apr 2020 15:32
Operator : CG/JU
Sample : L2065-10DL2 200X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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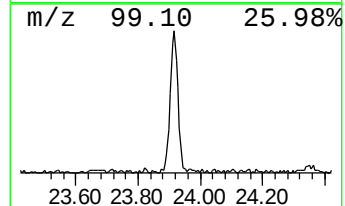
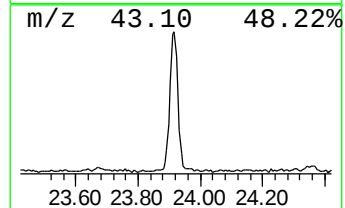
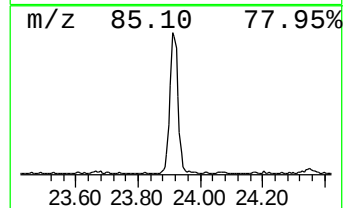
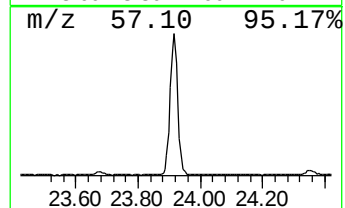
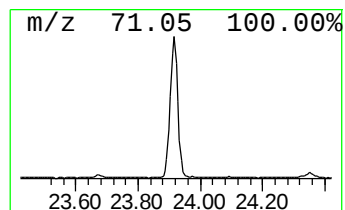
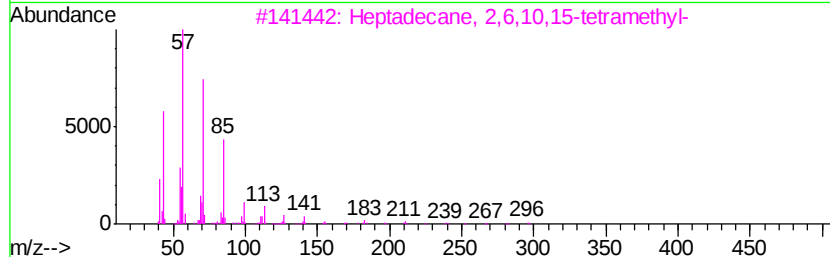
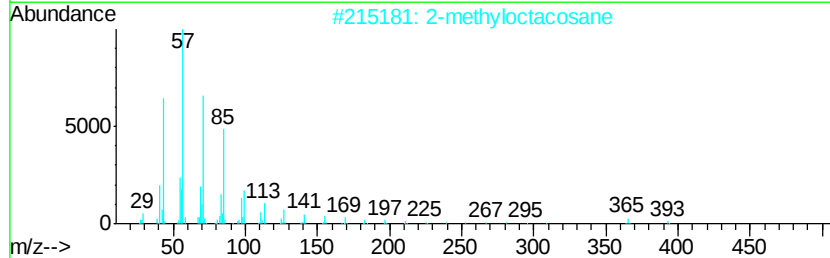
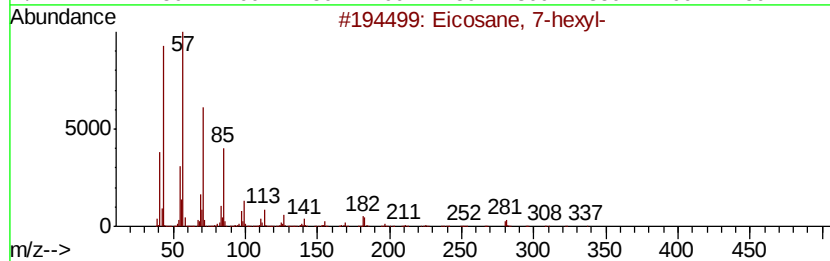
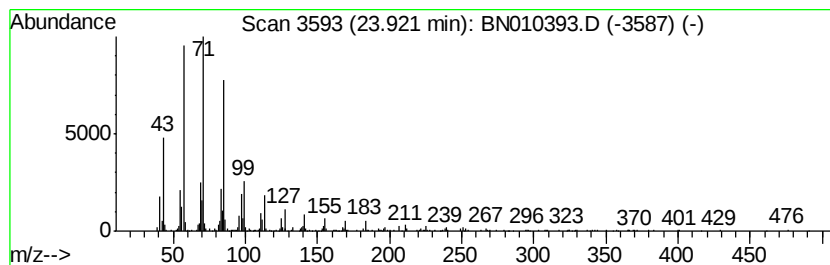
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.92	2.73 ng/ul	1335750	Perylene-d12	23.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
4		Hentriacontane	437	C31H64	000630-04-6	83
5		Tetratetracontane	619	C44H90	007098-22-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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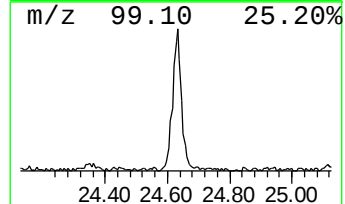
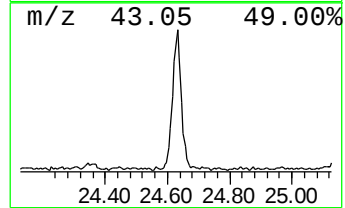
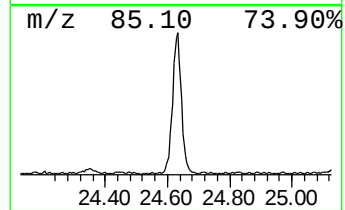
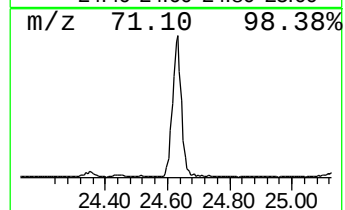
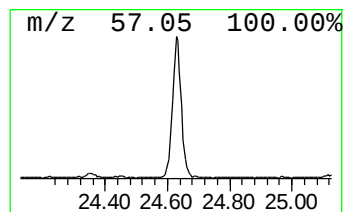
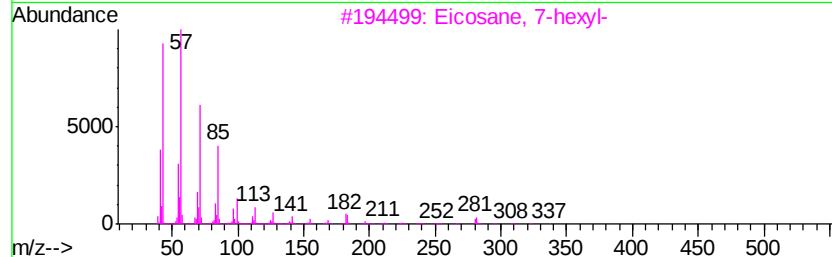
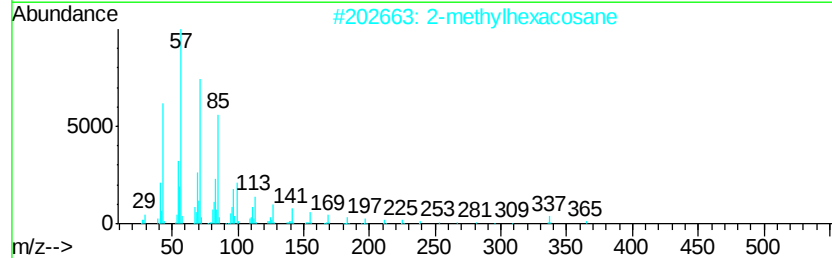
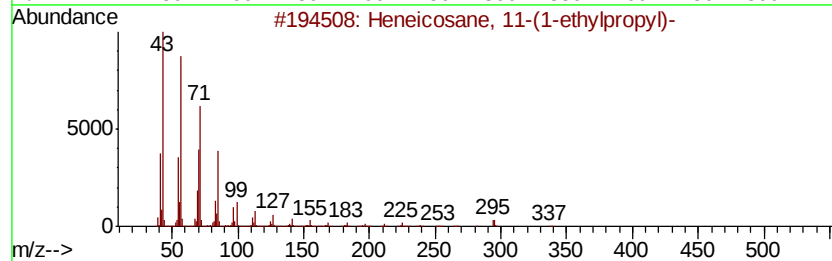
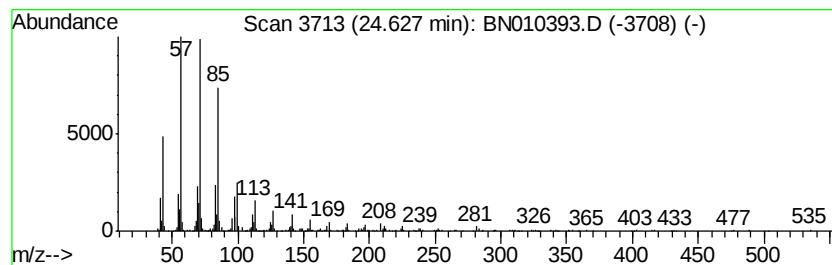
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.63	2.83 ng/ul	1388740	Perylene-d12	23.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
2		2-methylhexacosane	380	C27H56	1000376-72-7	91
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040120\
Data File : BN010393.D
Acq On : 02 Apr 2020 15:32
Operator : CG/JU
Sample : L2065-10DL2 200X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
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
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Benzene, 1-propynyl-	8.14	5.5	ng/ul	594267	1	7.61	2155610	20.0
Cyclohexanemethan...	9.75	8.0	ng/ul	1356190	2	10.37	3384020	20.0
Benzo[b]thiophene	10.53	2.2	ng/ul	369805	2	10.37	3384020	20.0
Naphthalene, 1-me...	12.26	5.4	ng/ul	914621	2	10.37	3384020	20.0
(DEL) Alkane: Str...	23.29	2.0	ng/ul	1001470	6	23.36	9798890	20.0
(DEL) Alkane: Str...	23.92	2.7	ng/ul	1335750	6	23.36	9798890	20.0
(DEL) Alkane: Str...	24.63	2.8	ng/ul	1388740	6	23.36	9798890	20.0