

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
 Data File : BN005384.D
 Acq On : 30 Apr 2019 22:32
 Operator : JU/SJ
 Sample : K2481-17ME
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GAHXOME

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-BN041919MA.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	5.293	386	392	408	rVB	1011615	2122611	3.07%	0.335%
2	5.634	443	450	451	rBV	709893	1042111	1.50%	0.164%
3	6.775	637	644	650	rVV2	1492066	2624891	3.79%	0.414%
4	6.940	665	672	678	rBV	1148165	1809726	2.61%	0.285%
5	7.128	695	704	711	rVV	1452928	2394692	3.46%	0.377%
6	7.252	719	725	734	rVV2	3980434	7234121	10.45%	1.140%
7	7.340	734	740	748	rVB	1095544	1630138	2.35%	0.257%
8	7.593	776	783	789	rBV	1240349	1967196	2.84%	0.310%
9	8.122	862	873	881	rBV	6728497	10779983	15.57%	1.699%
10	8.293	896	902	910	rVV	1624666	2604318	3.76%	0.410%
11	8.681	963	968	972	rVV	716172	1167925	1.69%	0.184%
12	8.734	972	977	987	rVV	1548554	2895767	4.18%	0.456%
13	8.851	992	997	1005	rVV	1116100	2194996	3.17%	0.346%
14	9.263	1061	1067	1074	rVB	671940	1144760	1.65%	0.180%
15	9.340	1074	1080	1085	rBV	1006350	1533943	2.22%	0.242%
16	9.451	1093	1099	1105	rVB	1245853	2179753	3.15%	0.344%
17	9.793	1147	1157	1163	rBV2	3585909	7709990	11.13%	1.215%
18	9.863	1163	1169	1173	rBV	2687054	5283432	7.63%	0.833%
19	9.981	1183	1189	1197	rVB	1893758	3282598	4.74%	0.517%
20	10.357	1246	1253	1256	rBV2	2037269	3825504	5.52%	0.603%
21	10.428	1256	1265	1271	rVB2	30681209	66272865	95.70%	10.445%
22	10.493	1271	1276	1280	rBV	1697684	2918257	4.21%	0.460%
23	11.398	1418	1430	1433	rBV	754149	1841783	2.66%	0.290%
24	11.440	1433	1437	1444	rVB4	532223	1297405	1.87%	0.204%
25	11.545	1450	1455	1464	rVB	759531	1451585	2.10%	0.229%
26	11.675	1468	1477	1484	rBV3	362314	1070715	1.55%	0.169%
27	11.804	1494	1499	1506	rBV2	857702	1485800	2.15%	0.234%
28	12.045	1529	1540	1546	rVB2	34077148	69249363	100.00%	10.914%
29	12.257	1564	1576	1584	rBV	19538027	33722192	48.70%	5.315%
30	13.051	1705	1711	1725	rBV2	3316843	5882930	8.50%	0.927%
31	13.251	1738	1745	1756	rVB	3242456	6806808	9.83%	1.073%
32	13.387	1761	1768	1770	rBV	8174207	13596873	19.63%	2.143%
33	13.551	1788	1796	1801	rVV	10168000	18748110	27.07%	2.955%
34	13.604	1801	1805	1809	rVV	6366015	9226377	13.32%	1.454%

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35	13.645	1809	1812	1815	rVV	3670828	5495407	7.94%	0.866%
36	13.681	1815	1818	1824	rVB2	7551218	10693800	15.44%	1.685%
37	13.787	1830	1836	1840	rBV	4851092	7809563	11.28%	1.231%
38	13.922	1852	1859	1860	rBV	4099716	5631005	8.13%	0.887%
39	13.951	1860	1864	1875	rVB	11490650	17516588	25.29%	2.761%
40	14.210	1901	1908	1916	rVV3	3485835	8451978	12.21%	1.332%
41	14.286	1916	1921	1924	rVV2	1454017	2289181	3.31%	0.361%
42	14.322	1924	1927	1931	rVB	1403480	1825528	2.64%	0.288%
43	14.439	1941	1947	1956	rVB2	2718153	5164997	7.46%	0.814%
44	14.634	1970	1980	1987	rVV3	2700634	5565332	8.04%	0.877%
45	14.710	1987	1993	1998	rVB	2179812	3032273	4.38%	0.478%
46	14.839	2008	2015	2020	rBV2	3821352	7096396	10.25%	1.118%
47	14.904	2023	2026	2030	rVB	1427407	1837004	2.65%	0.290%
48	14.998	2037	2042	2045	rBV	1160588	1863549	2.69%	0.294%
49	15.104	2055	2060	2066	rVB	2661955	3770911	5.45%	0.594%
50	15.169	2067	2071	2074	rBV	740386	1104294	1.59%	0.174%
51	15.228	2075	2081	2086	rVV2	7372343	10806697	15.61%	1.703%
52	15.286	2086	2091	2100	rVV	6676432	11489063	16.59%	1.811%
53	15.363	2100	2104	2108	rVV2	1508055	2117913	3.06%	0.334%
54	15.404	2108	2111	2115	rVB2	910867	1124619	1.62%	0.177%
55	15.486	2121	2125	2131	rBV2	687630	1291109	1.86%	0.203%
56	15.575	2135	2140	2144	rBV2	1120882	1956309	2.83%	0.308%
57	15.692	2155	2160	2164	rBV	2780852	4055825	5.86%	0.639%
58	15.963	2200	2206	2211	rBV2	781849	1479265	2.14%	0.233%
59	16.286	2254	2261	2266	rVV2	2087958	5069045	7.32%	0.799%
60	16.339	2266	2270	2275	rVV	2197893	2968889	4.29%	0.468%
61	16.439	2282	2287	2292	rVV3	1082281	1675991	2.42%	0.264%
62	16.645	2318	2322	2329	rVB	1136830	1689107	2.44%	0.266%
63	16.798	2343	2348	2352	rVB2	1577299	2335925	3.37%	0.368%
64	16.980	2373	2379	2382	rBV	3629607	5369646	7.75%	0.846%
65	17.033	2382	2388	2392	rVV	21341392	33345950	48.15%	5.256%
66	17.080	2392	2396	2399	rVV2	5883357	8396788	12.13%	1.323%
67	17.116	2399	2402	2407	rVB	4982748	6523767	9.42%	1.028%
68	17.175	2408	2412	2417	rVB2	744266	1205463	1.74%	0.190%
69	17.563	2474	2478	2484	rVB2	804871	1283727	1.85%	0.202%
70	17.857	2523	2528	2532	rVV	6755165	9397718	13.57%	1.481%
71	17.910	2532	2537	2545	rVV	7537511	10888284	15.72%	1.716%

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72	17.986	2545	2550	2555	rVV	2293305	3313272	4.78%	0.522%
73	18.051	2555	2561	2565	rVV2	5117167	9575758	13.83%	1.509%
74	18.092	2565	2568	2573	rVB	3432600	4439797	6.41%	0.700%
75	18.363	2610	2614	2621	rVB	2579286	3440949	4.97%	0.542%
76	18.616	2653	2657	2661	rBV2	766031	1057040	1.53%	0.167%
77	18.786	2681	2686	2691	rBV2	2245059	3787154	5.47%	0.597%
78	18.851	2691	2697	2700	rBV2	1162200	2298404	3.32%	0.362%
79	18.880	2700	2702	2706	rVB	1400337	1473146	2.13%	0.232%
80	19.051	2726	2731	2740	rBV2	5481195	7932652	11.46%	1.250%
81	19.204	2753	2757	2762	rVB	2941820	3744422	5.41%	0.590%
82	19.392	2783	2789	2791	rBV	7695936	10957763	15.82%	1.727%
83	19.422	2791	2794	2803	rVV	7964478	10976951	15.85%	1.730%
84	19.751	2845	2850	2860	rVB	1074080	2057839	2.97%	0.324%
85	19.892	2868	2874	2876	rBV3	1035949	1872442	2.70%	0.295%
86	19.921	2876	2879	2884	rVB	2402704	3154037	4.55%	0.497%
87	20.021	2892	2896	2902	rBV	1969395	2641283	3.81%	0.416%
88	20.080	2902	2906	2910	rVB2	1318161	1599876	2.31%	0.252%
89	20.221	2926	2930	2934	rVV	1075298	1450944	2.10%	0.229%
90	20.269	2934	2938	2947	rVB	1631894	2754044	3.98%	0.434%
91	20.639	2997	3001	3008	rBV2	416332	1105172	1.60%	0.174%
92	20.933	3047	3051	3066	rVB5	1579033	2481279	3.58%	0.391%
93	21.198	3089	3096	3100	rBV3	6638795	12622126	18.23%	1.989%
94	21.239	3100	3103	3110	rVB	2934451	3790633	5.47%	0.597%
95	21.751	3187	3190	3195	rVV	1294265	1871709	2.70%	0.295%
96	21.974	3225	3228	3235	rVB2	883485	1178433	1.70%	0.186%
97	22.392	3295	3299	3304	rVB	1610113	2149498	3.10%	0.339%
98	22.762	3354	3362	3377	rVB3	1117084	3485090	5.03%	0.549%
99	23.262	3441	3447	3451	rBV2	5018998	8864646	12.80%	1.397%
100	23.398	3463	3470	3475	rBV	3084105	5795640	8.37%	0.913%

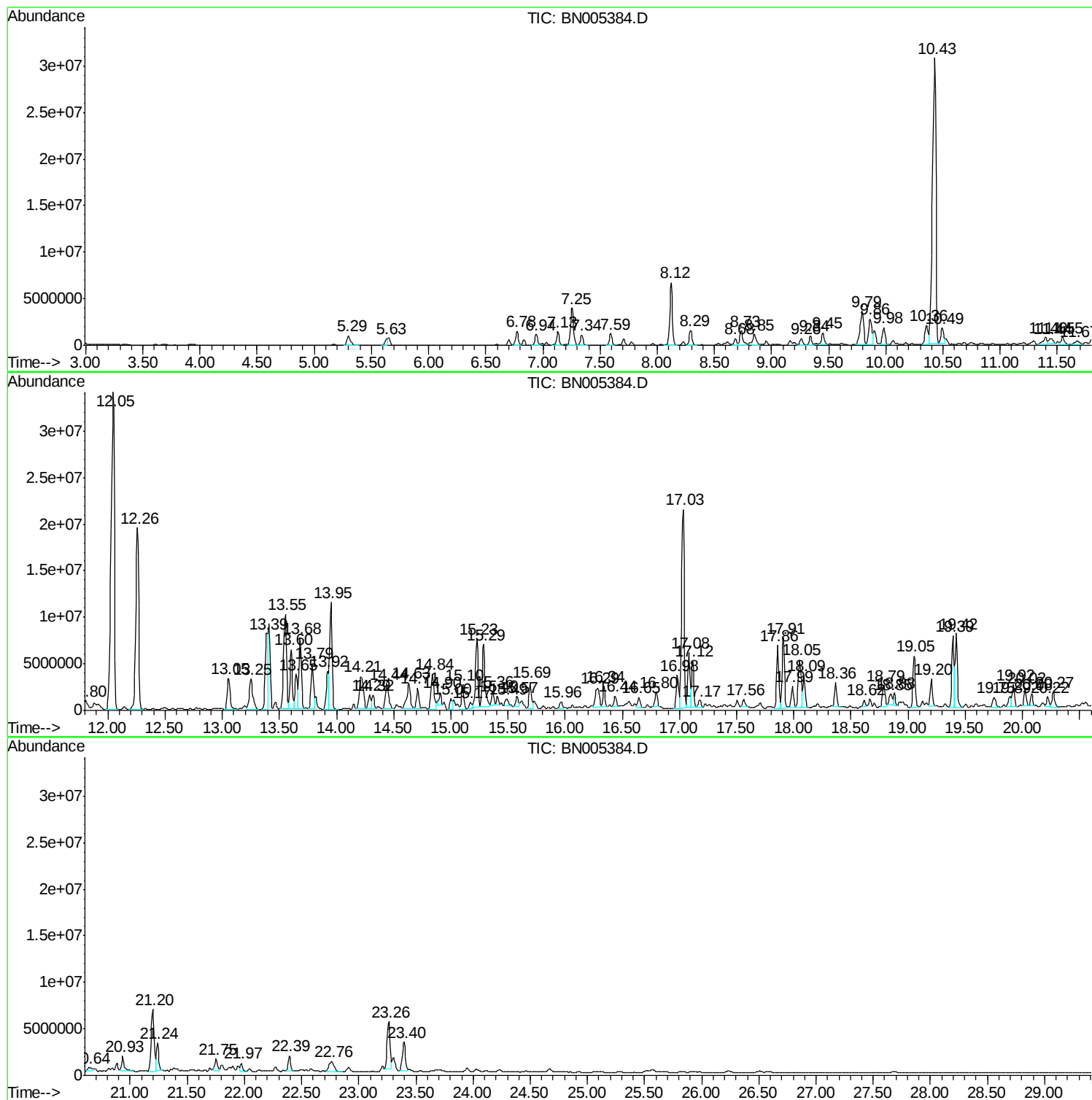
Sum of corrected areas: 634490423

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Instrument :
BNA_N
ClientSampled :
GAHXOME

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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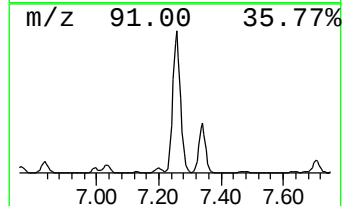
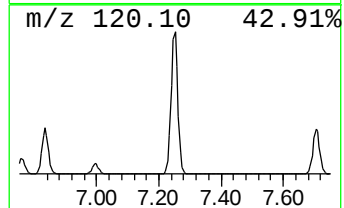
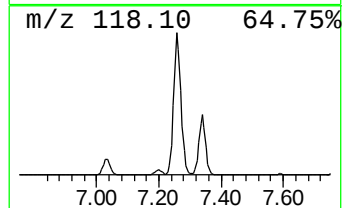
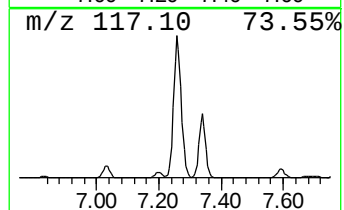
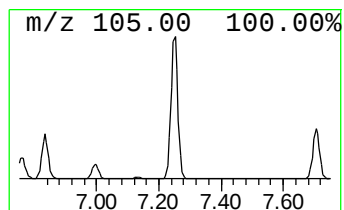
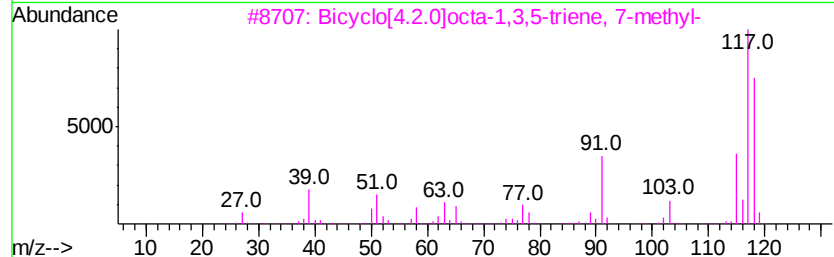
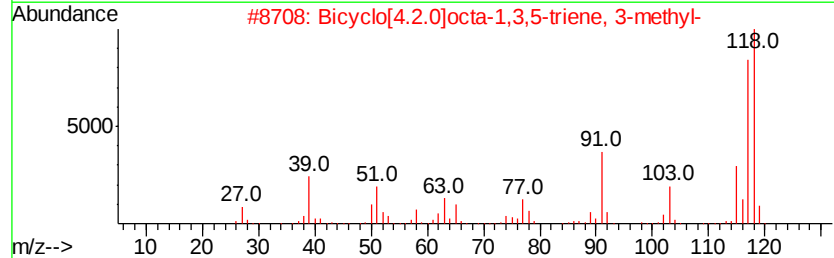
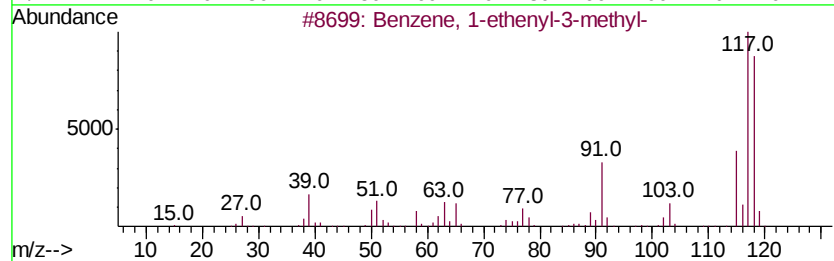
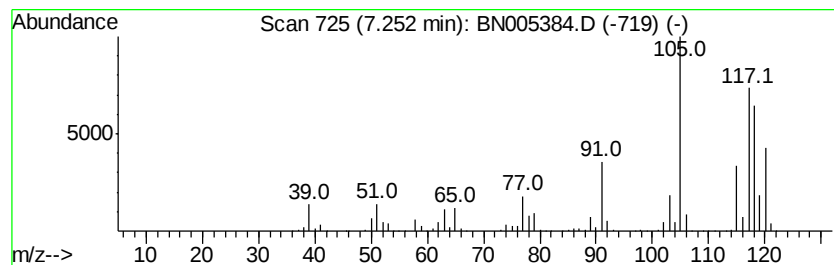
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethenyl-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.25	73.55 ng/ul	7234120	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	60
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	55
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	55
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	55



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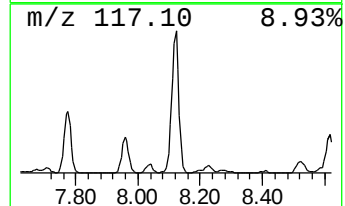
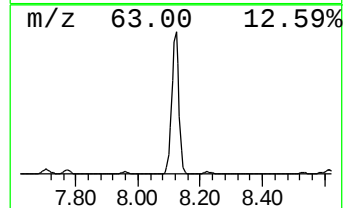
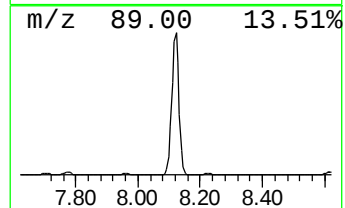
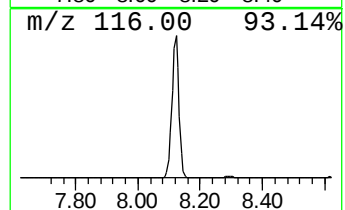
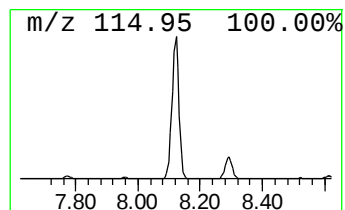
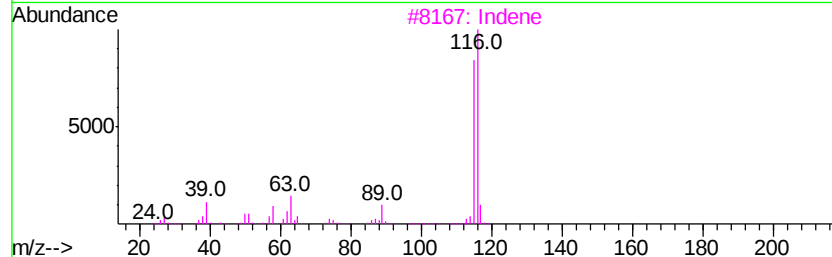
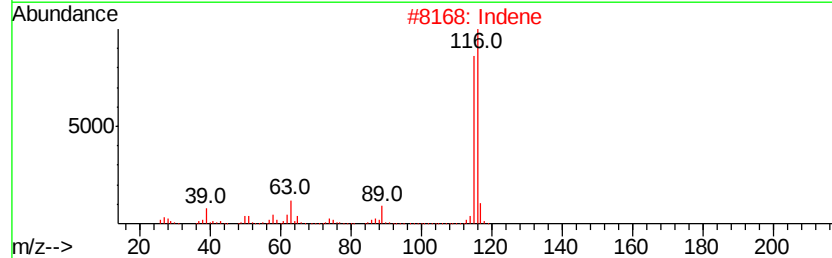
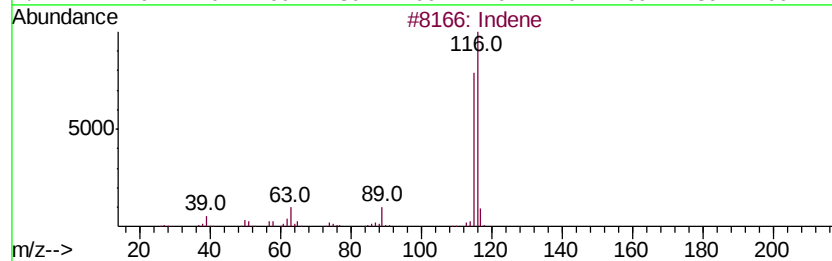
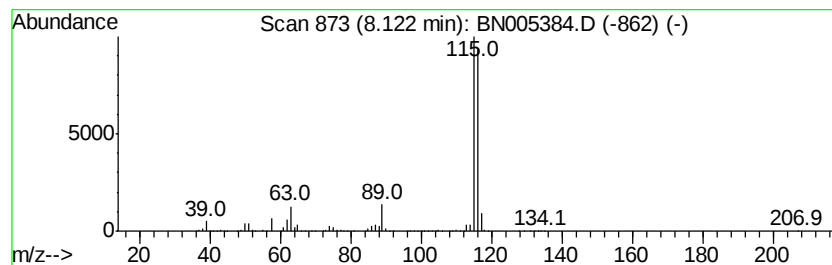
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	109.60 ng/ul	10780000	1,4-Dichlorobenzene-d4	7.59

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene		116	C9H8	000095-13-6	97
2	Indene		116	C9H8	000095-13-6	95
3	Indene		116	C9H8	000095-13-6	94
4	Benzene, 1-propynyl-		116	C9H8	000673-32-5	91
5	Benzene, 1-ethynyl-4-methyl-		116	C9H8	000766-97-2	91



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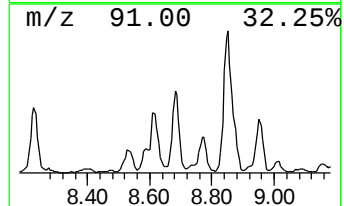
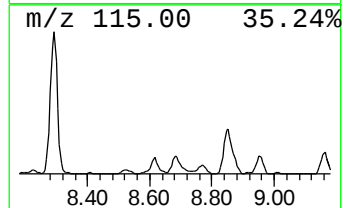
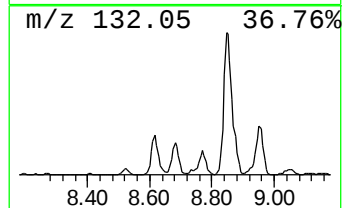
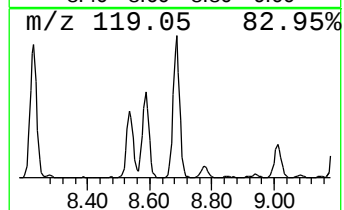
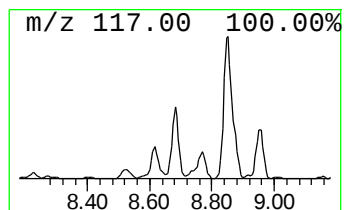
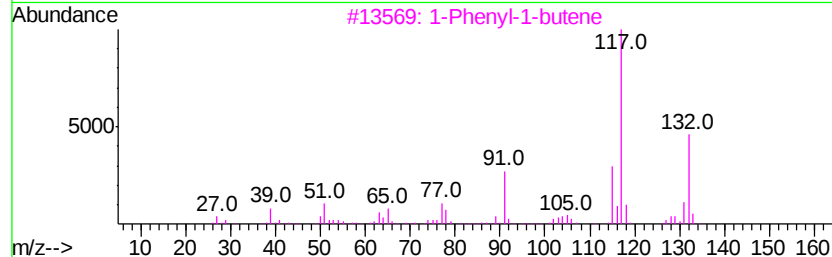
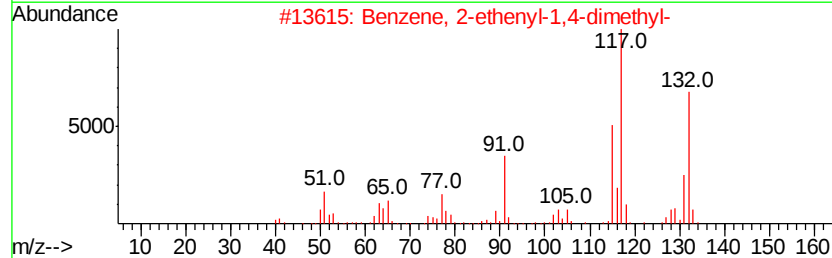
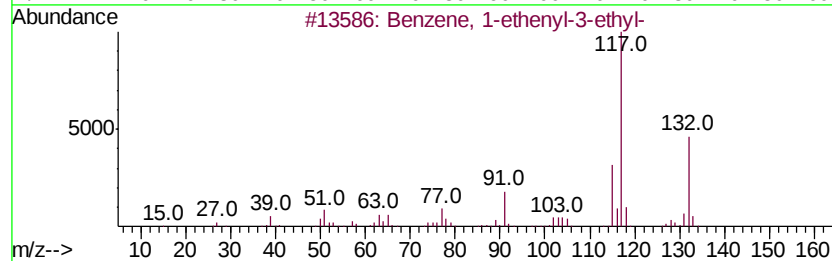
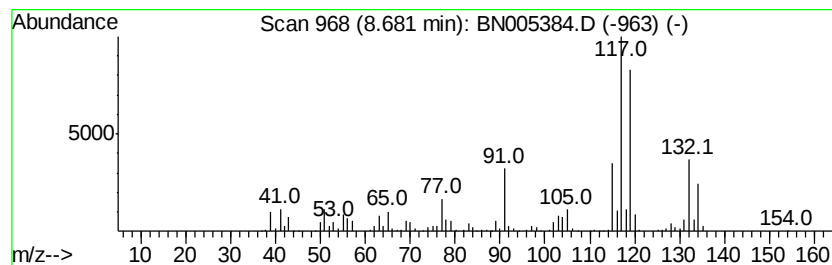
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.68	11.87 ng/ul	1167930	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	89
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	55
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	55
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	55
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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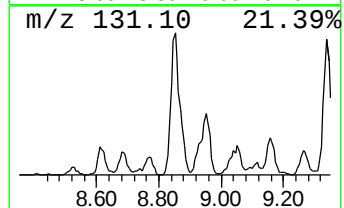
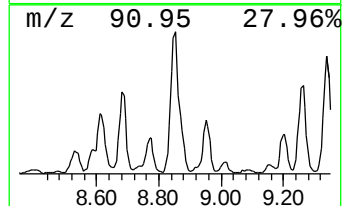
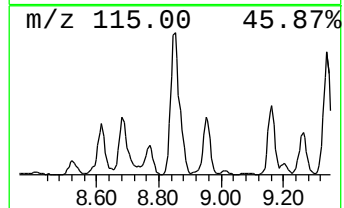
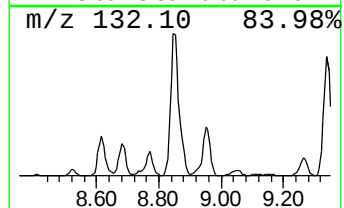
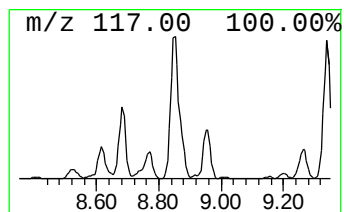
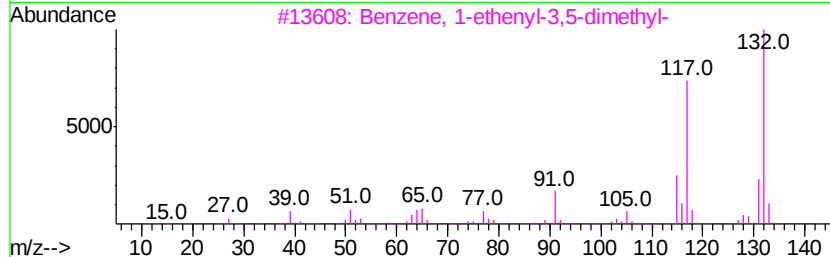
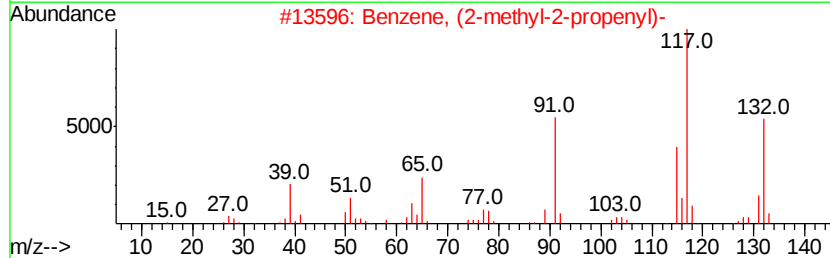
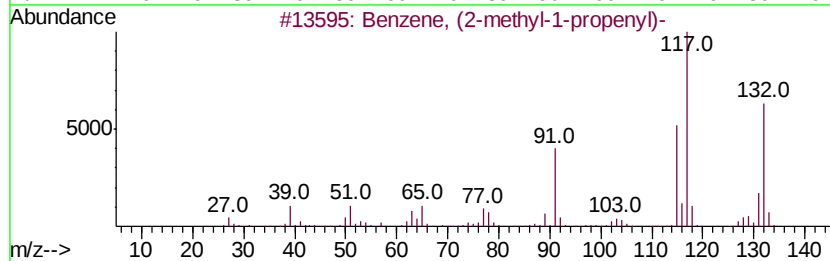
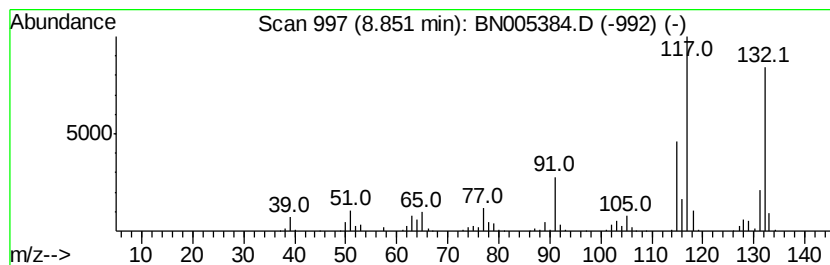
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, (2-methyl-1-propen... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.85	22.32 ng/ul	2195000	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	96
2		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	95
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	95
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	95
5		Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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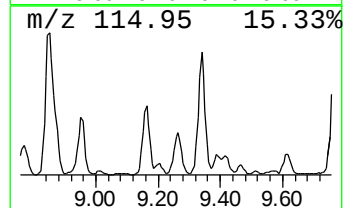
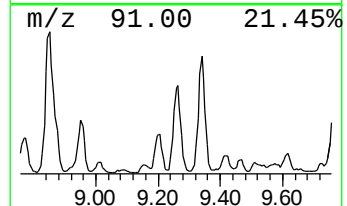
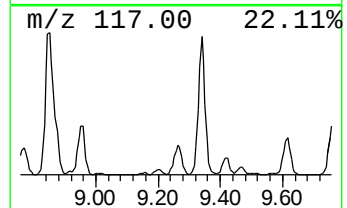
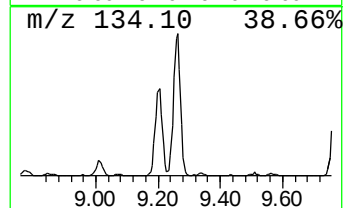
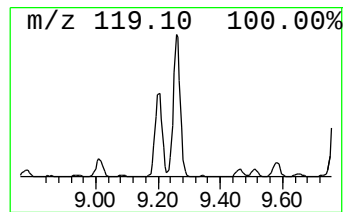
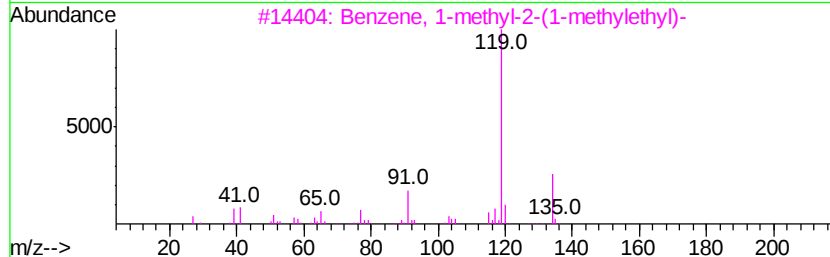
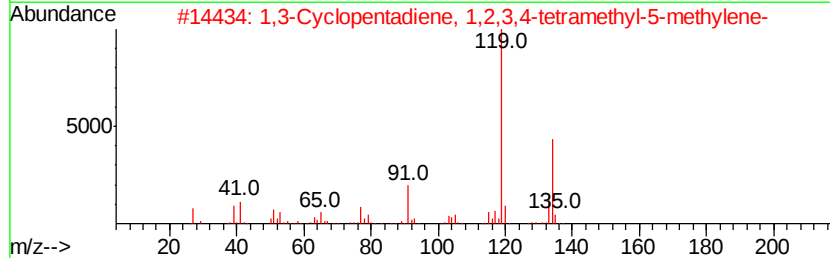
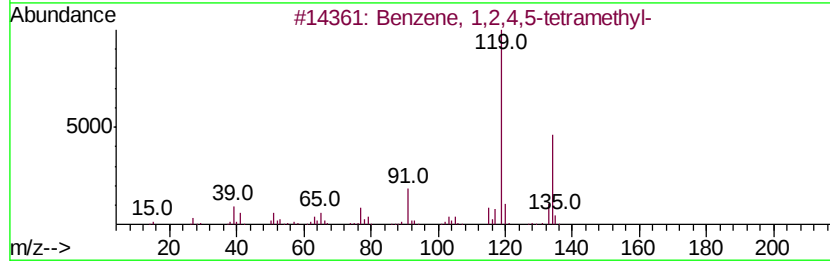
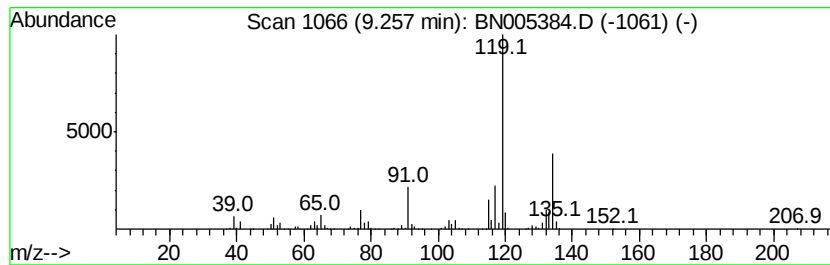
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.26	5.98 ng/ul	1144760	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81
2		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81
3		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	76
4		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	76
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
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Misc :
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ClientSampleId :
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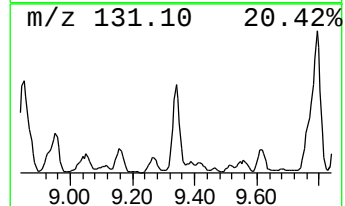
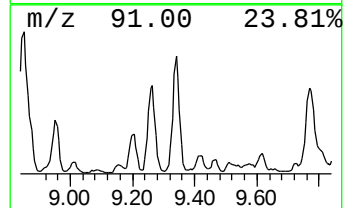
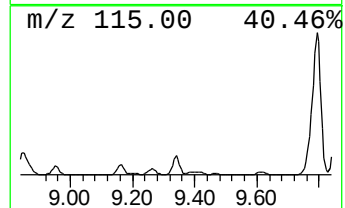
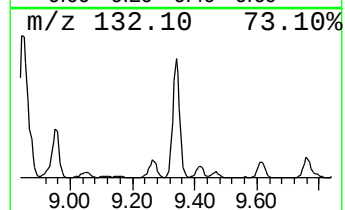
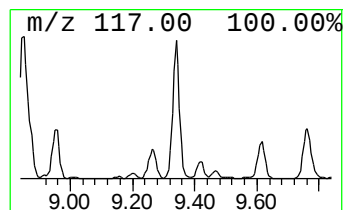
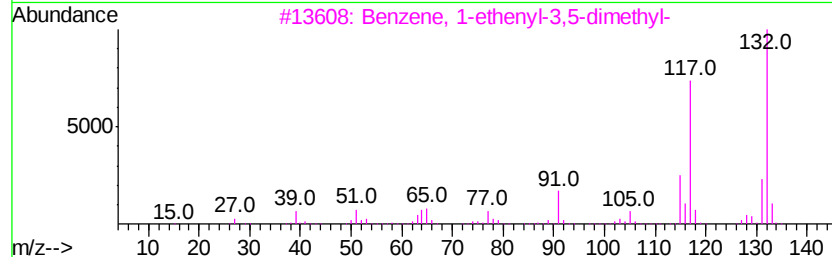
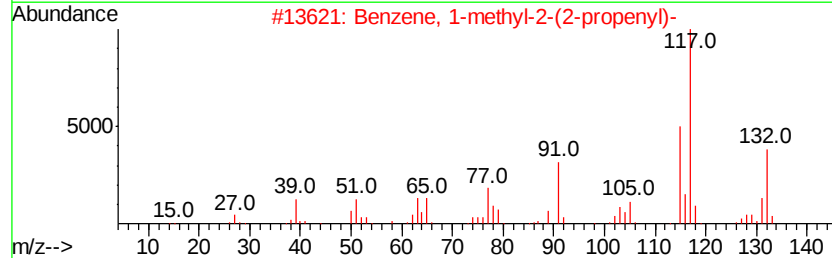
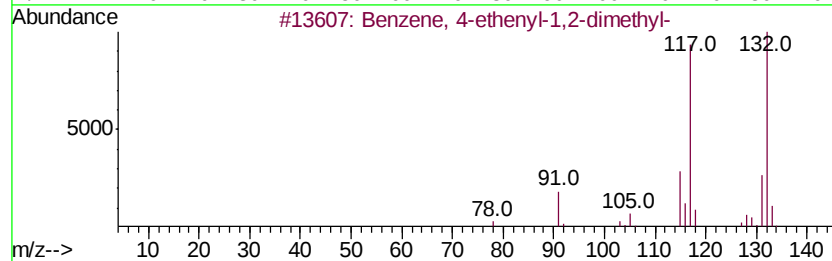
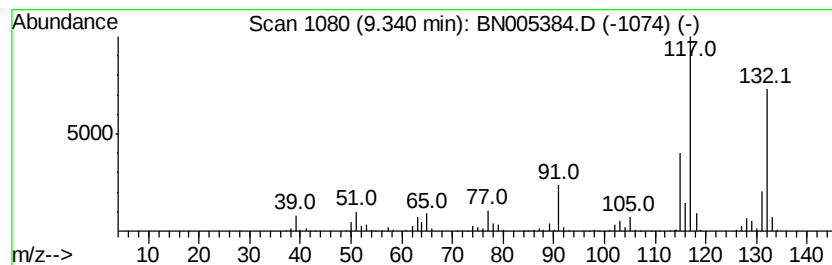
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 4-ethenyl-1,2-dime... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	8.02 ng/ul	1533940	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	97
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	94
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	94
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	94
5		Benzene, 1-butenyl-, (E)-	132	C10H12	001005-64-7	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
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Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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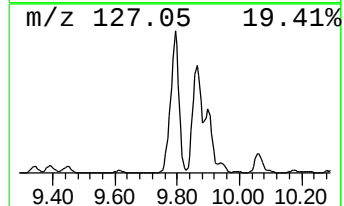
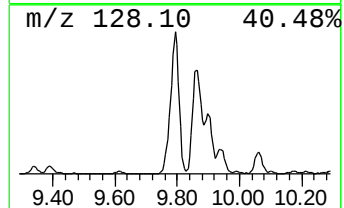
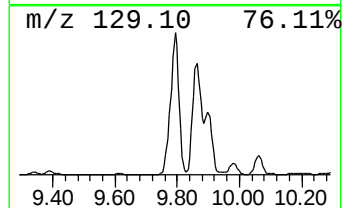
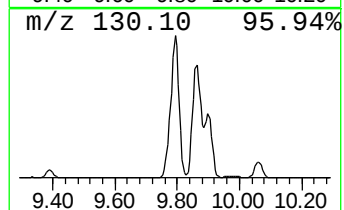
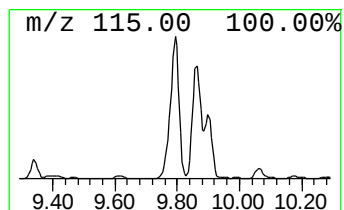
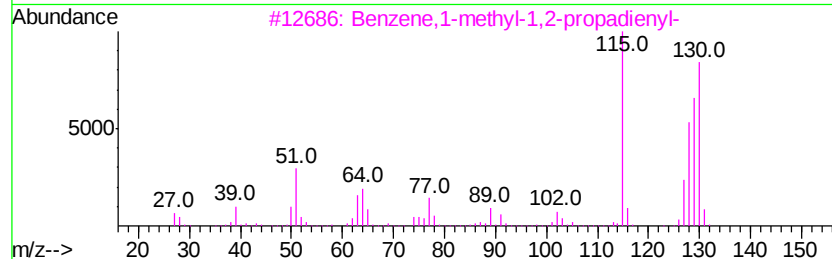
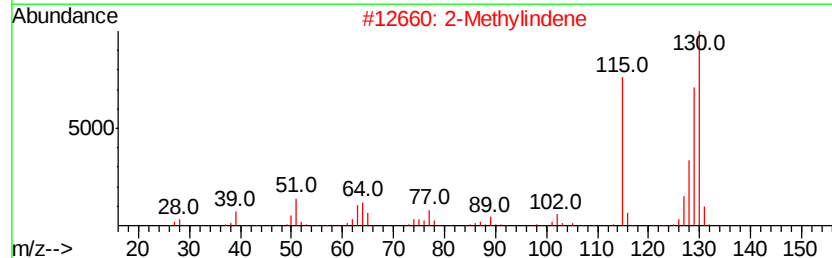
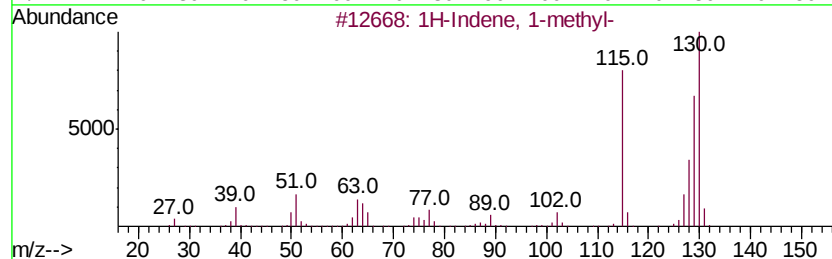
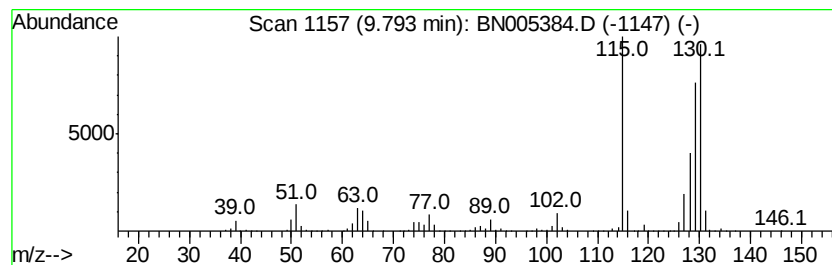
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1H-Indene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	40.31 ng/ul	7709990	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
2		2-Methylindene	130	C10H10	002177-47-1	95
3		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
4		Benzene, 1-butenyl-	130	C10H10	000622-76-4	94
5		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
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Acq On : 30 Apr 2019 22:32
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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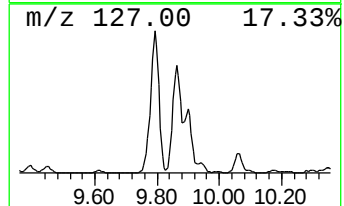
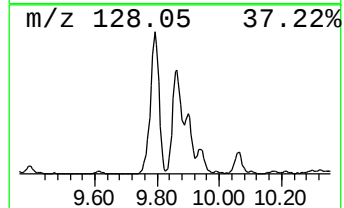
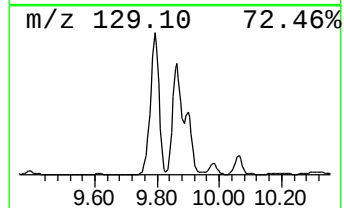
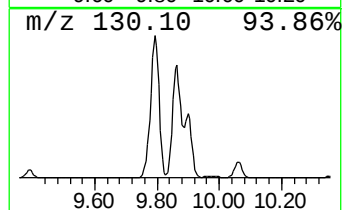
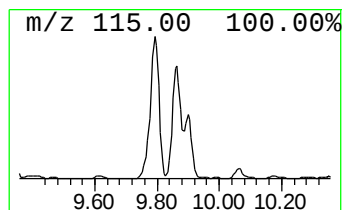
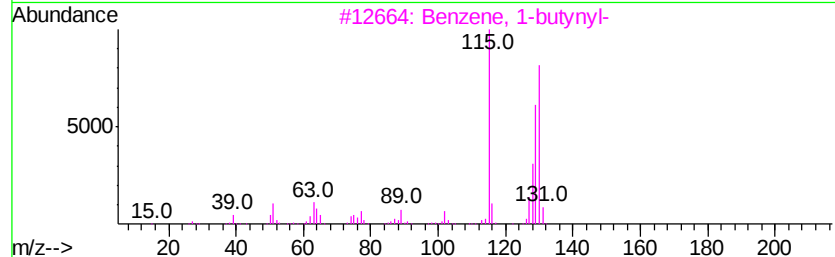
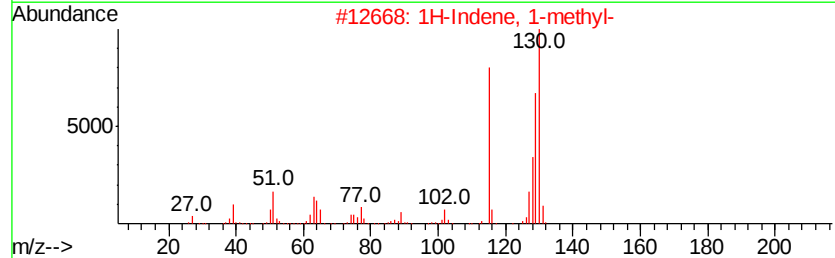
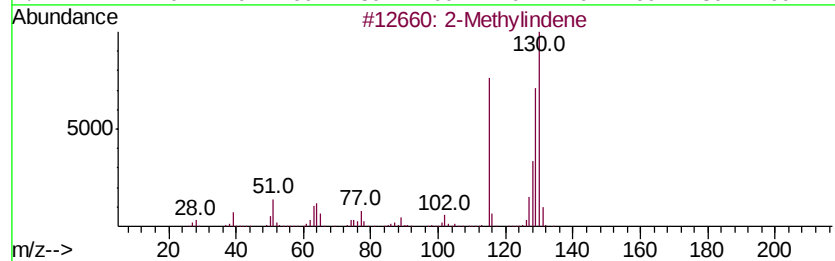
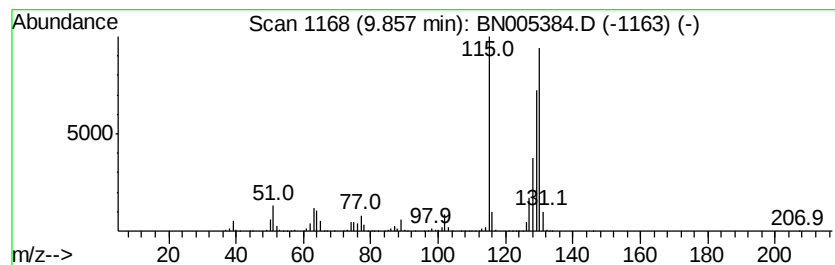
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 2-Methylindene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	27.62 ng/ul	5283430	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylindene	130	C10H10	002177-47-1	96
2		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
3		Benzene, 1-butynyl-	130	C10H10	000622-76-4	93
4		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	93
5		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
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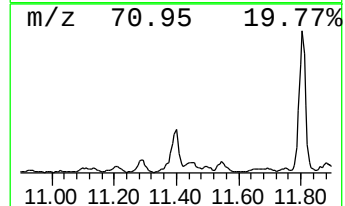
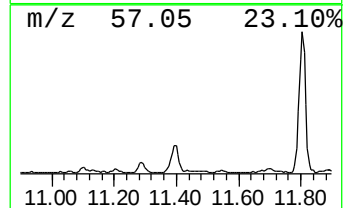
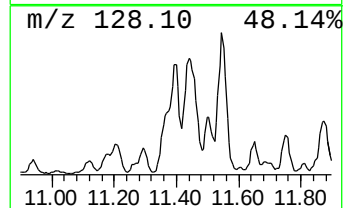
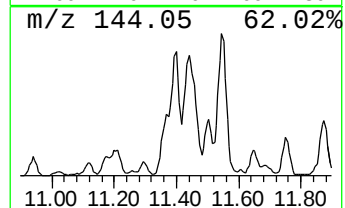
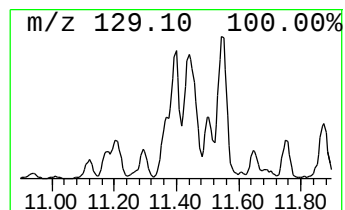
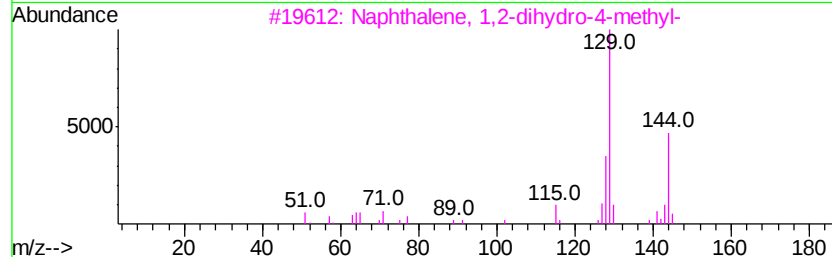
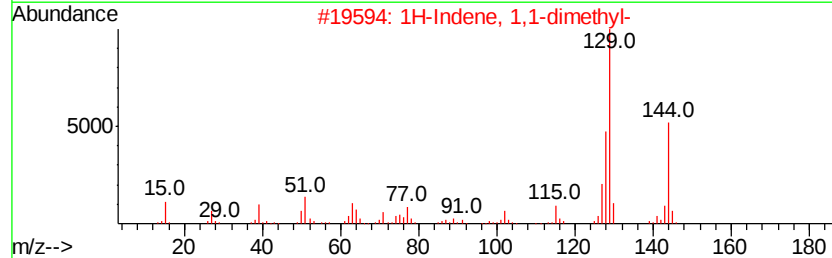
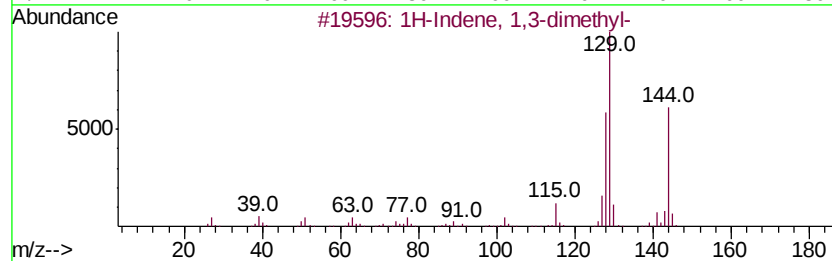
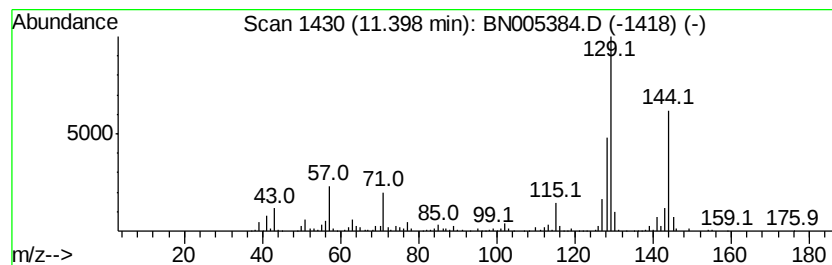
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 1H-Indene, 1,3-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	9.63 ng/ul	1841780	Naphthalene-d8	10.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	96
2			1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	91
3			Naphthalene, 1,2-dihydro-4-methyl-	144	C11H12	004373-13-1	90
4			1H-Indene, 2,3-dimethyl-	144	C11H12	004773-82-4	83
5			1H-Indene, 4,7-dimethyl-	144	C11H12	006974-97-6	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
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Sample : K2481-17ME
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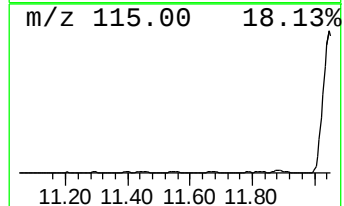
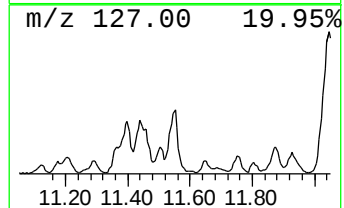
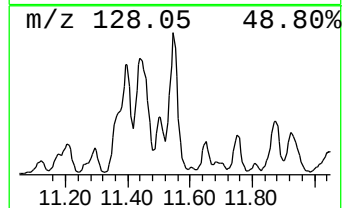
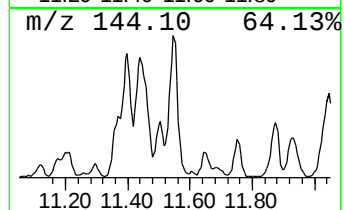
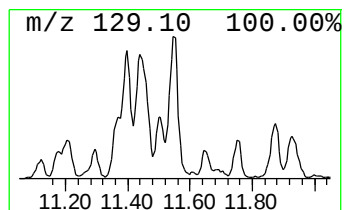
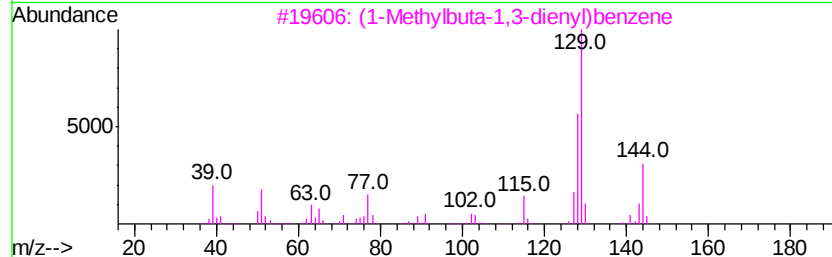
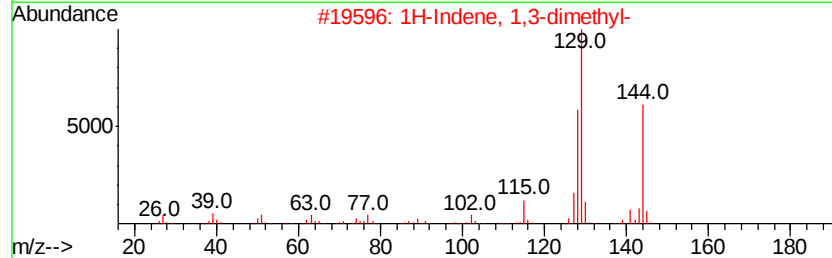
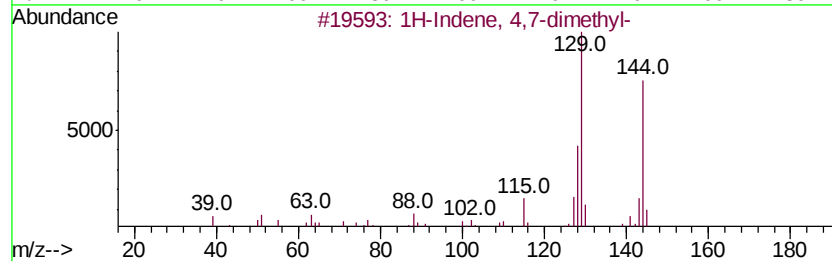
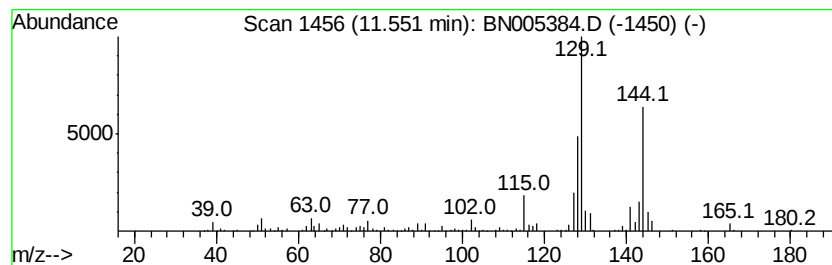
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 1H-Indene, 4,7-dimethyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	7.59 ng/ul	1451590	Naphthalene-d8	10.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 4,7-dimethyl-	144	C11H12	006974-97-6	94
2			1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	93
3			(1-Methylbuta-1,3-dienyl)benzene	144	C11H12	054758-36-0	91
4			1H-Cyclopropa[blnaphthalene, 1a,...	144	C11H12	006571-72-8	90
5			1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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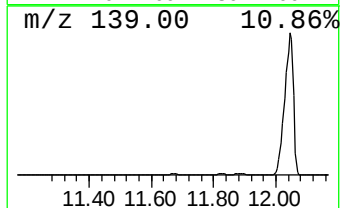
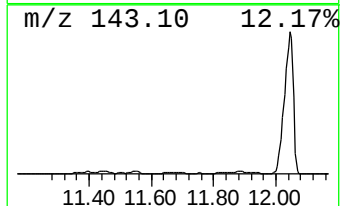
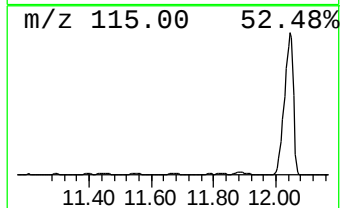
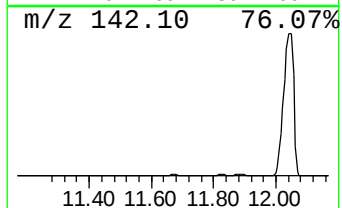
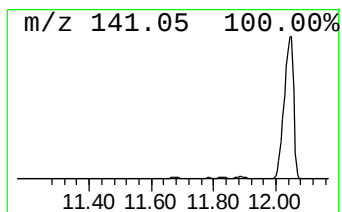
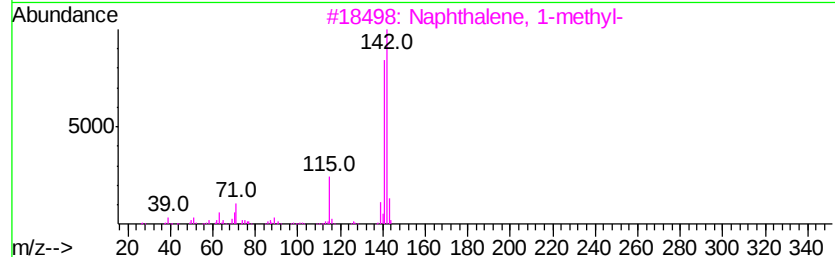
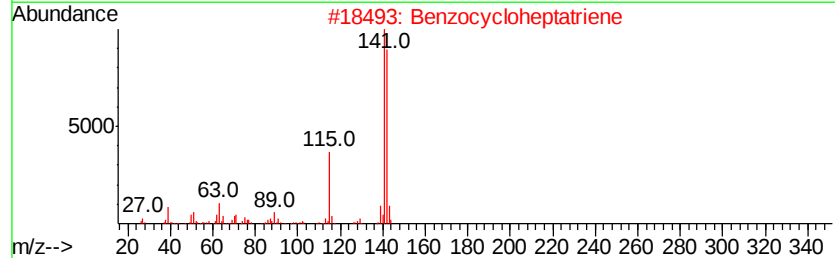
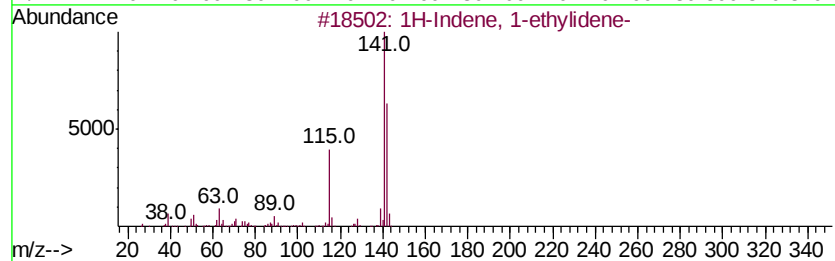
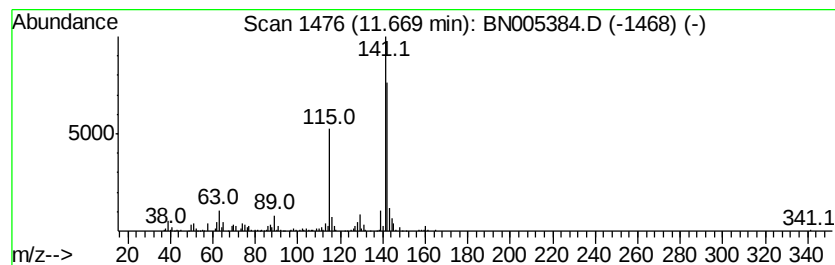
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 1H-Indene, 1-ethylidene- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	5.60 ng/ul	1070720	Naphthalene-d8	10.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	93
2			Benzocycloheptatriene	142	C11H10	000264-09-5	90
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90
4			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	90
5			Benzocycloheptatriene	142	C11H10	000264-09-5	87



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Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
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Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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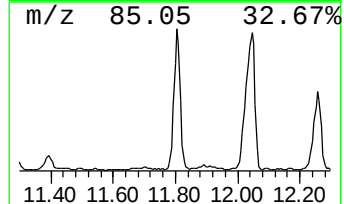
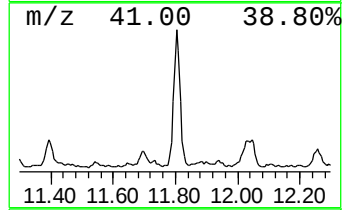
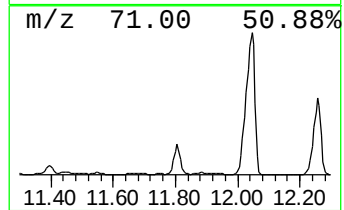
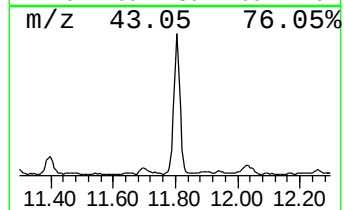
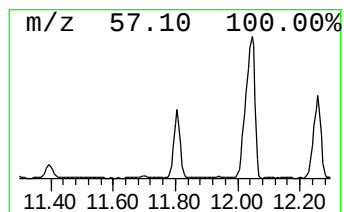
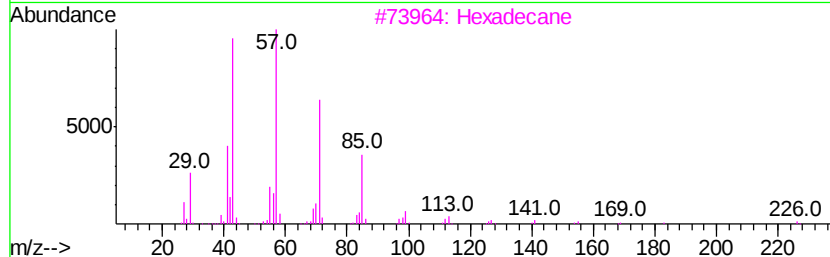
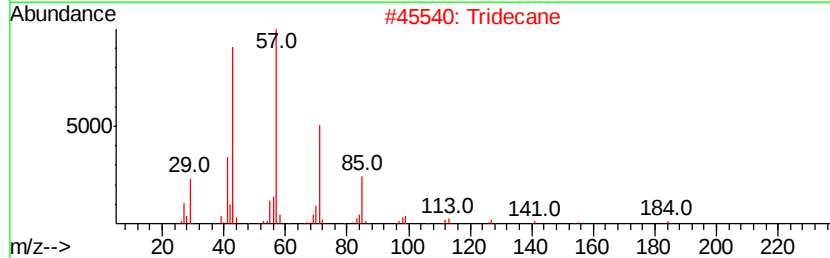
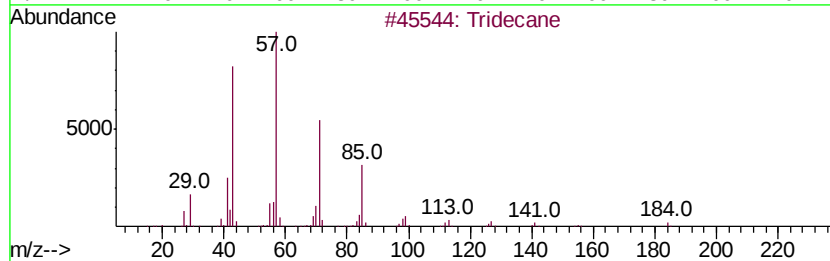
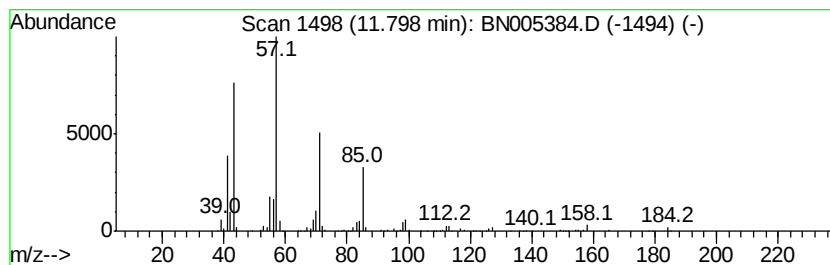
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	7.77 ng/ul	1485800	Naphthalene-d8	10.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	94
2			Tridecane	184	C13H28	000629-50-5	87
3			Hexadecane	226	C16H34	000544-76-3	86
4			Tetradecane	198	C14H30	000629-59-4	86
5			Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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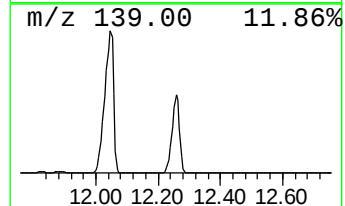
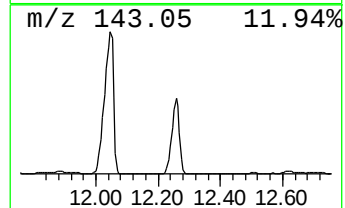
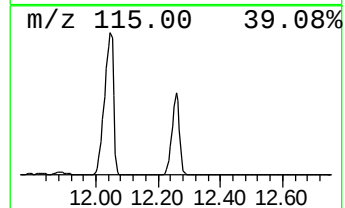
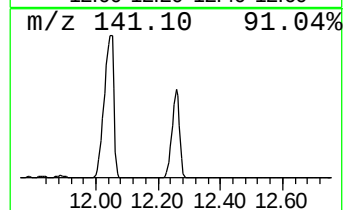
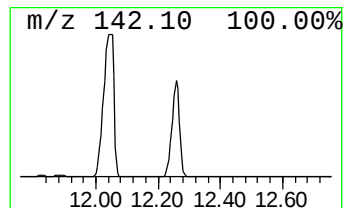
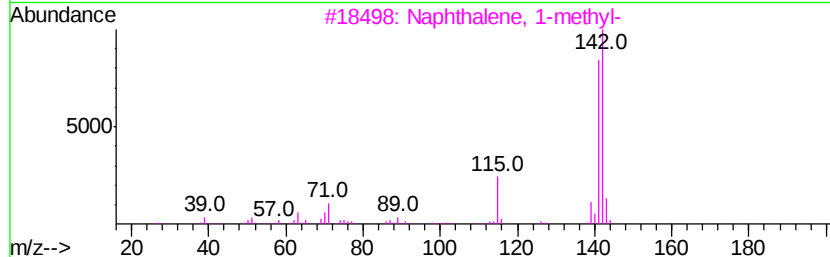
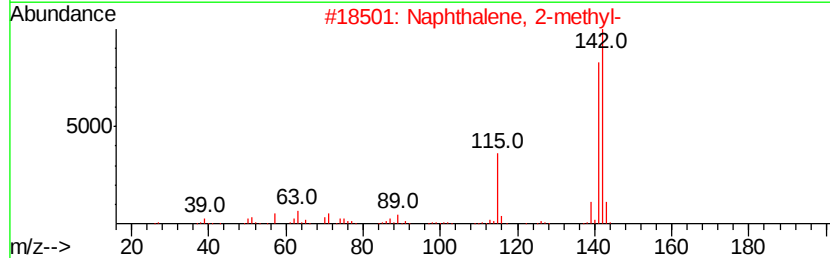
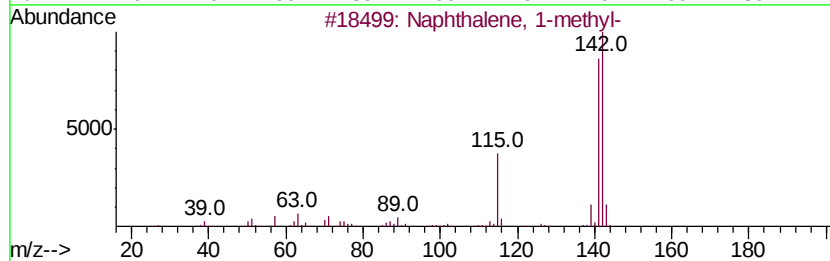
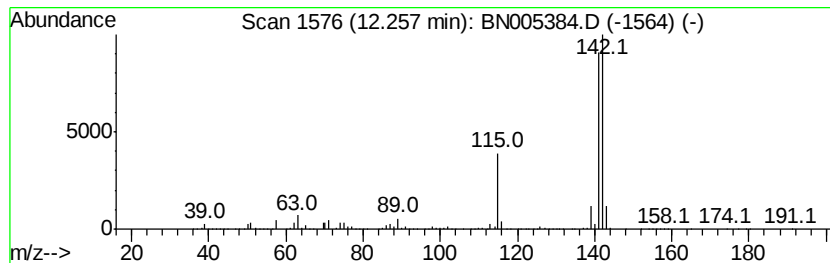
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	176.30 ng/ul	33722200	Naphthalene-d8	10.36

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, 1-methyl-	142	C11H10	000090-12-0	94
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
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Sample : K2481-17ME
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ALS Vial : 18 Sample Multiplier: 1

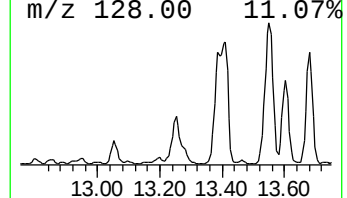
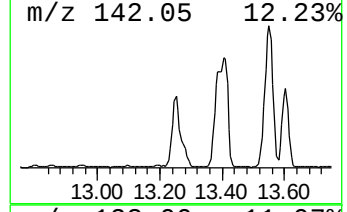
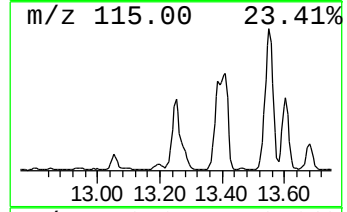
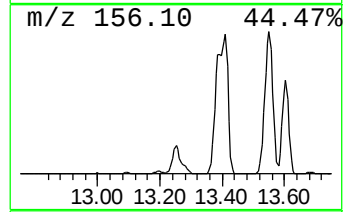
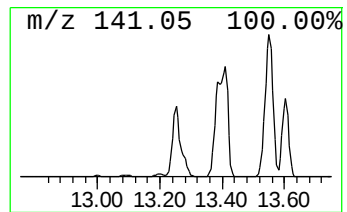
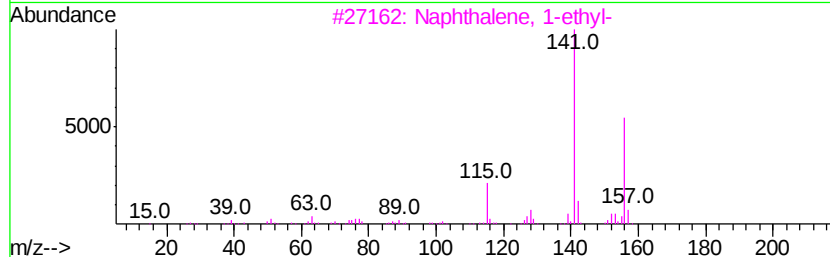
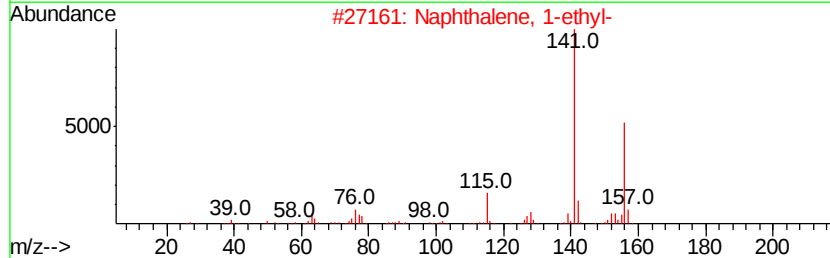
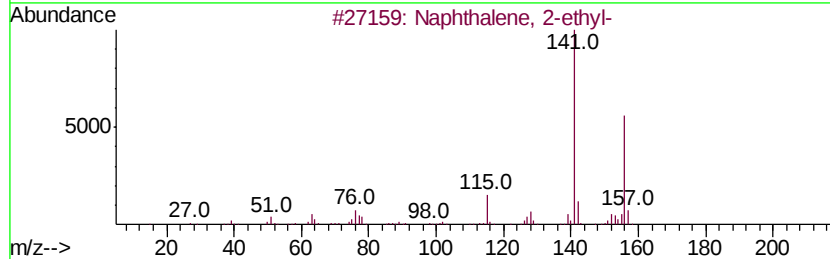
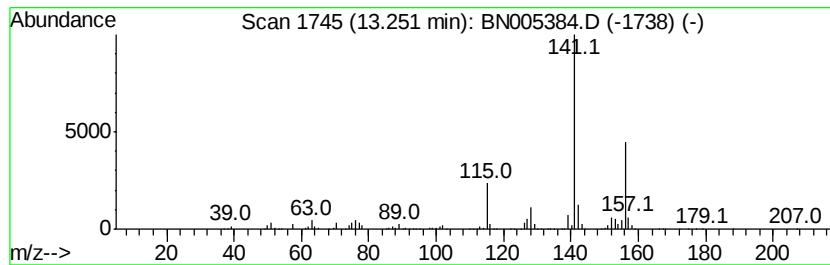
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Naphthalene, 2-ethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.25	16.11 ng/ul	6806810	Acenaphthene-d10	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 96
2		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 96
3		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 96
4		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 95
5		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
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Acq On : 30 Apr 2019 22:32
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Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

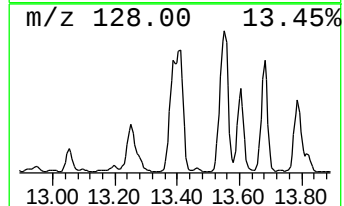
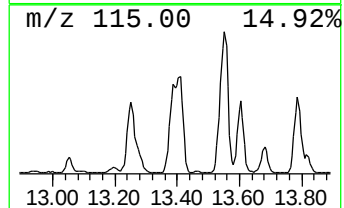
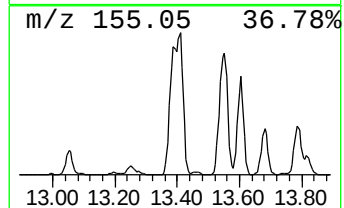
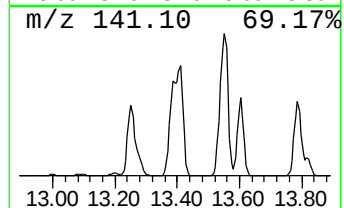
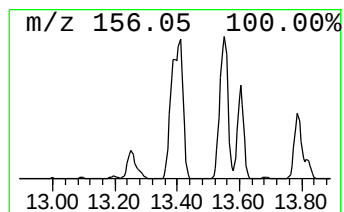
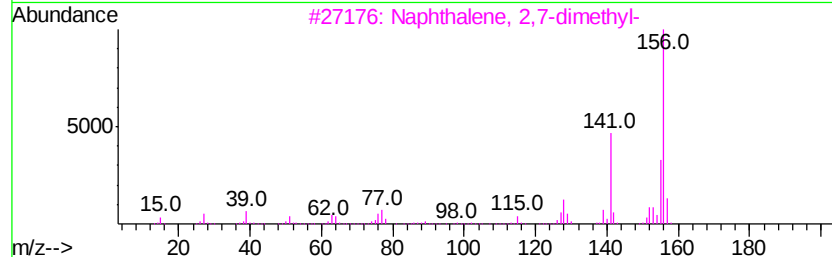
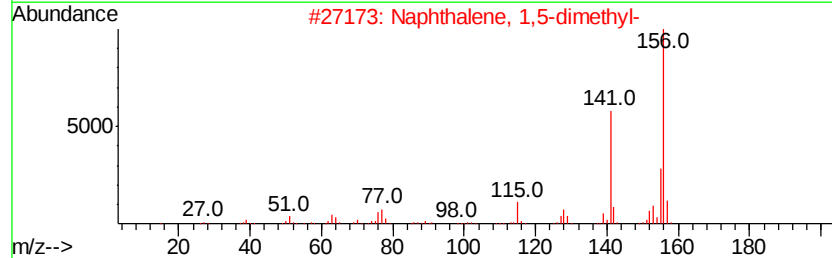
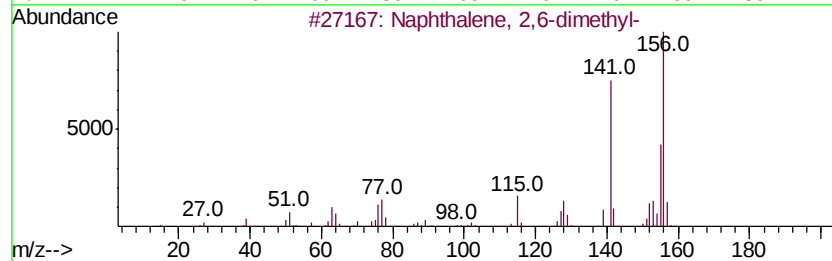
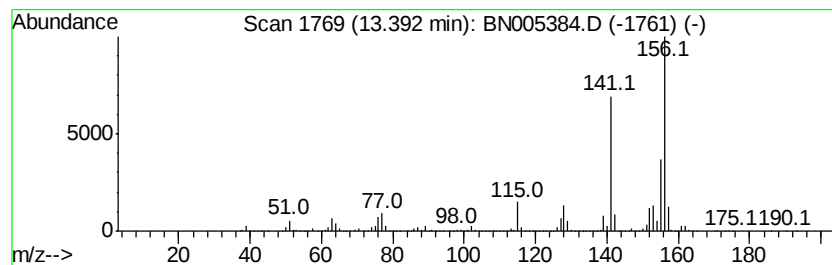
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Naphthalene, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.39	32.17 ng/ul	13596900	Acenaphthene-d10	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
2		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
3		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97
4		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
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ClientSampleId :
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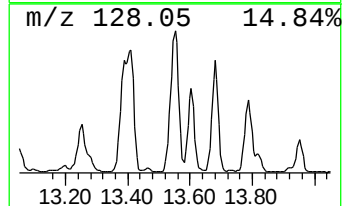
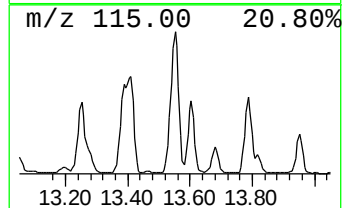
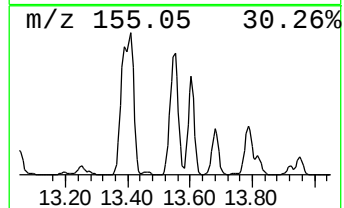
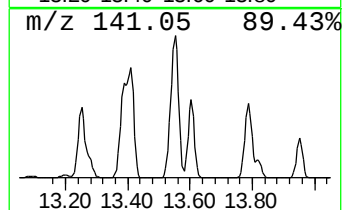
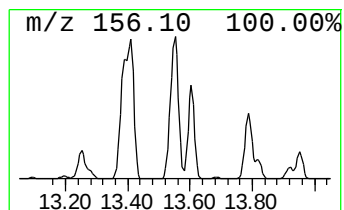
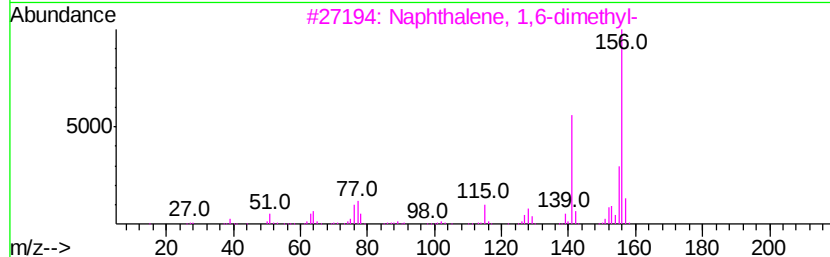
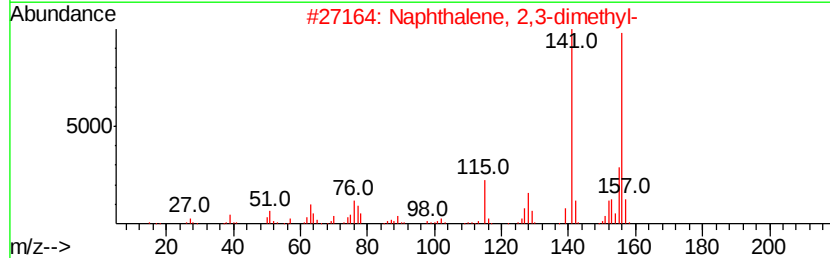
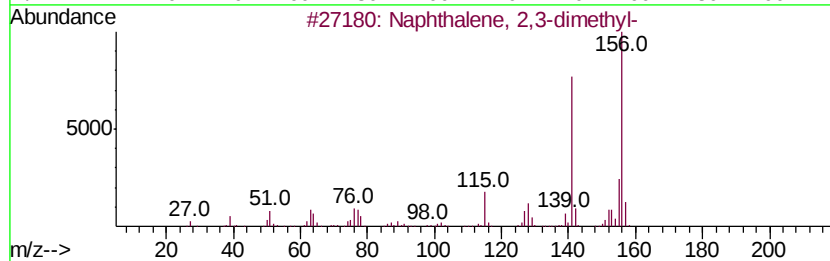
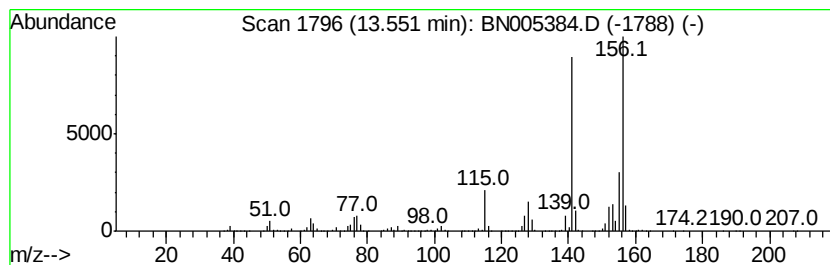
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Naphthalene, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.55	44.36 ng/ul	18748100	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

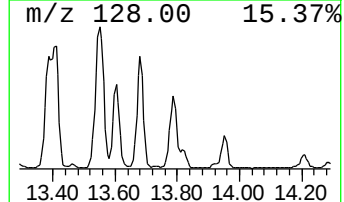
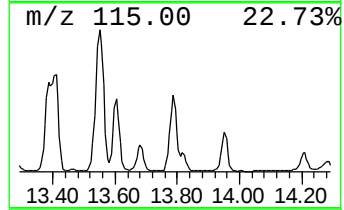
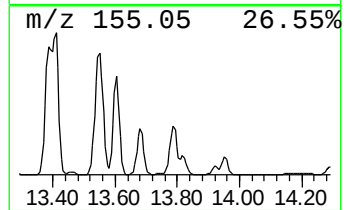
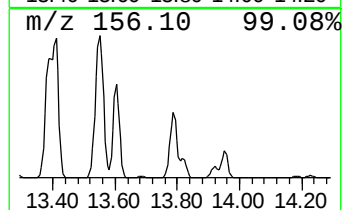
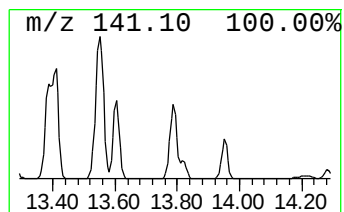
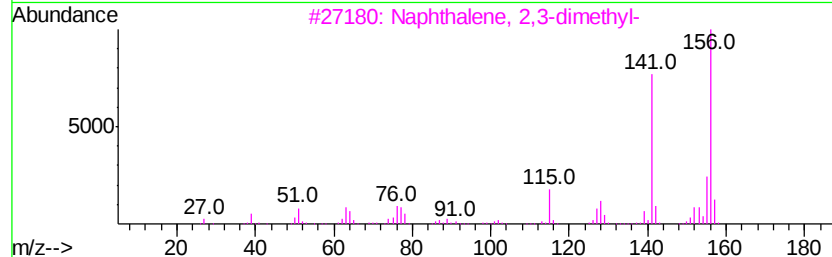
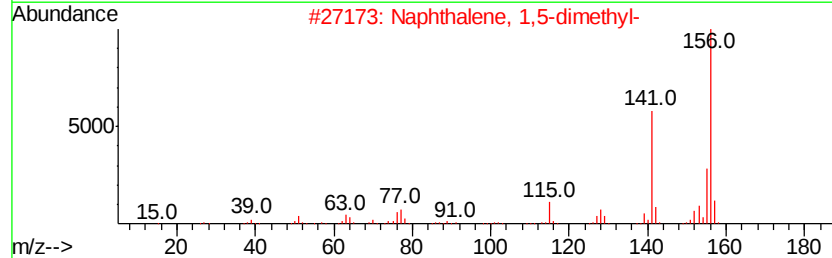
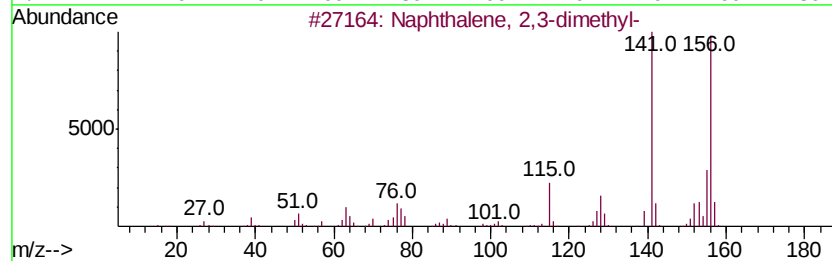
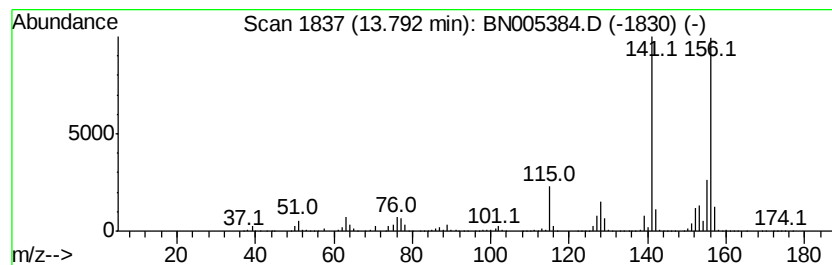
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BNA_N
ClientSampleId :
GAHXOME

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 Naphthalene, 1,5-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	18.48 ng/ul	7809560	Acenaphthene-d10	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
3		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
4		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
5		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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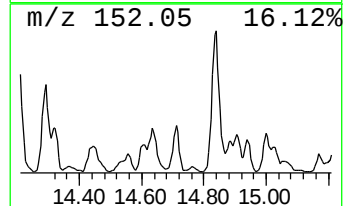
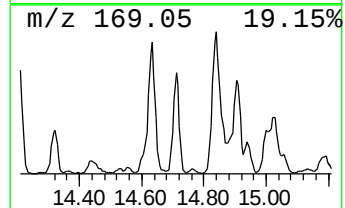
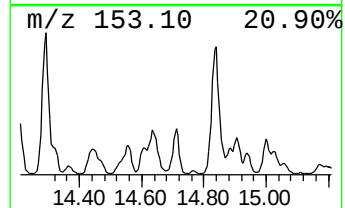
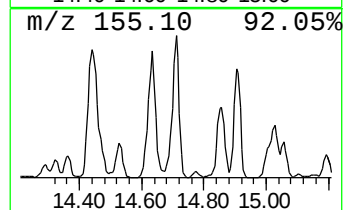
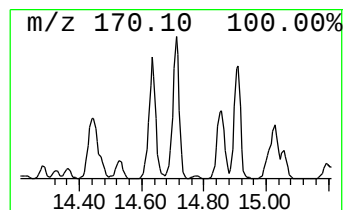
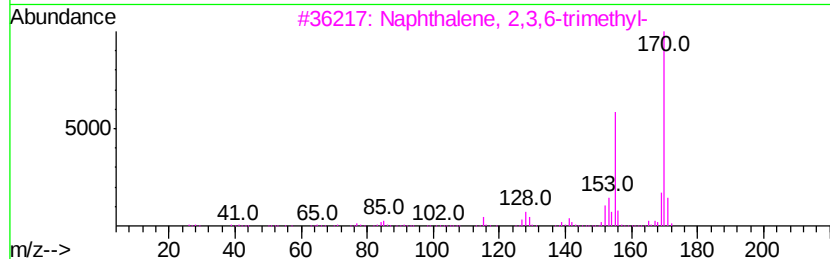
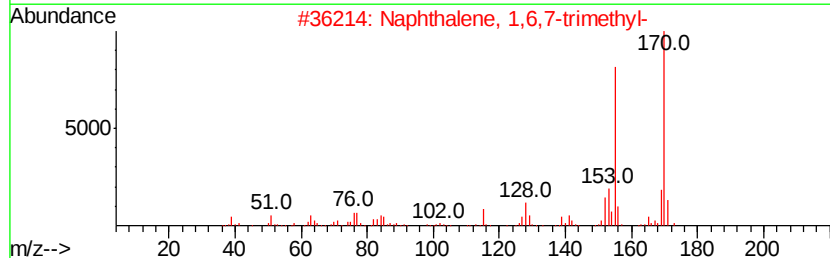
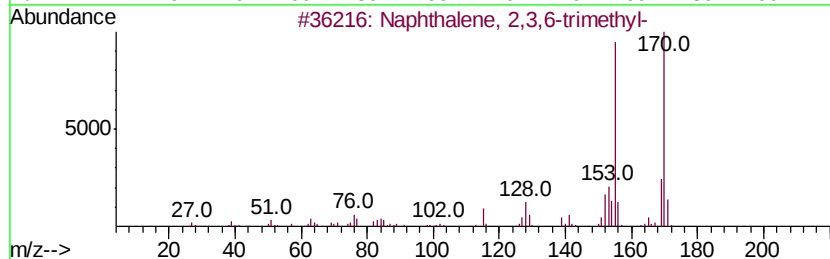
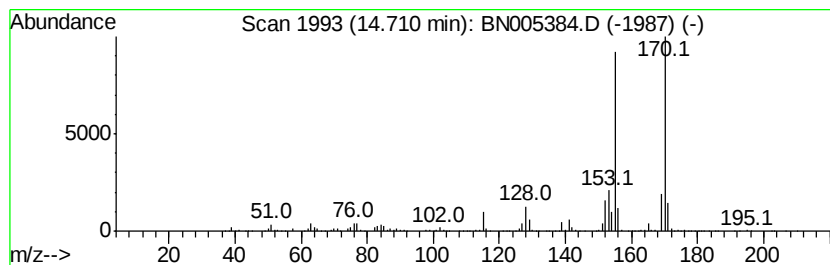
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Naphthalene, 2,3,6-trimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	7.18 ng/ul	3032270	Acenaphthene-d10	14.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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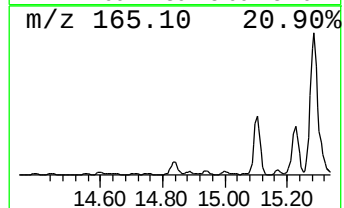
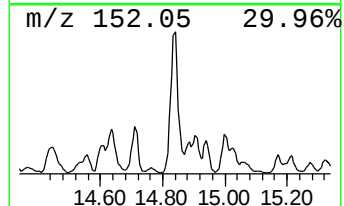
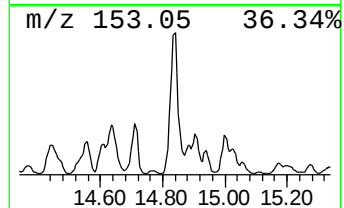
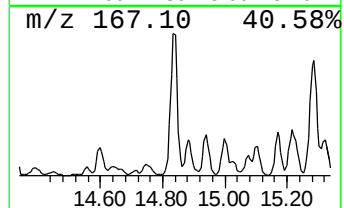
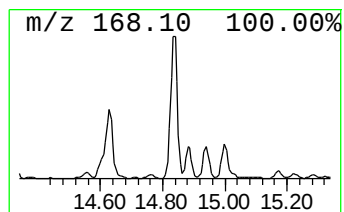
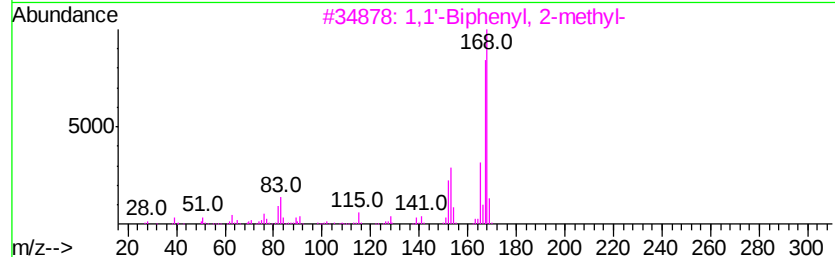
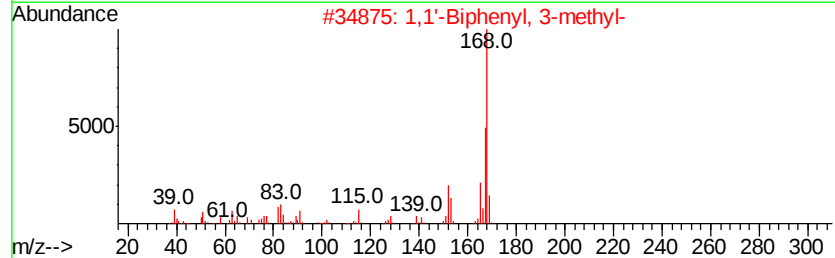
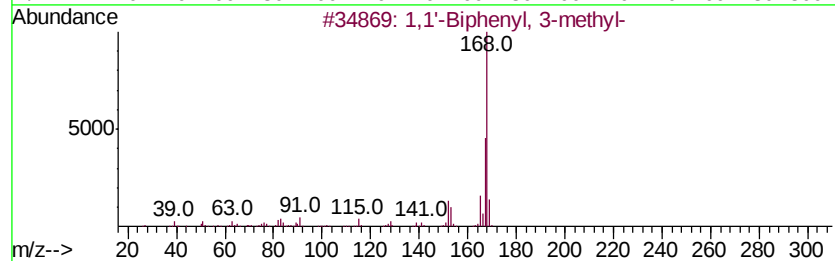
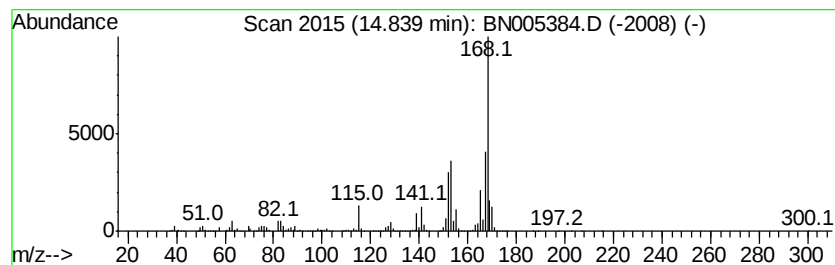
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 1,1'-Biphenyl, 3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	16.79 ng/ul	7096400	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	72
2		1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	70
3		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	53
4		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	53
5		Naphthalene, 2-(1-methylethenyl)-	168	C13H12	003710-23-4	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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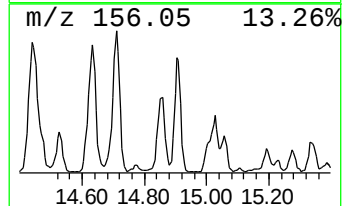
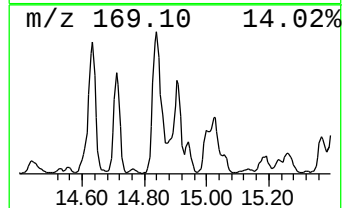
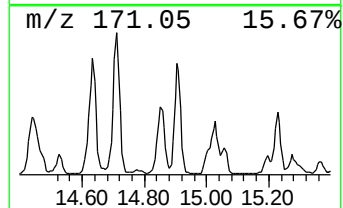
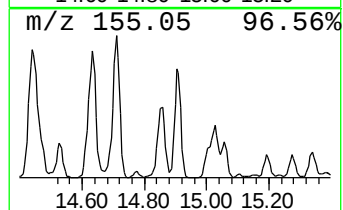
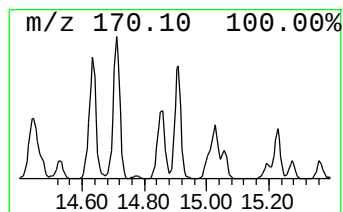
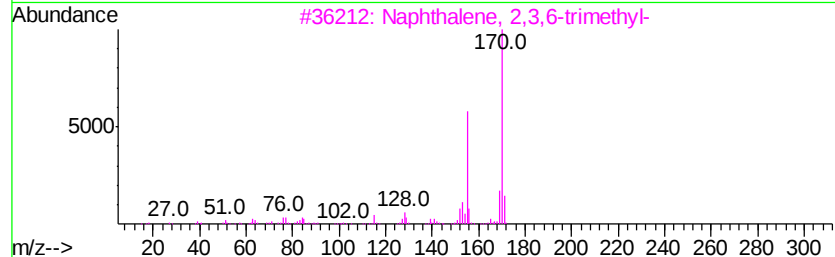
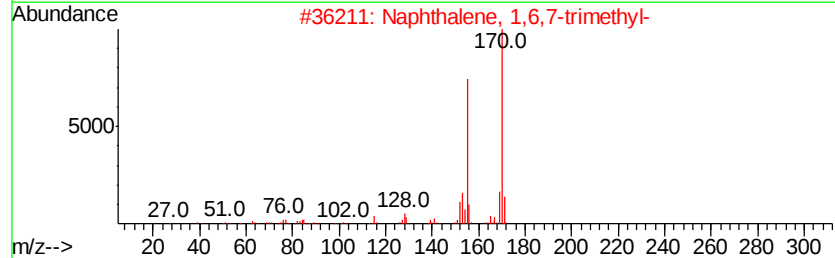
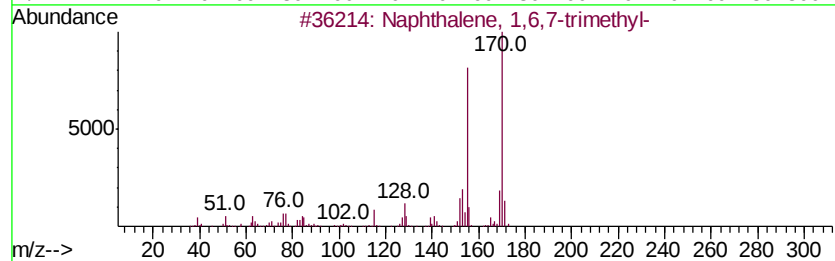
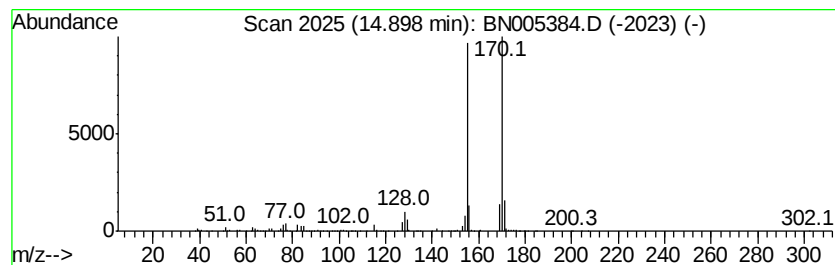
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Naphthalene, 1,6,7-trimethyl- Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	4.35 ng/ul	1837000	Acenaphthene-d10	14.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
4			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	91
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
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Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

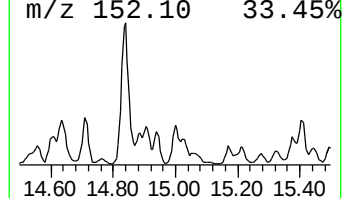
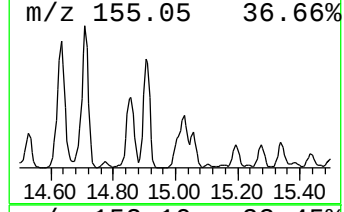
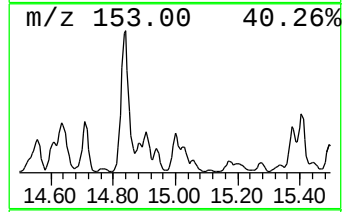
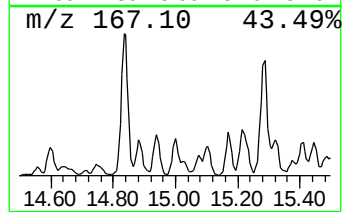
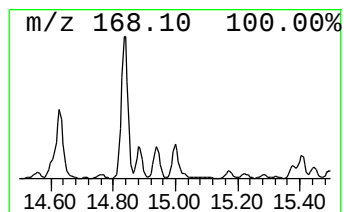
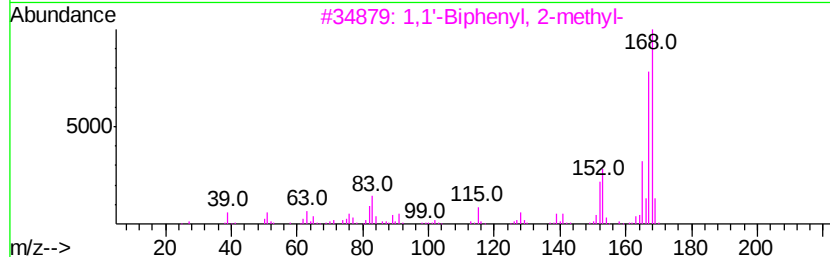
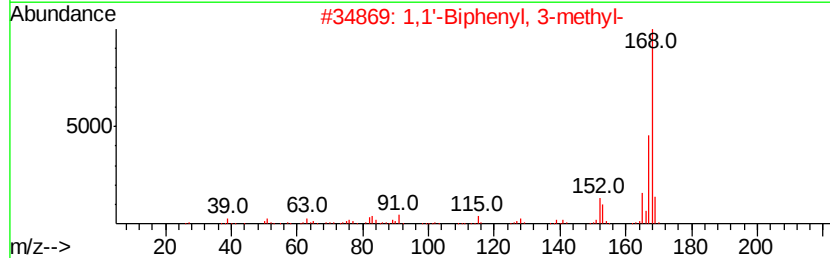
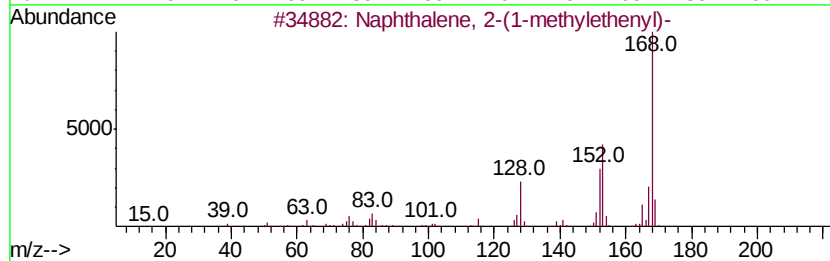
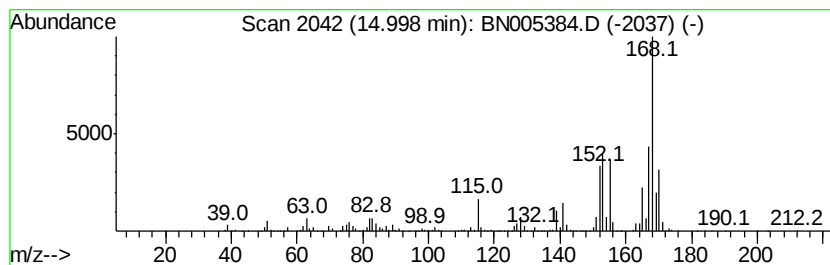
Instrument :
BNA_N
ClientSampleId :
GAHXOME

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 Naphthalene, 2-(1-methyleth... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.00	4.41 ng/ul	1863550	Acenaphthene-d10	14.23	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-(1-methylethenyl)-	168	C13H12	003710-23-4	58
2	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	47
3	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	46
4	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	46
5	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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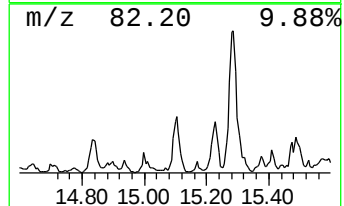
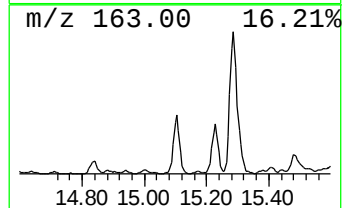
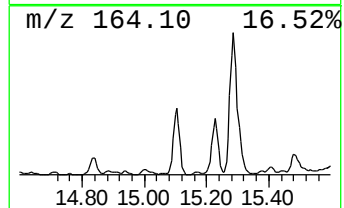
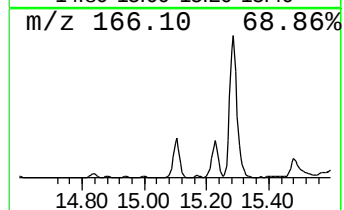
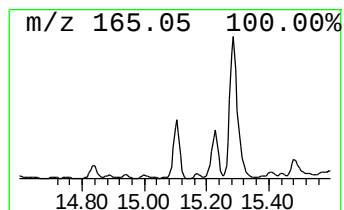
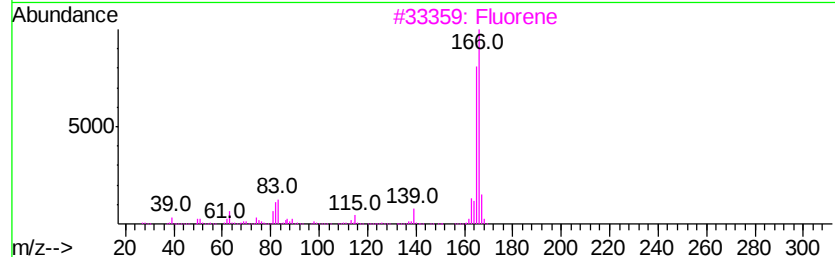
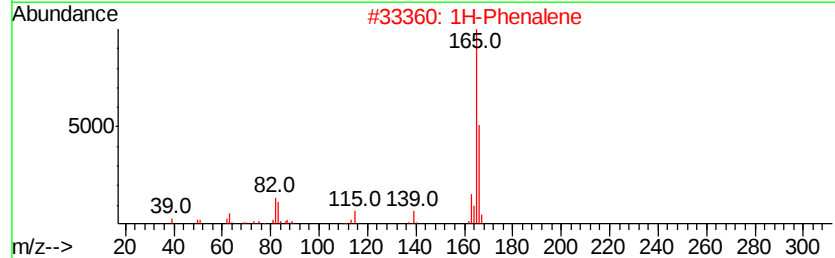
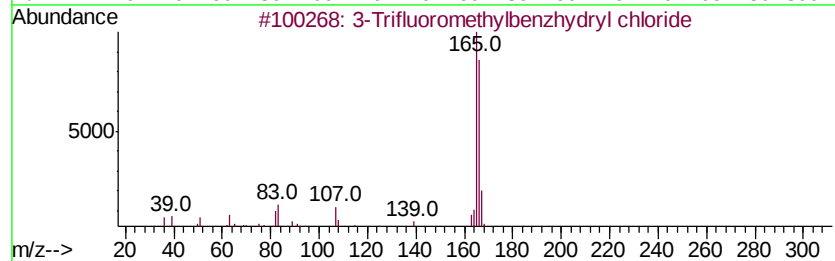
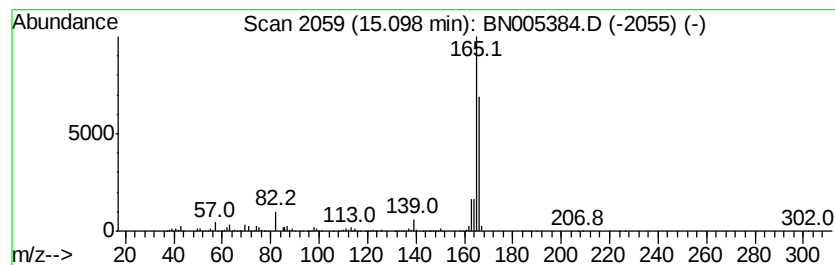
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 3-Trifluoromethylbenzhydryl... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	8.92 ng/ul	3770910	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Trifluoromethylbenzhvdrv1 chlo...	270	C14H10ClF3	067240-79-3	78
2		1H-Phenylene	166	C13H10	000203-80-5	72
3		Fluorene	166	C13H10	000086-73-7	58
4		Fluorene	166	C13H10	000086-73-7	50
5		9H-Fluorene, 9-bromo-	244	C13H9Br	001940-57-4	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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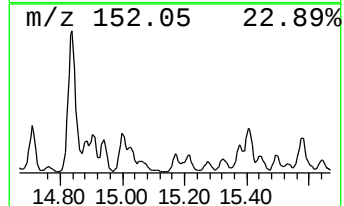
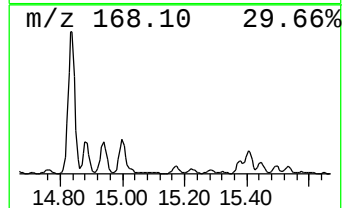
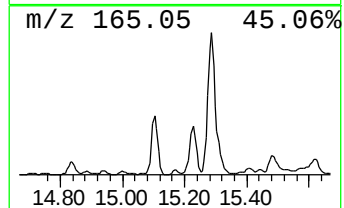
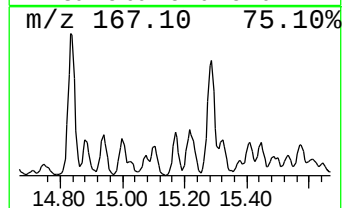
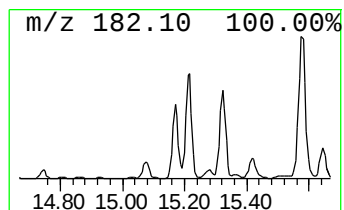
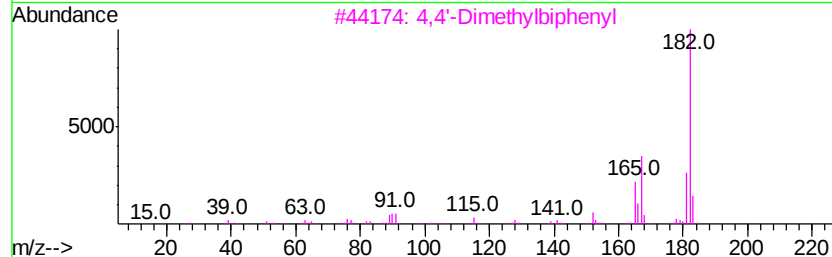
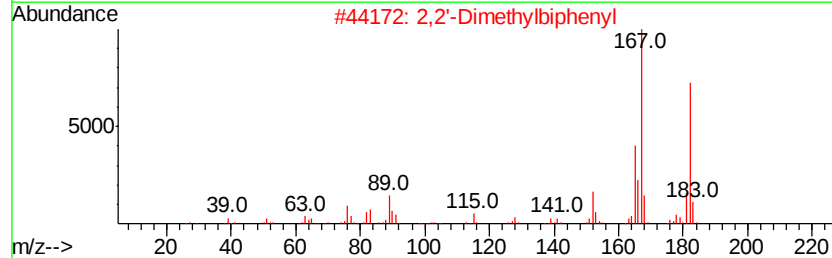
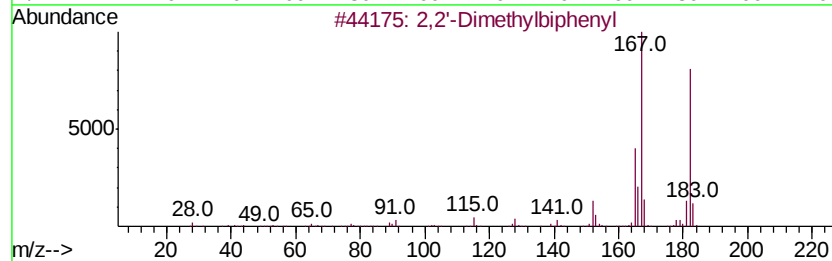
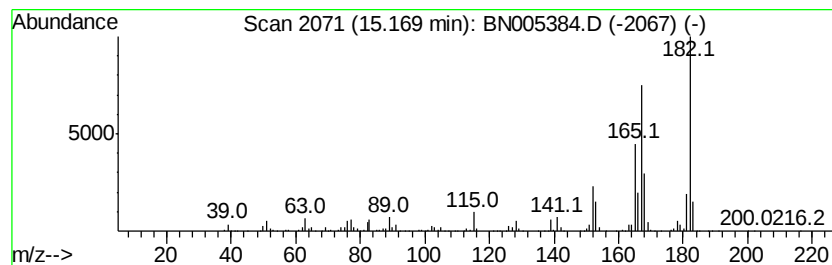
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 2,2'-Dimethylbiphenyl Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	2.61 ng/ul	1104290	Acenaphthene-d10	14.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2'-Dimethylbiphenyl	182	C14H14	000605-39-0	94
2			2,2'-Dimethylbiphenyl	182	C14H14	000605-39-0	93
3			4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	90
4			1,1'-Biphenyl, 3,4'-dimethyl-	182	C14H14	007383-90-6	81
5			4,4'-Dimethylbiphenyl	182	C14H14	000613-33-2	81



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Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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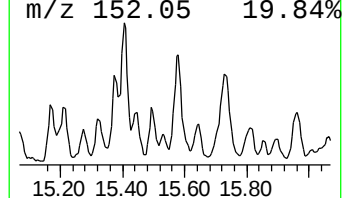
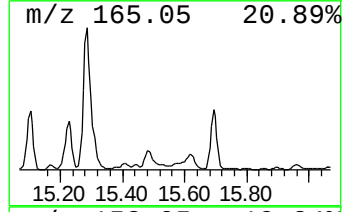
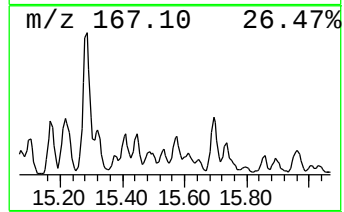
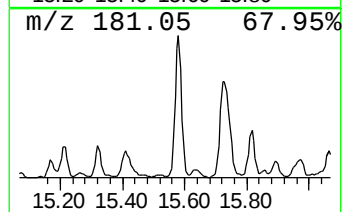
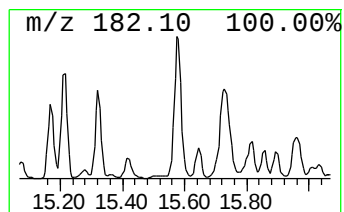
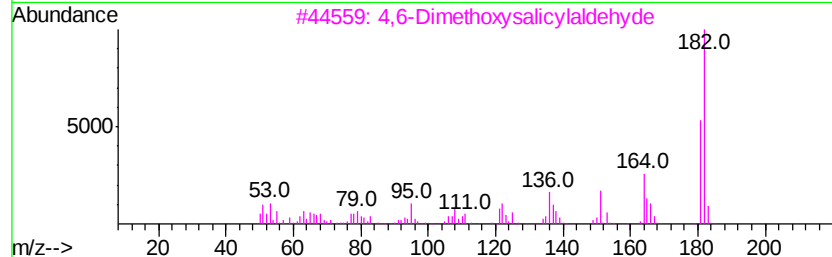
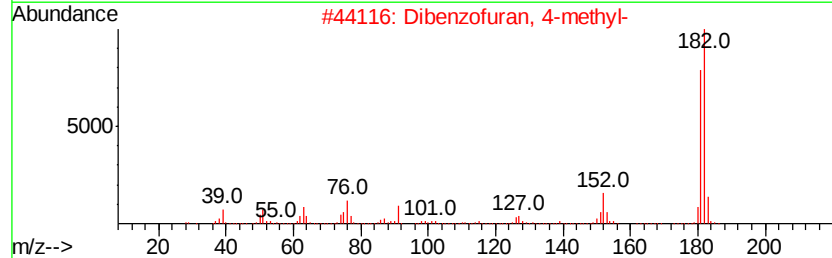
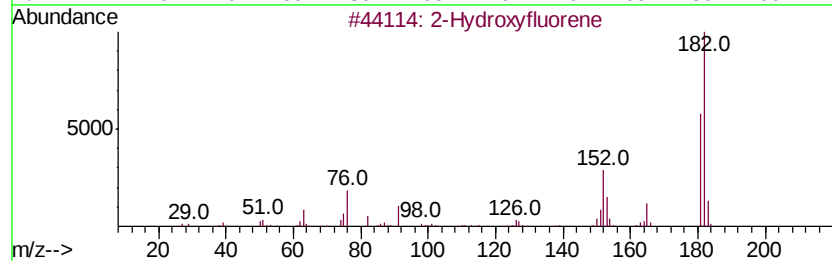
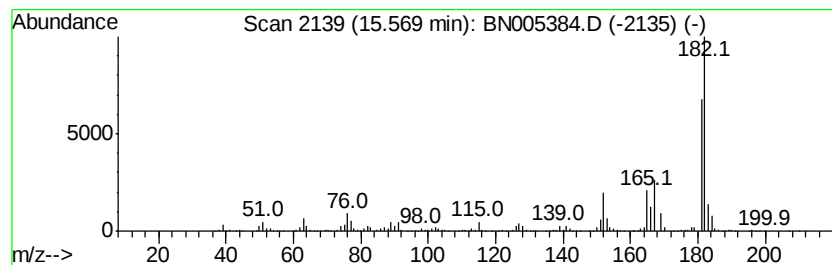
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 2-Hydroxyfluorene Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	4.63 ng/ul	1956310	Acenaphthene-d10	14.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hydroxyfluorene	182	C13H10O	002443-58-5	87
2			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	87
3			4,6-Dimethoxysalicylaldehyde	182	C9H10O4	000708-76-9	72
4			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	68
5			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GAHXOME

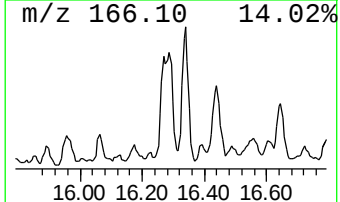
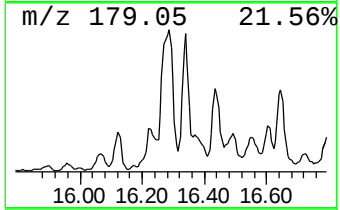
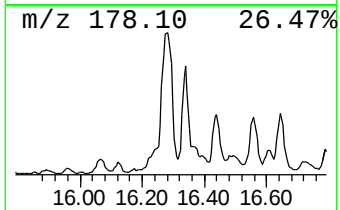
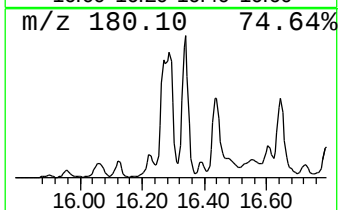
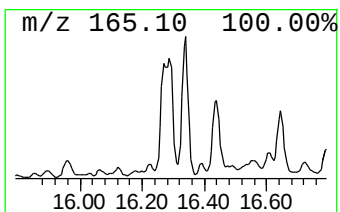
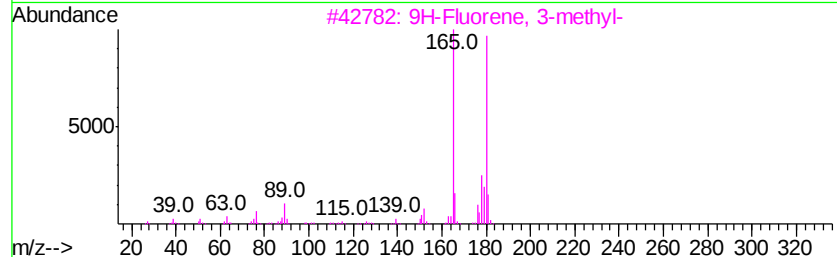
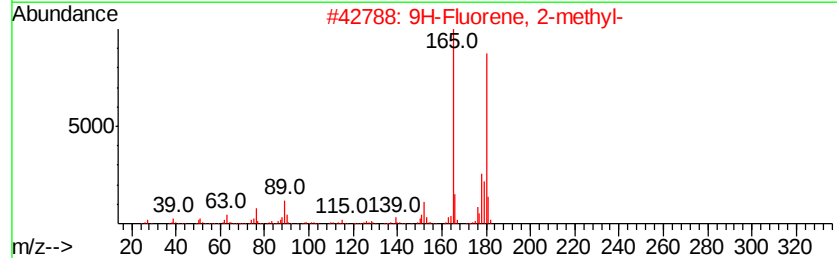
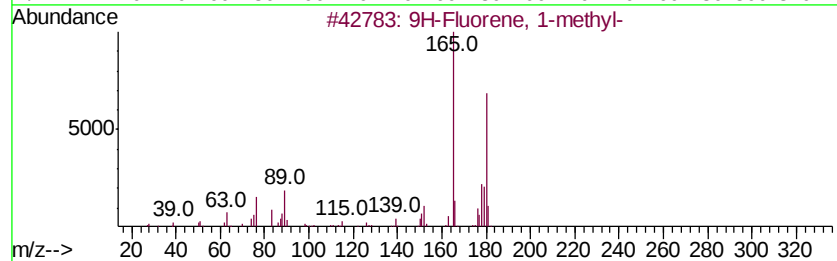
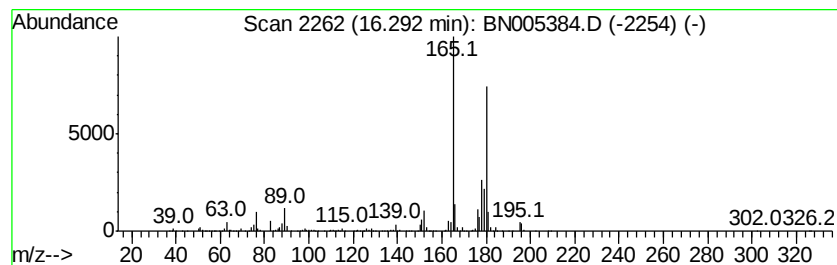
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 9H-Fluorene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.29	18.88 ng/ul	5069050	Phenanthrene-d10	16.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	96
2		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	96
3		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	93
4		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	91
5		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN043019\
Data File : BN005384.D
Acq On : 30 Apr 2019 22:32
Operator : JU/SJ
Sample : K2481-17ME
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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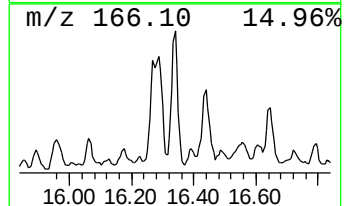
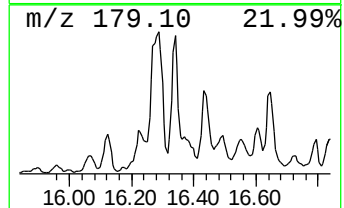
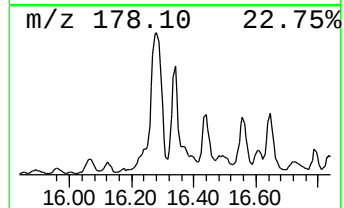
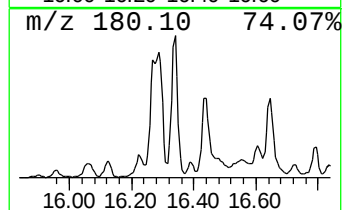
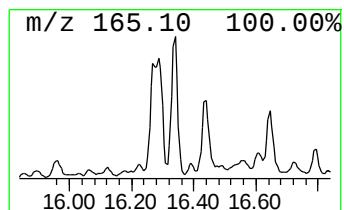
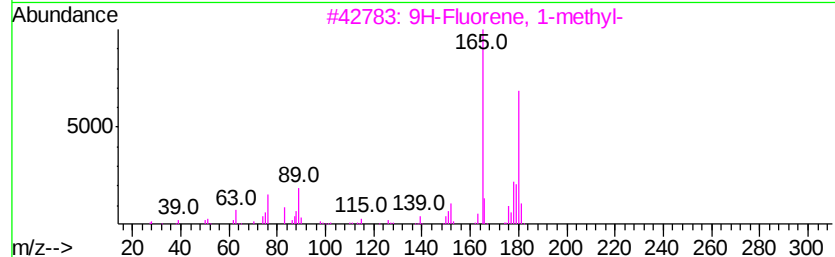
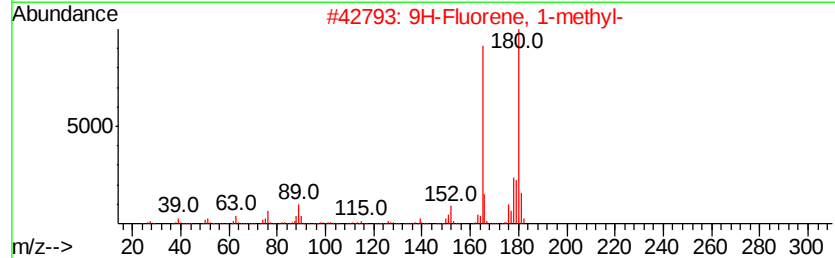
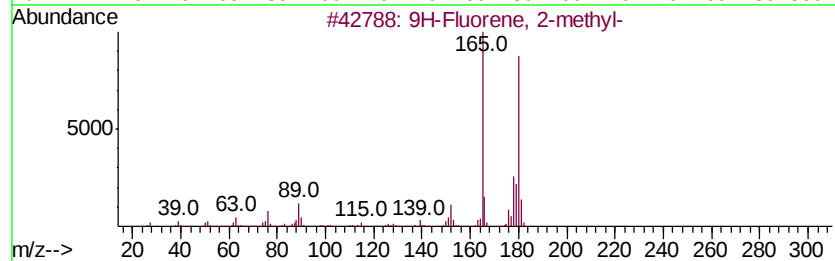
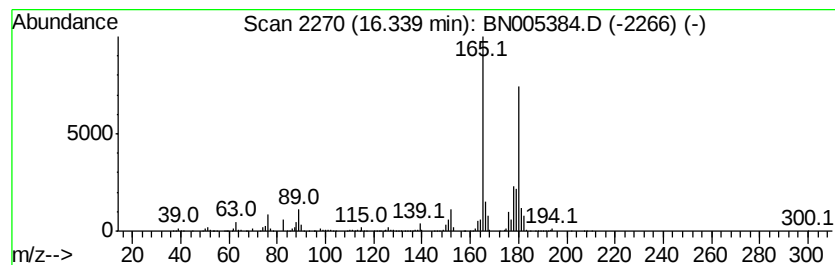
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 9H-Fluorene, 2-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	11.06 ng/ul	2968890	Phenanthrene-d10	16.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	97
2		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	96
3		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	95
4		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	95
5		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN043019\
 Data File : BN005384.D
 Acq On : 30 Apr 2019 22:32
 Operator : JU/SJ
 Sample : K2481-17ME
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-etheny...	7.25	73.5	ng/ul	7234120	1	7.59	1967200	20.0
Indene	8.12	109.6	ng/ul	10780000	1	7.59	1967200	20.0
Benzene, 2-etheny...	8.68	11.9	ng/ul	1167930	1	7.59	1967200	20.0
Benzene, (2-methy...	8.85	22.3	ng/ul	2195000	1	7.59	1967200	20.0
Benzene, 1,2,4,5-...	9.26	6.0	ng/ul	1144760	2	10.36	3825500	20.0
Benzene, 4-etheny...	9.34	8.0	ng/ul	1533940	2	10.36	3825500	20.0
1H-Indene, 1-methyl-	9.79	40.3	ng/ul	7709990	2	10.36	3825500	20.0
2-Methylindene	9.86	27.6	ng/ul	5283430	2	10.36	3825500	20.0
1H-Indene, 1,3-di...	11.40	9.6	ng/ul	1841780	2	10.36	3825500	20.0
1H-Indene, 4,7-di...	11.55	7.6	ng/ul	1451590	2	10.36	3825500	20.0
1H-Indene, 1-ethy...	11.67	5.6	ng/ul	1070720	2	10.36	3825500	20.0
(DEL) Alkane: Str...	11.80	7.8	ng/ul	1485800	2	10.36	3825500	20.0
Naphthalene, 1-me...	12.26	176.3	ng/ul	33722200	2	10.36	3825500	20.0
Naphthalene, 2-et...	13.25	16.1	ng/ul	6806810	3	14.23	8451980	20.0
Naphthalene, 2,6-...	13.39	32.2	ng/ul	13596900	3	14.23	8451980	20.0
Naphthalene, 2,3-...	13.55	44.4	ng/ul	18748100	3	14.23	8451980	20.0
Naphthalene, 1,5-...	13.79	18.5	ng/ul	7809560	3	14.23	8451980	20.0
Naphthalene, 2,3,...	14.71	7.2	ng/ul	3032270	3	14.23	8451980	20.0
1,1'-Biphenyl, 3-...	14.84	16.8	ng/ul	7096400	3	14.23	8451980	20.0
Naphthalene, 1,6,...	14.90	4.3	ng/ul	1837000	3	14.23	8451980	20.0
Naphthalene, 2-(1...	15.00	4.4	ng/ul	1863550	3	14.23	8451980	20.0
3-Trifluoromethyl...	15.10	8.9	ng/ul	3770910	3	14.23	8451980	20.0
2,2'-Dimethylbiph...	15.17	2.6	ng/ul	1104290	3	14.23	8451980	20.0
2-Hydroxyfluorene	15.57	4.6	ng/ul	1956310	3	14.23	8451980	20.0
9H-Fluorene, 1-me...	16.29	18.9	ng/ul	5069050	4	16.98	5369650	20.0
9H-Fluorene, 2-me...	16.34	11.1	ng/ul	2968890	4	16.98	5369650	20.0