

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
 Data File : BN005470.D
 Acq On : 04 May 2019 02:26
 Operator : JU/SJ
 Sample : K2603-15DL 25X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GAJ67DL

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	3.911	151	157	176	rVB	2783720	4423298	3.58%	0.250%
2	5.293	386	392	407	rVB	3885671	7899808	6.39%	0.446%
3	5.628	442	449	461	rBV2	5889783	12723047	10.29%	0.719%
4	7.252	718	725	733	rVV2	8211252	15347457	12.42%	0.867%
5	7.334	733	739	748	rVB	2218614	3533107	2.86%	0.200%
6	8.122	862	873	880	rBV	16600262	28832681	23.33%	1.628%
7	8.846	992	996	1005	rVB	1869198	3397336	2.75%	0.192%
8	9.793	1147	1157	1163	rBV3	4853003	11441126	9.26%	0.646%
9	9.863	1163	1169	1173	rBV	4177894	8093711	6.55%	0.457%
10	10.351	1246	1252	1255	rBV2	1923135	3374774	2.73%	0.191%
11	10.451	1255	1269	1274	rVV2	38840559	123605236	100.00%	6.981%
12	11.393	1418	1429	1434	rBV2	1436901	3658940	2.96%	0.207%
13	11.757	1485	1491	1495	rBV	2173457	3872572	3.13%	0.219%
14	12.057	1530	1542	1546	rBV2	36394010	85357382	69.06%	4.821%
15	12.269	1563	1578	1584	rBV	28971421	59880575	48.45%	3.382%
16	13.057	1705	1712	1725	rBV2	12211448	20153179	16.30%	1.138%
17	13.251	1739	1745	1757	rVB	3220854	8037524	6.50%	0.454%
18	13.392	1761	1769	1770	rBV	9718661	15418169	12.47%	0.871%
19	13.557	1788	1797	1802	rBV	16283311	31930016	25.83%	1.803%
20	13.610	1802	1806	1810	rVV	10346753	14119072	11.42%	0.797%
21	13.687	1814	1819	1824	rVB	12256951	18240037	14.76%	1.030%
22	13.792	1830	1837	1841	rBV	8357810	14673737	11.87%	0.829%
23	13.969	1855	1867	1872	rVB3	30731962	60083757	48.61%	3.394%
24	14.151	1893	1898	1901	rBV	3123785	4149347	3.36%	0.234%
25	14.210	1902	1908	1916	rBV2	6448302	11570715	9.36%	0.654%
26	14.292	1917	1922	1925	rBV2	4677026	7697878	6.23%	0.435%
27	14.445	1942	1948	1957	rVB	2209203	4845680	3.92%	0.274%
28	14.634	1970	1980	1988	rBV3	6589604	13902201	11.25%	0.785%
29	14.710	1988	1993	1998	rBV	4940664	6624343	5.36%	0.374%
30	14.845	2009	2016	2021	rBV3	6704322	14666301	11.87%	0.828%
31	14.910	2024	2027	2030	rVB	3174044	3528193	2.85%	0.199%
32	15.004	2038	2043	2045	rBV	2800353	4095203	3.31%	0.231%
33	15.110	2056	2061	2066	rVB	10151261	13879056	11.23%	0.784%
34	15.228	2077	2081	2086	rVB2	7826741	12591927	10.19%	0.711%

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35	15.298	2086	2093	2103	rVB	22911016	45801477	37.05%	2.587%
36	15.410	2108	2112	2115	rVB2	2809918	3360266	2.72%	0.190%
37	15.492	2121	2126	2135	rVB4	2772520	6986557	5.65%	0.395%
38	15.581	2136	2141	2145	rBV2	3381367	5967756	4.83%	0.337%
39	15.698	2156	2161	2164	rBV	6596680	9342698	7.56%	0.528%
40	15.969	2201	2207	2210	rBV2	4957108	7889143	6.38%	0.446%
41	16.292	2255	2262	2267	rVV2	7429559	16250892	13.15%	0.918%
42	16.345	2267	2271	2275	rVB	5343882	7116433	5.76%	0.402%
43	16.439	2282	2287	2293	rVB	4731212	7195674	5.82%	0.406%
44	16.645	2319	2322	2329	rVB2	2701836	4075373	3.30%	0.230%
45	16.769	2339	2343	2345	rVV	2702261	3578591	2.90%	0.202%
46	16.804	2345	2349	2360	rVB	9692469	16109652	13.03%	0.910%
47	17.063	2376	2393	2397	rBV2	43093916	116329481	94.11%	6.570%
48	17.098	2397	2399	2401	rVV	3161718	3674690	2.97%	0.208%
49	17.139	2401	2406	2410	rVV	23607572	33626417	27.20%	1.899%
50	17.180	2410	2413	2418	rVB2	2362606	3901511	3.16%	0.220%
51	17.510	2465	2469	2473	rVB	6232511	7624175	6.17%	0.431%
52	17.569	2475	2479	2485	rVB2	6167868	9357890	7.57%	0.529%
53	17.710	2496	2503	2510	rVB3	4345175	10514087	8.51%	0.594%
54	17.875	2525	2531	2535	rVV	19133815	31690767	25.64%	1.790%
55	17.927	2535	2540	2548	rVV	22228962	33393467	27.02%	1.886%
56	18.004	2548	2553	2558	rVV	8069796	12931878	10.46%	0.730%
57	18.075	2558	2565	2569	rVV2	28034105	56692193	45.87%	3.202%
58	18.110	2569	2571	2578	rVB	15812474	17022221	13.77%	0.961%
59	18.374	2611	2616	2620	rBV	11980796	17193982	13.91%	0.971%
60	18.427	2622	2625	2629	rVB	3736127	4589587	3.71%	0.259%
61	18.622	2655	2658	2663	rVV3	3434505	5653006	4.57%	0.319%
62	18.674	2663	2667	2670	rVV	4655915	6670100	5.40%	0.377%
63	18.710	2670	2673	2678	rVB3	2933350	4239166	3.43%	0.239%
64	18.798	2683	2688	2693	rVV2	11646499	21236507	17.18%	1.199%
65	18.863	2693	2699	2702	rVV3	8238207	18538267	15.00%	1.047%
66	18.892	2702	2704	2709	rVV	6272207	8653263	7.00%	0.489%
67	18.957	2709	2715	2721	rVV4	5398157	13559287	10.97%	0.766%
68	19.086	2728	2737	2741	rBV2	34982736	68016729	55.03%	3.842%
69	19.139	2742	2746	2749	rVV	2515661	3560668	2.88%	0.201%
70	19.227	2755	2761	2765	rVB	18125325	26261694	21.25%	1.483%
71	19.463	2787	2801	2805	rBV2	40235359	98429300	79.63%	5.559%

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72	19.510	2805	2809	2815	rVB3	2742036	4610142	3.73%	0.260%
73	19.610	2822	2826	2833	rVB2	1988284	3499936	2.83%	0.198%
74	19.774	2847	2854	2862	rVB	6850297	14952861	12.10%	0.845%
75	19.904	2870	2876	2878	rBV3	6612340	11601388	9.39%	0.655%
76	19.939	2879	2882	2886	rVB	13973131	19205274	15.54%	1.085%
77	20.045	2895	2900	2905	rBV	11542580	15952201	12.91%	0.901%
78	20.104	2905	2910	2913	rBV	9247019	11842060	9.58%	0.669%
79	20.192	2920	2925	2928	rBV4	1920810	3373622	2.73%	0.191%
80	20.239	2929	2933	2937	rVV	7332915	10876730	8.80%	0.614%
81	20.286	2937	2941	2949	rVB	10789058	15217593	12.31%	0.859%
82	20.857	3035	3038	3042	rVV	4718087	6259727	5.06%	0.354%
83	20.898	3042	3045	3048	rVB	6673665	7359062	5.95%	0.416%
84	20.939	3048	3052	3055	rBV	4296461	5270161	4.26%	0.298%
85	21.210	3092	3098	3103	rBV	27351032	48419315	39.17%	2.735%
86	21.268	3104	3108	3112	rVB	21498163	28873124	23.36%	1.631%
87	21.380	3123	3127	3131	rBV3	1960053	3930216	3.18%	0.222%
88	21.427	3131	3135	3148	rVB3	5912152	9669861	7.82%	0.546%
89	21.768	3187	3193	3198	rVV	6944028	12524512	10.13%	0.707%
90	21.821	3198	3202	3206	rVB	3981552	5369605	4.34%	0.303%
91	21.963	3222	3226	3228	rBV2	3069651	4060453	3.29%	0.229%
92	21.992	3229	3231	3238	rVB2	4416636	5657162	4.58%	0.320%
93	22.780	3358	3365	3369	rBV	9554862	22426287	18.14%	1.267%
94	22.939	3386	3392	3398	rVB	3924278	6825865	5.52%	0.386%
95	23.239	3436	3443	3448	rBV	7341900	13833142	11.19%	0.781%
96	23.345	3454	3461	3466	rVB	12596747	24143045	19.53%	1.364%
97	23.468	3478	3482	3490	rVB2	2364660	4803072	3.89%	0.271%
98	24.268	3612	3618	3630	rVB	1991517	4829785	3.91%	0.273%
99	25.615	3838	3847	3856	rVB	4326168	12213526	9.88%	0.690%
100	26.286	3952	3961	3977	rVB	3257113	10267112	8.31%	0.580%

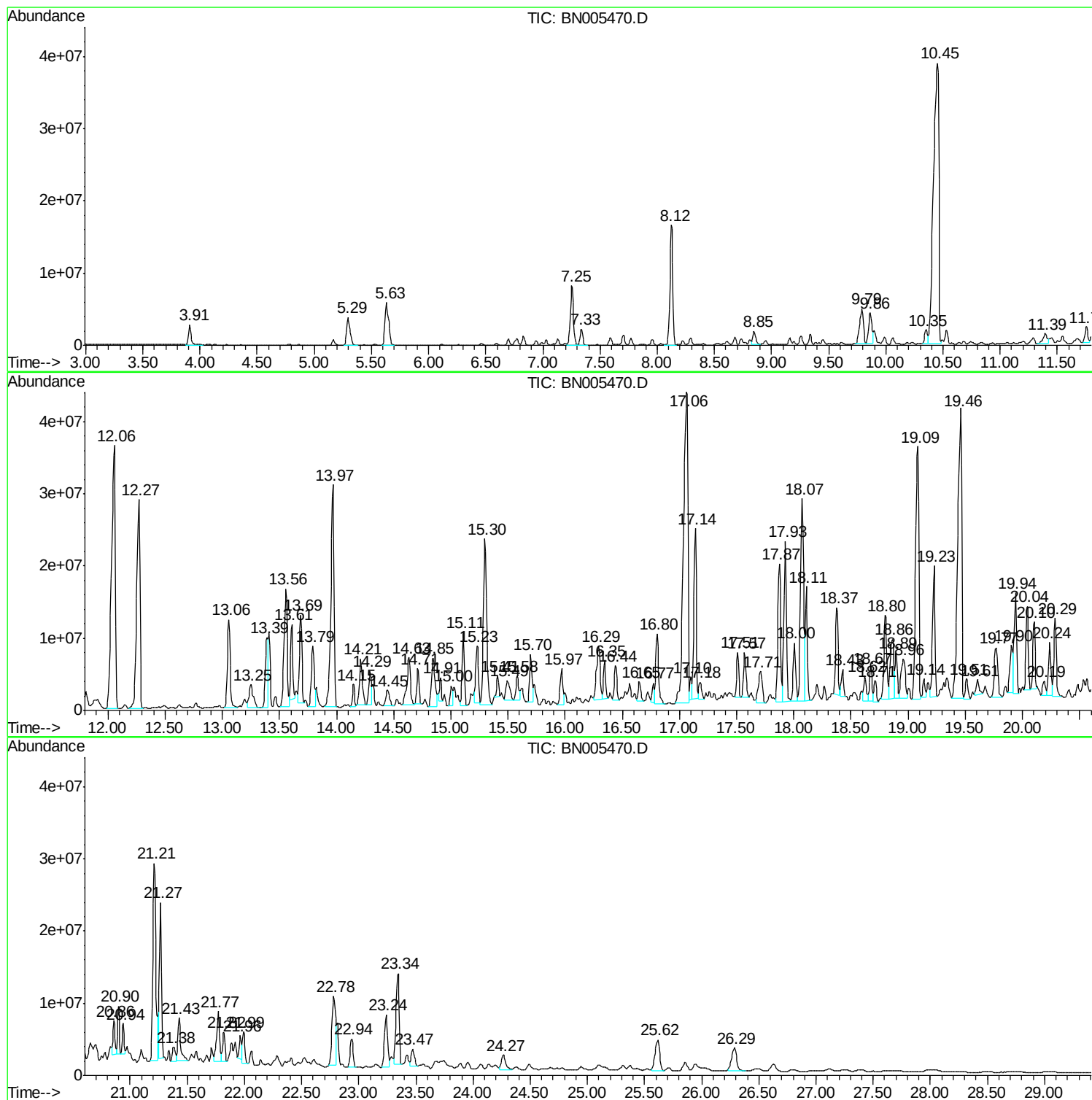
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Instrument :
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ClientSampleId :
GAJ67DL

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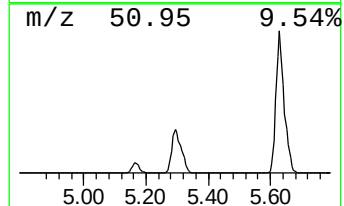
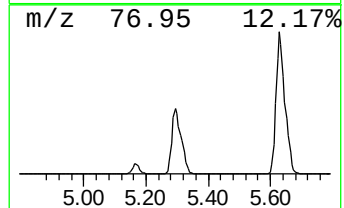
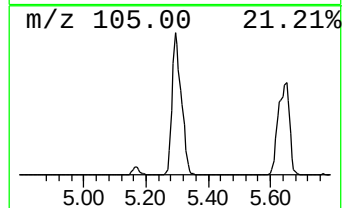
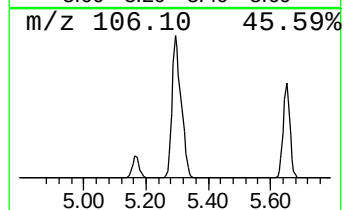
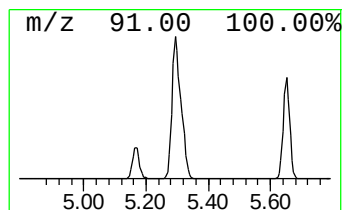
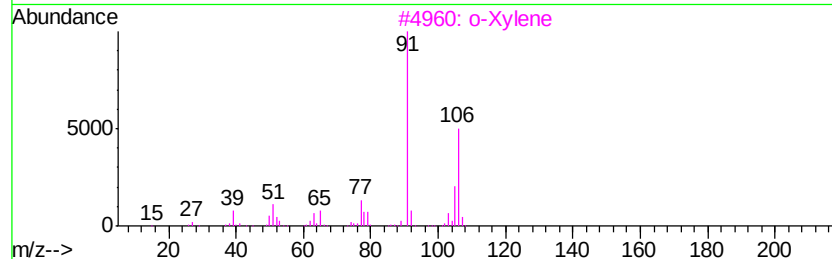
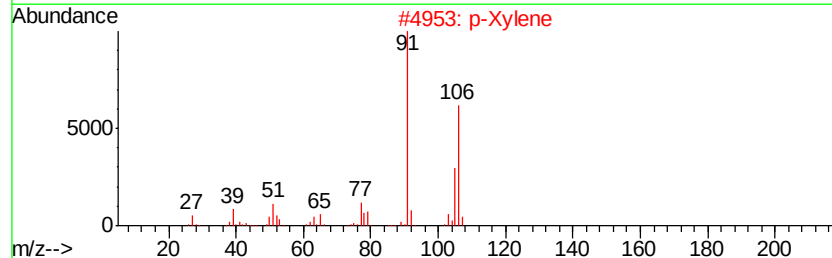
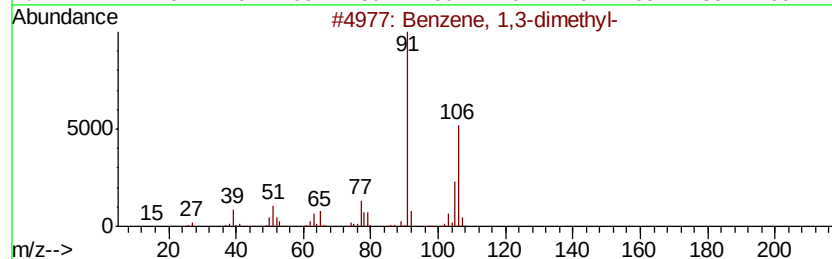
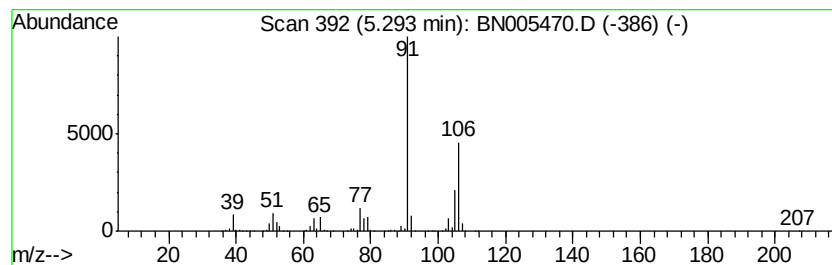
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.29	44.72 ng/ul	7899810	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		p-Xylene	106	C8H10	000106-42-3	97
3		o-Xylene	106	C8H10	000095-47-6	97
4		p-Xylene	106	C8H10	000106-42-3	97
5		p-Xylene	106	C8H10	000106-42-3	95



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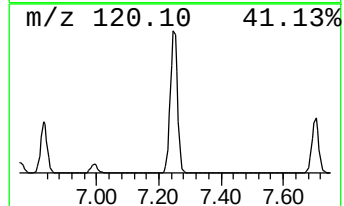
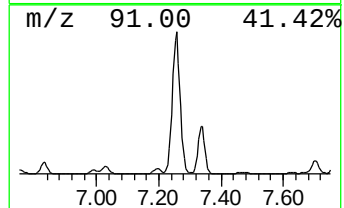
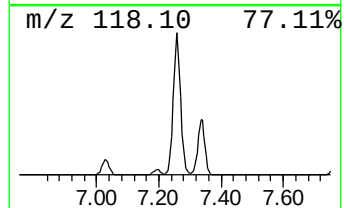
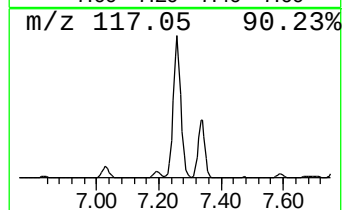
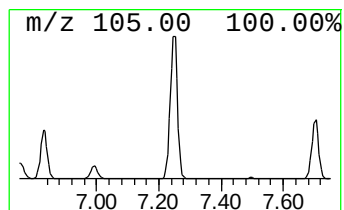
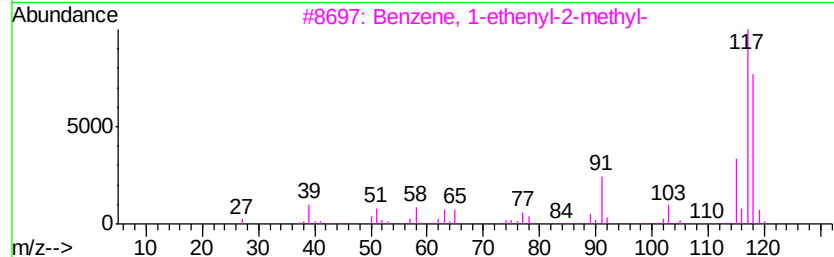
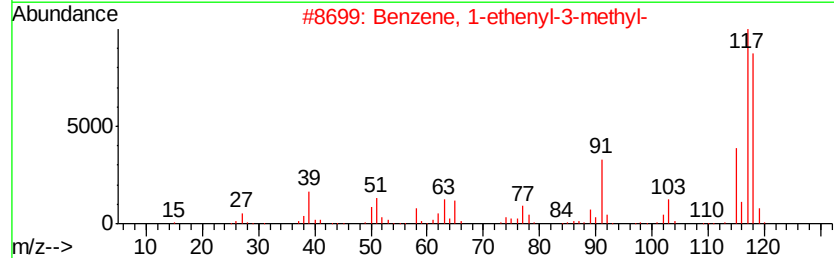
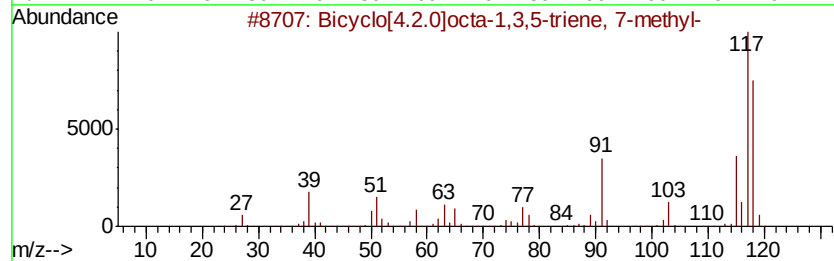
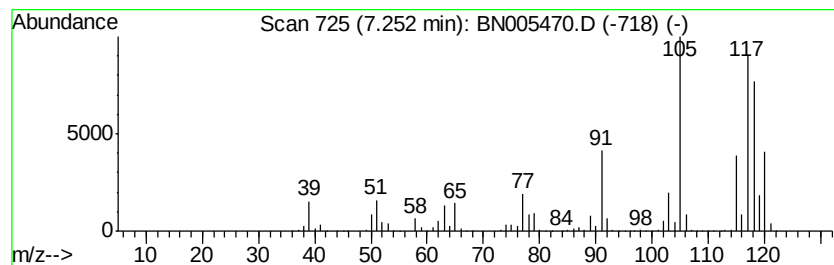
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.25	86.88 ng/ul	15347500	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	64
2		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	60
3		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	58
5		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	55



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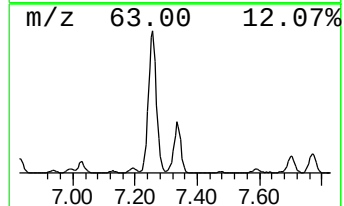
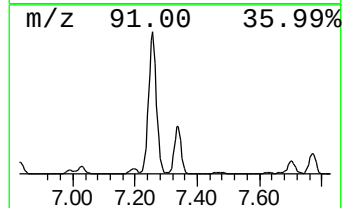
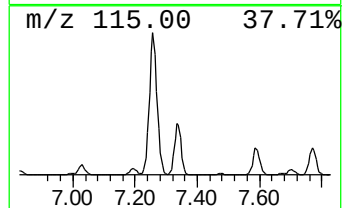
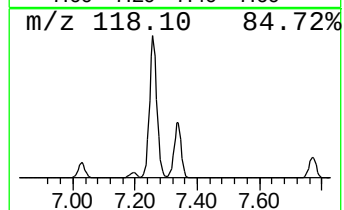
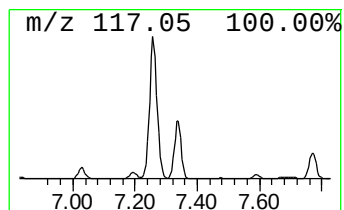
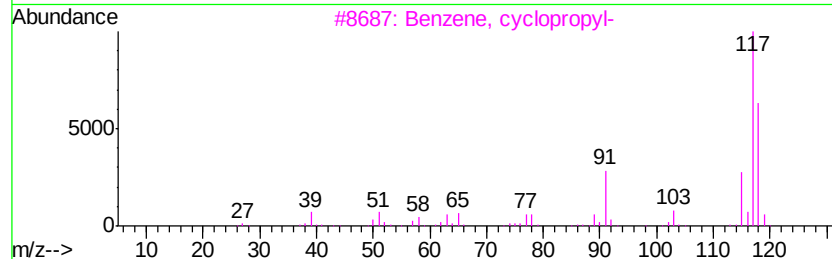
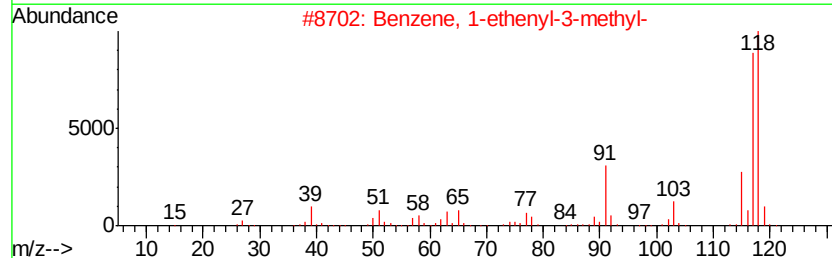
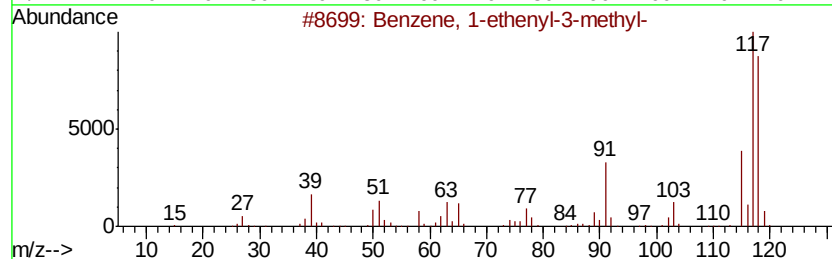
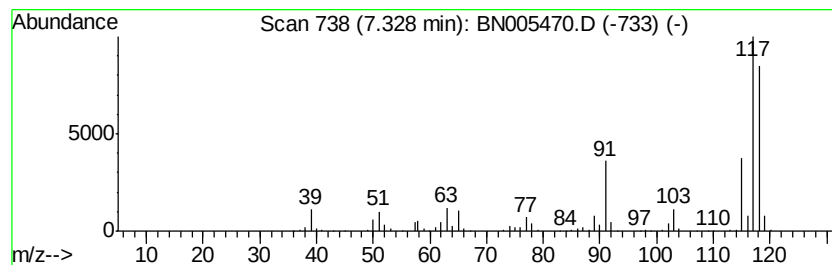
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-ethenyl-3-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.33	20.00 ng/ul	3533110	1,4-Dichlorobenzene-d4	7.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
5			Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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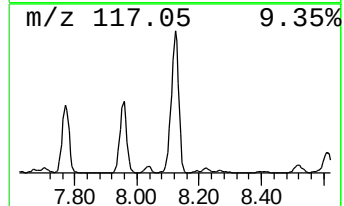
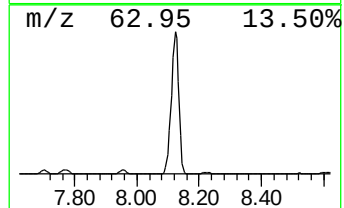
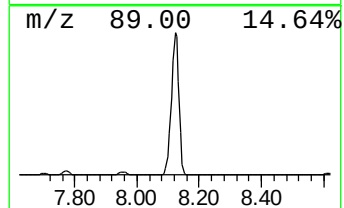
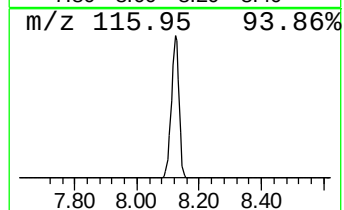
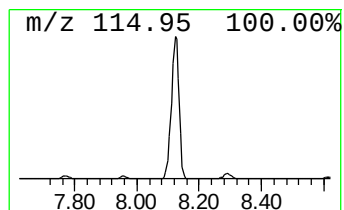
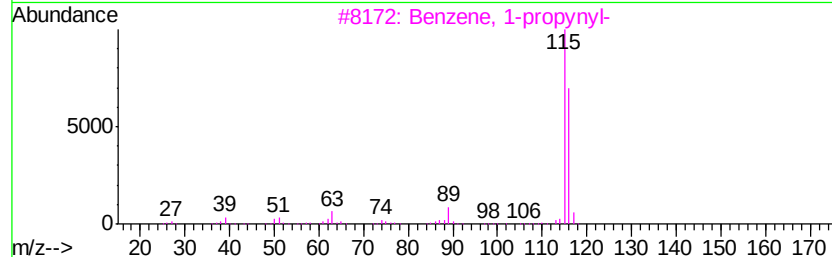
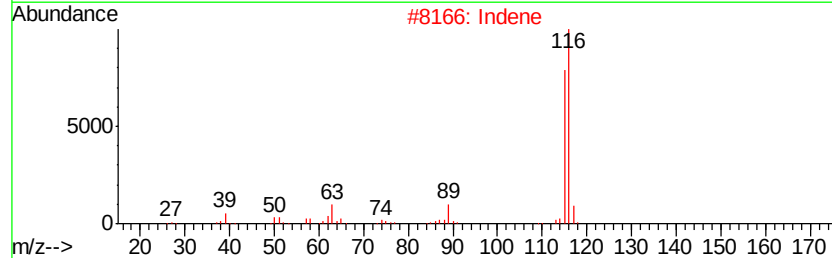
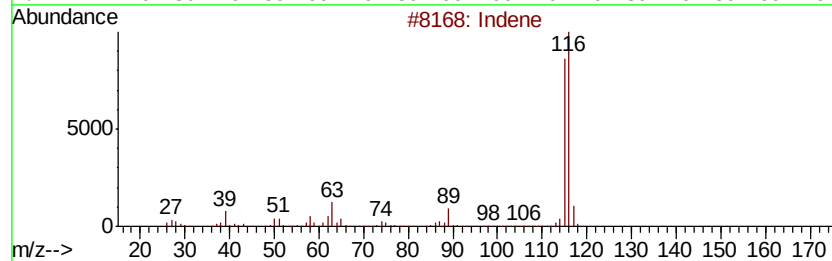
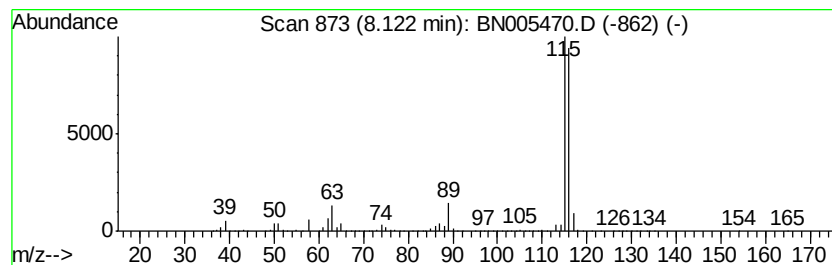
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	163.21 ng/ul	28832700	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	97
2		Indene	116	C9H8	000095-13-6	95
3		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



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Acq On : 04 May 2019 02:26
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Misc :
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ClientSampleId :
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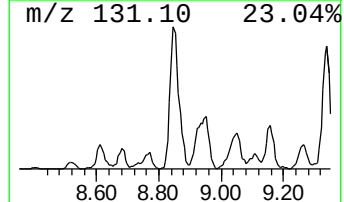
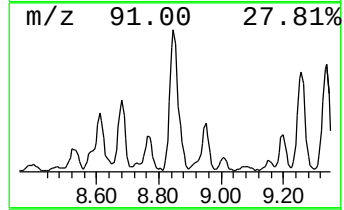
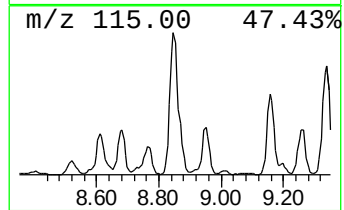
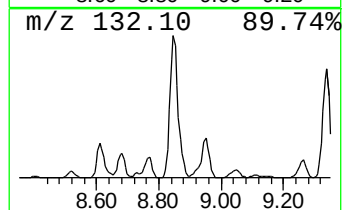
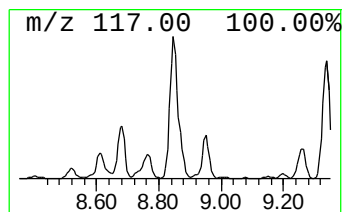
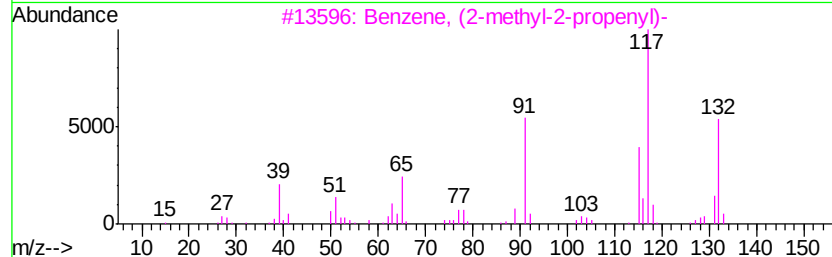
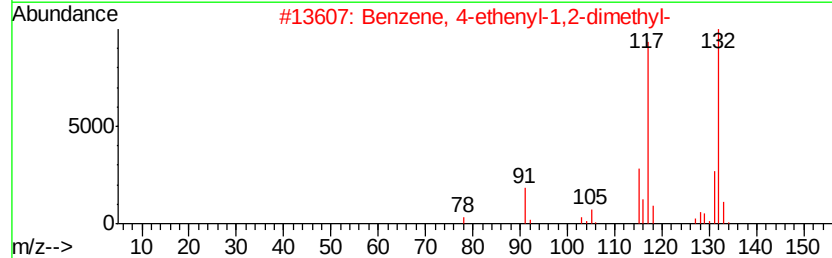
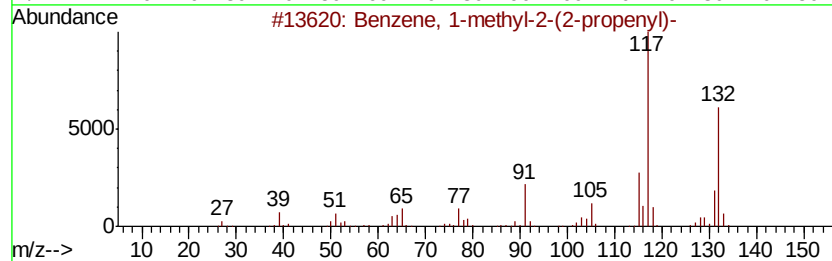
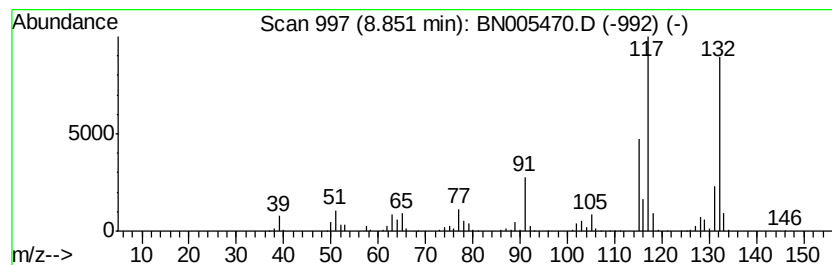
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.85	19.23 ng/ul	3397340	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	95
2		Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	95
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	95
4		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	95
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
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Instrument :
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ClientSampleId :
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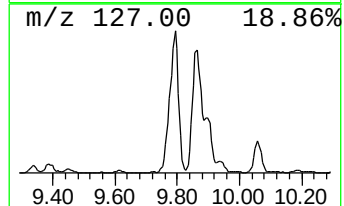
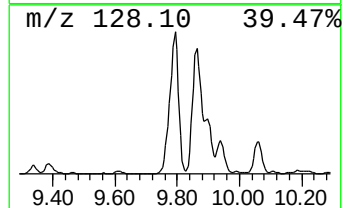
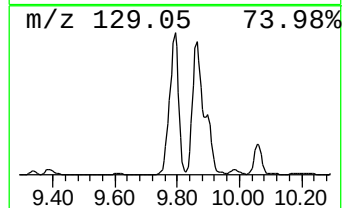
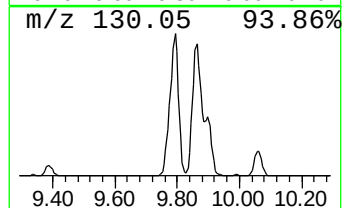
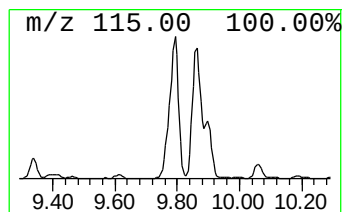
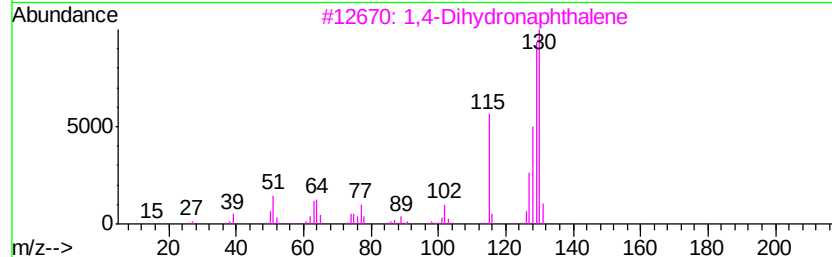
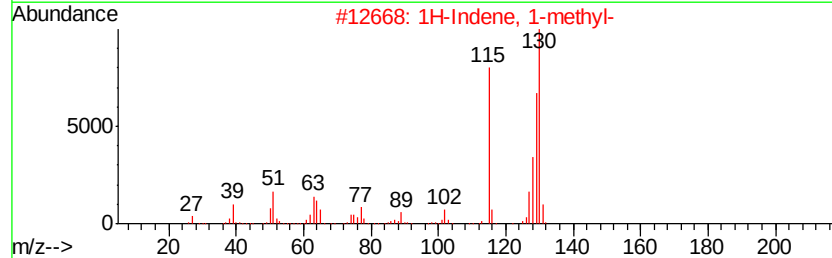
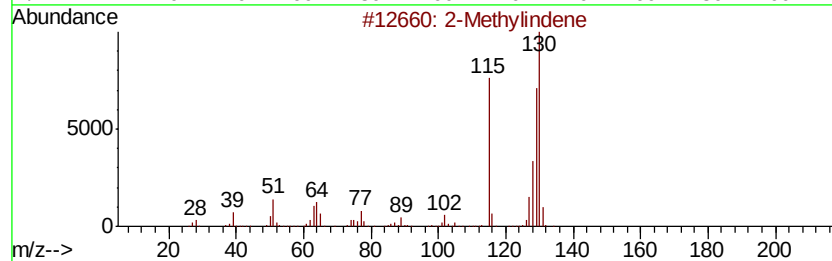
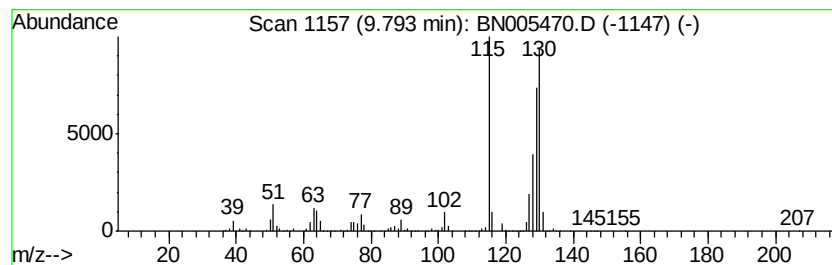
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Methylindene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	67.80 ng/ul	11441100	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		1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95
4		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
5		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
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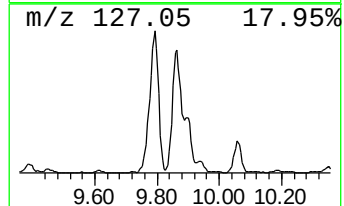
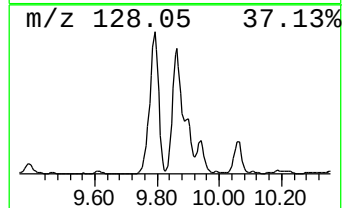
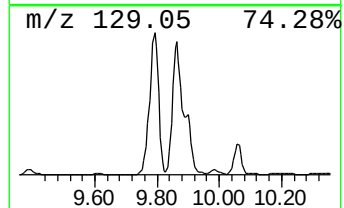
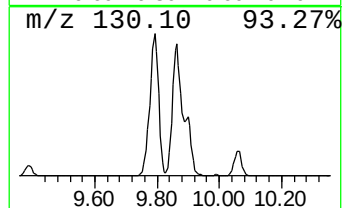
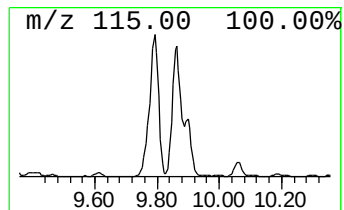
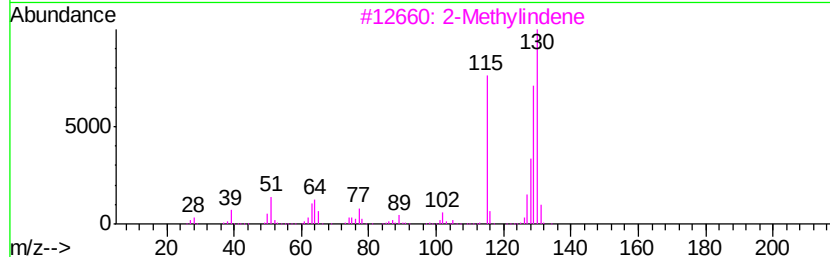
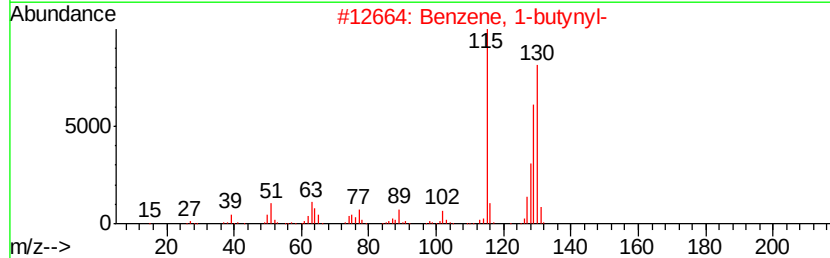
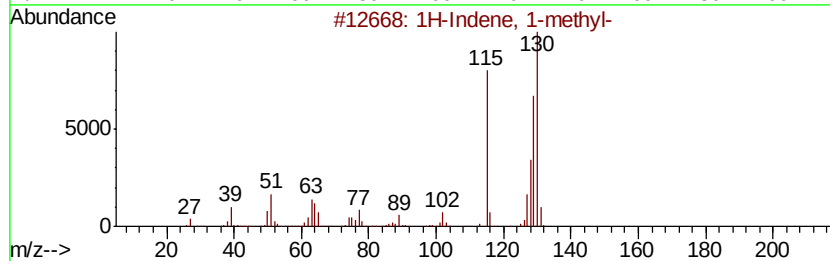
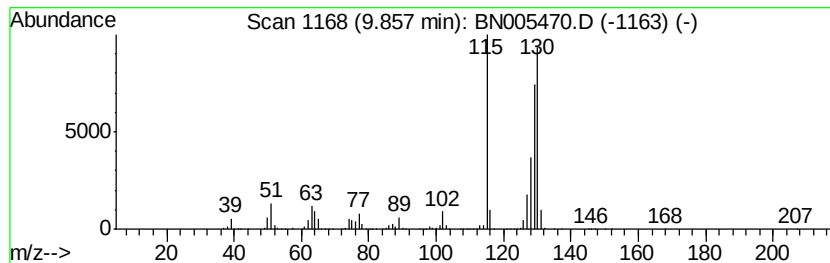
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1H-Indene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	47.97 ng/ul	8093710	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		Benzene, 1-butynyl-	130	C10H10	000622-76-4	95
3		2-Methylindene	130	C10H10	002177-47-1	94
4		Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93
5		2-Methylindene	130	C10H10	002177-47-1	93



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Data File : BN005470.D
Acq On : 04 May 2019 02:26
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Misc :
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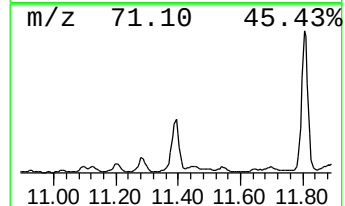
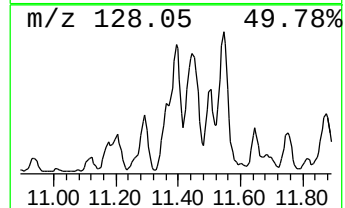
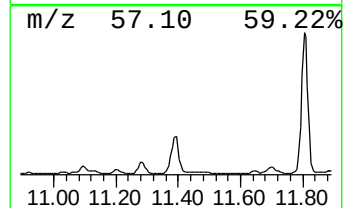
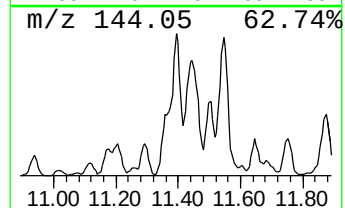
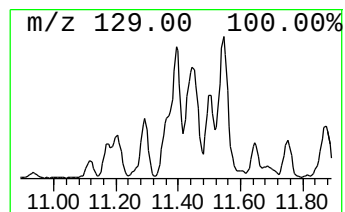
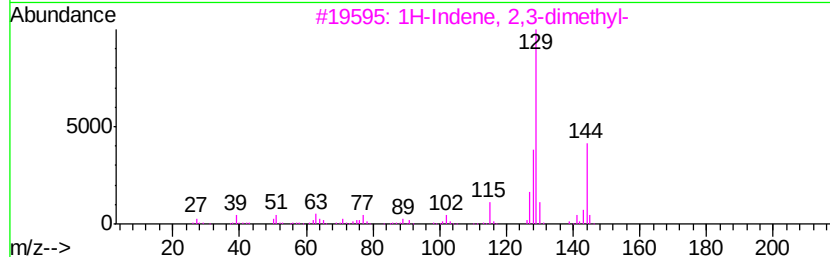
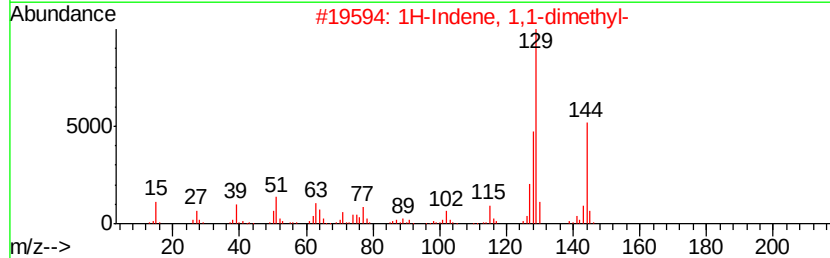
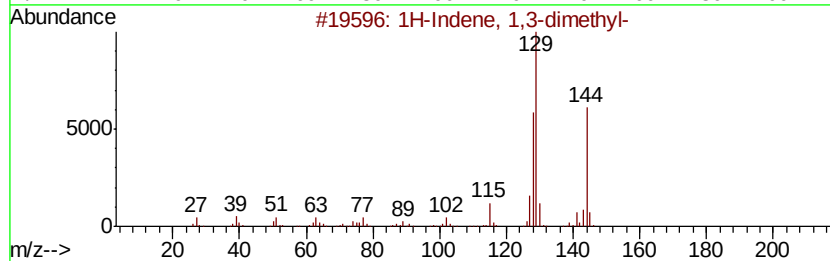
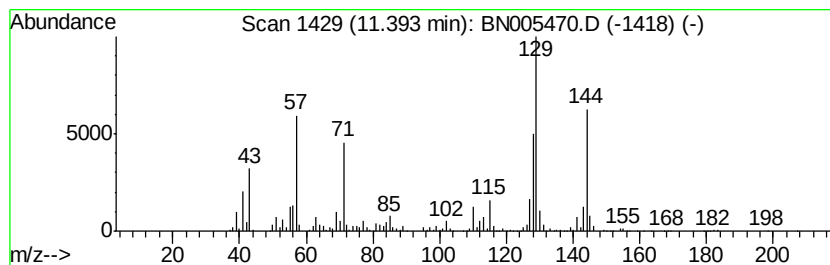
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1H-Indene, 1,3-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.39	21.68 ng/ul	3658940	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	95
2		1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	94
3		1H-Indene, 2,3-dimethyl-	144	C11H12	004773-82-4	90
4		1H-Cyclopropa[b]naphthalene, 1a,...	144	C11H12	006571-72-8	89
5		1H-Indene, 4,7-dimethyl-	144	C11H12	006974-97-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
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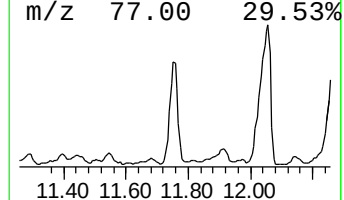
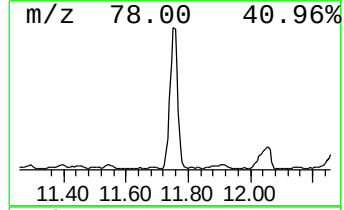
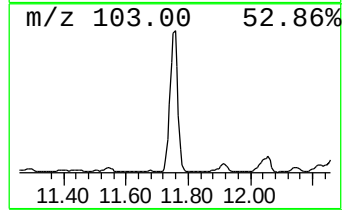
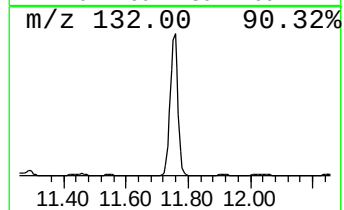
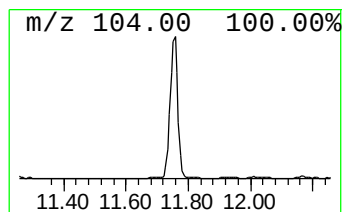
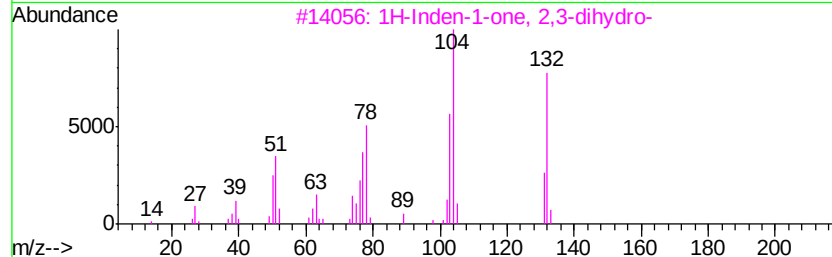
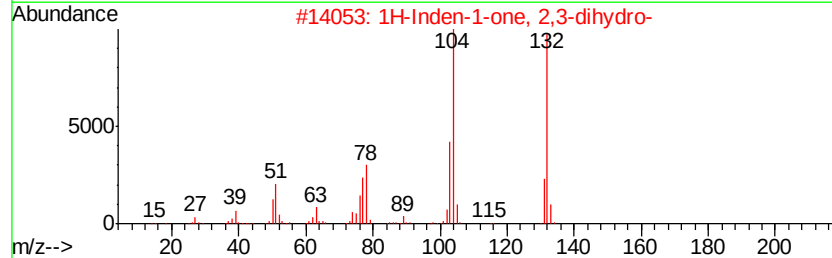
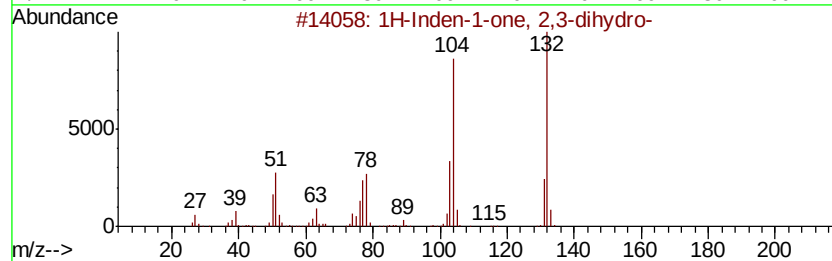
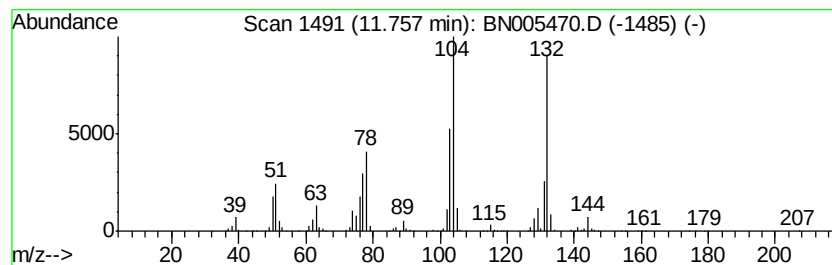
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	22.95 ng/ul	3872570	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	97
2		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	96
3		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	94
4		Styrene	104	C8H8	000100-42-5	91
5		Styrene	104	C8H8	000100-42-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
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Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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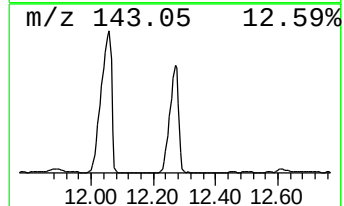
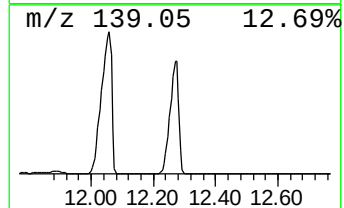
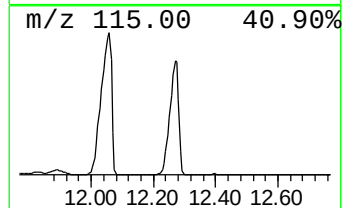
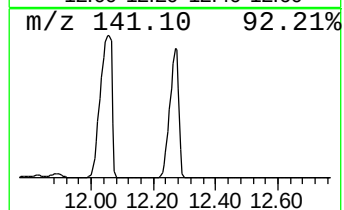
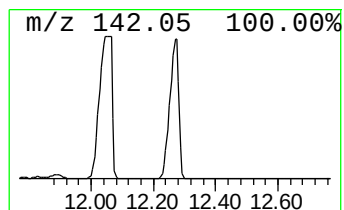
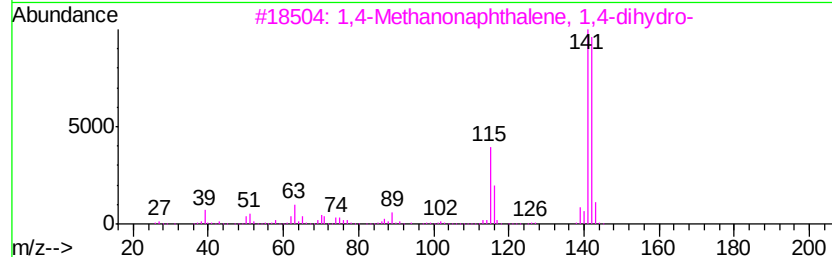
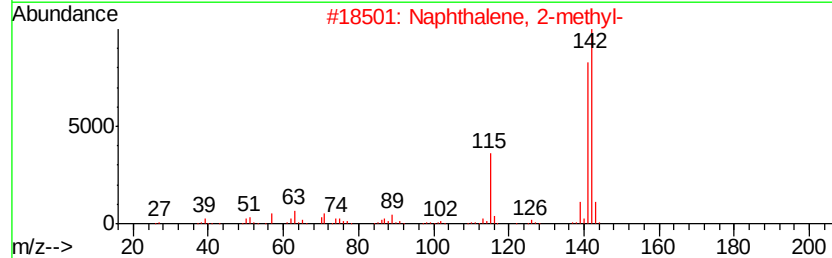
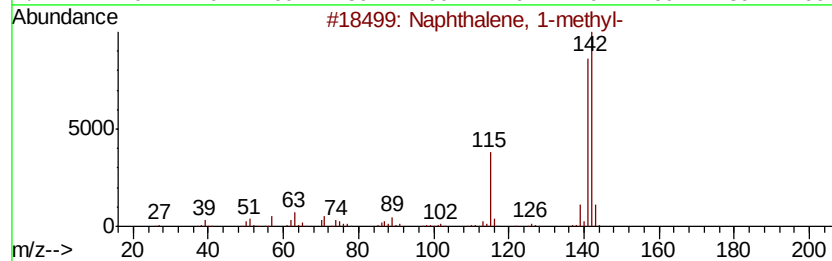
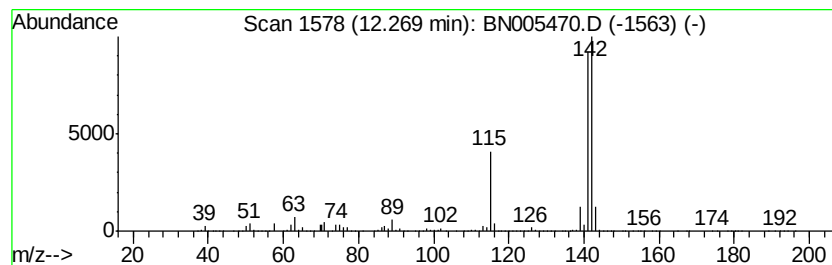
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	354.87 ng/ul	59880600	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		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	94
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

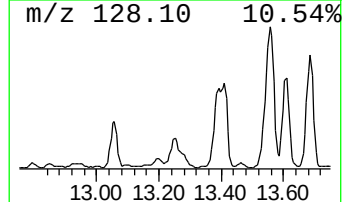
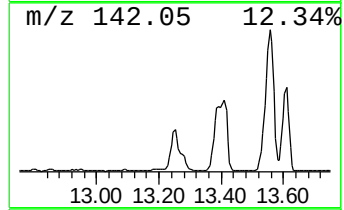
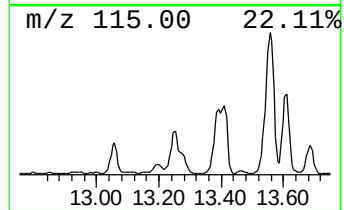
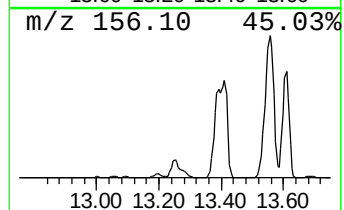
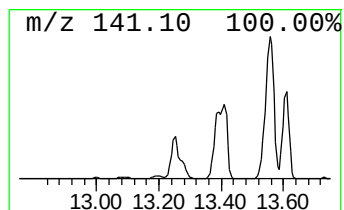
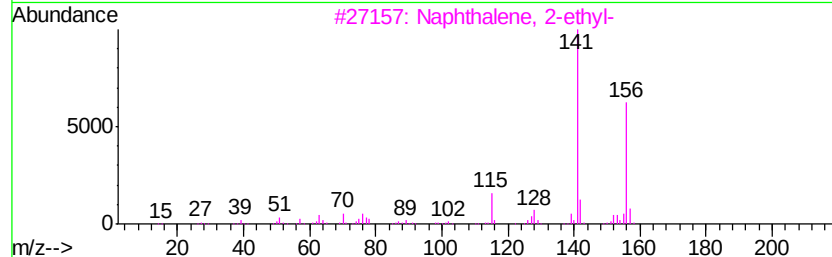
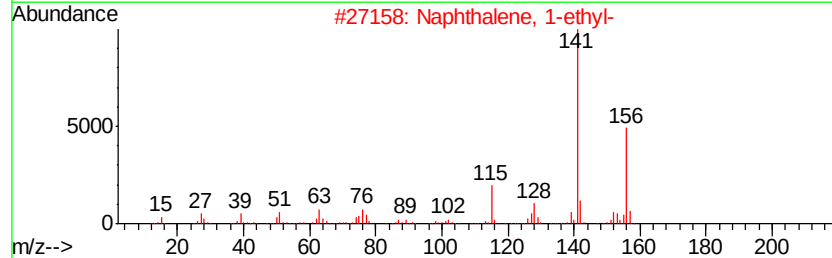
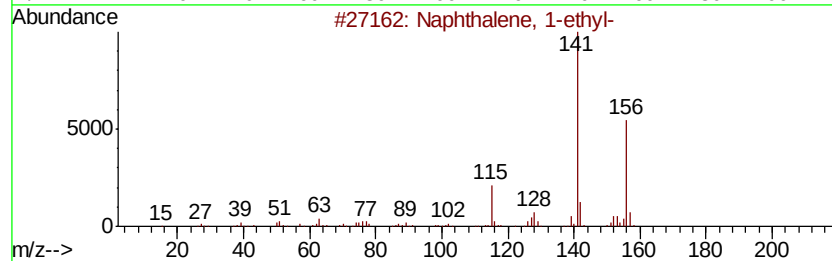
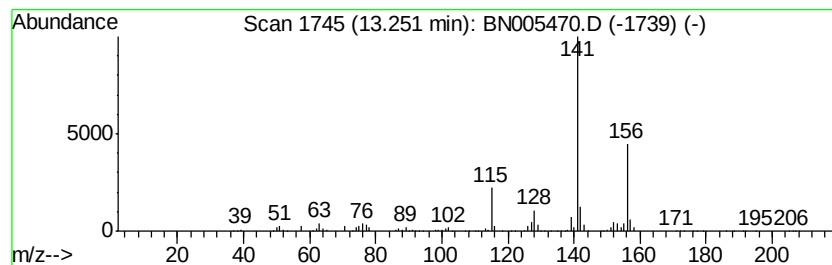
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Naphthalene, 1-ethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.25	13.89 ng/ul	8037520	Acenaphthene-d10	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 96
2		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 95
3		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 95
4		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 94
5		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 94



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
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Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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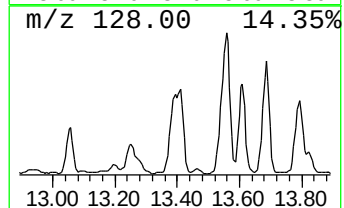
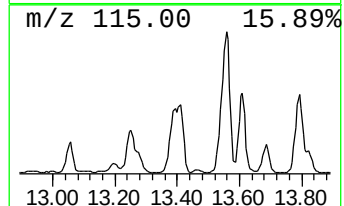
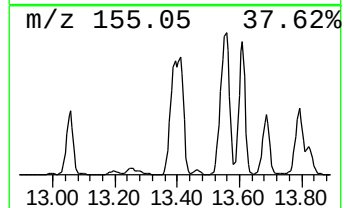
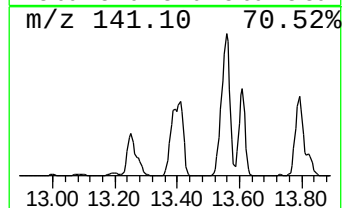
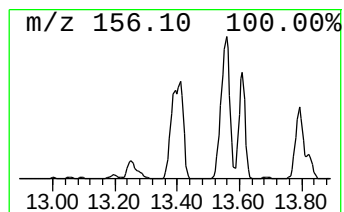
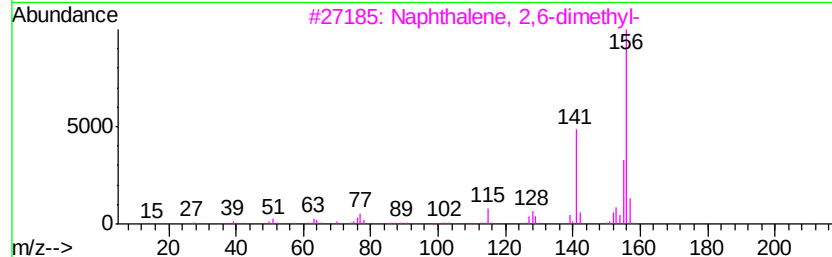
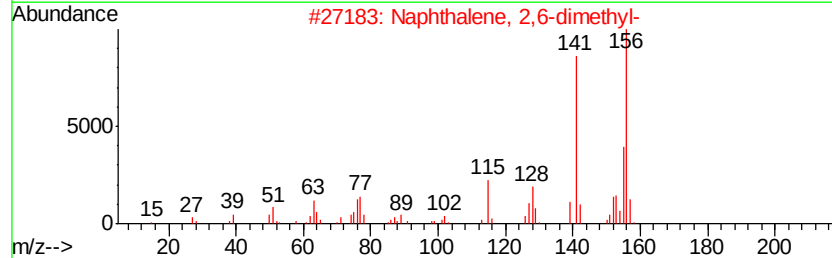
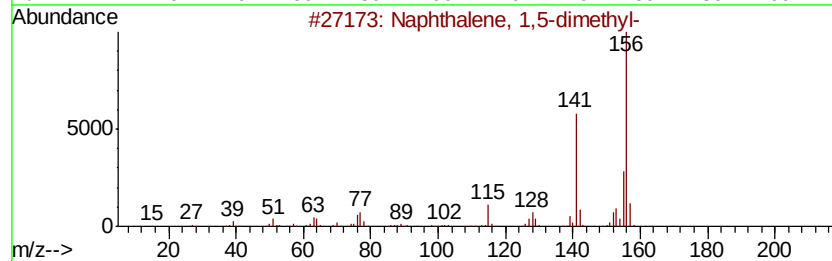
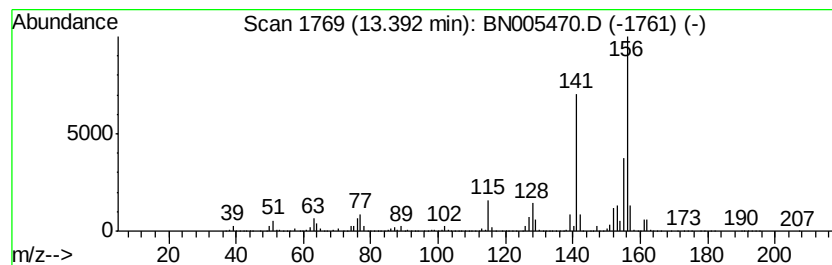
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Naphthalene, 1,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.39	26.65 ng/ul	15418200	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

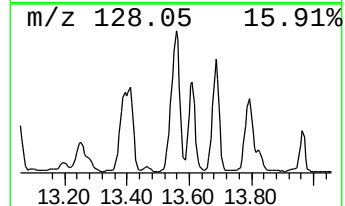
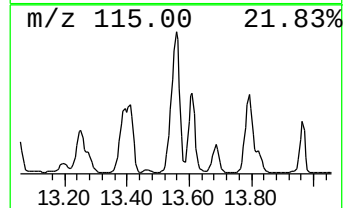
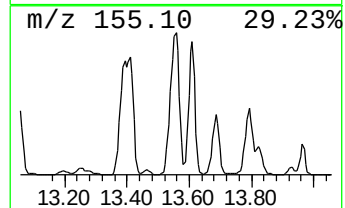
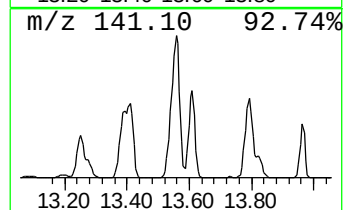
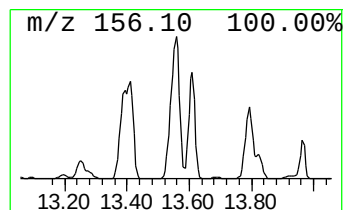
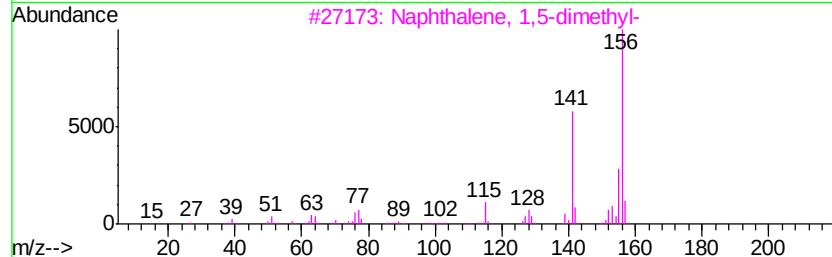
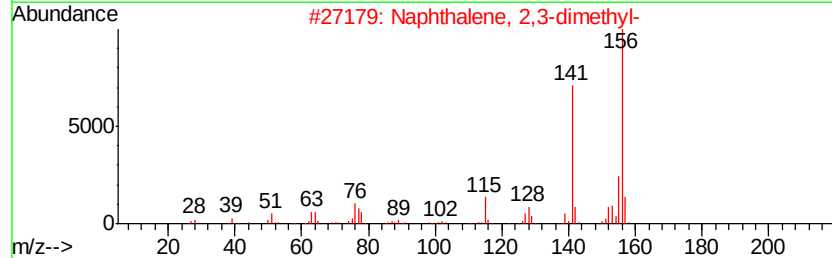
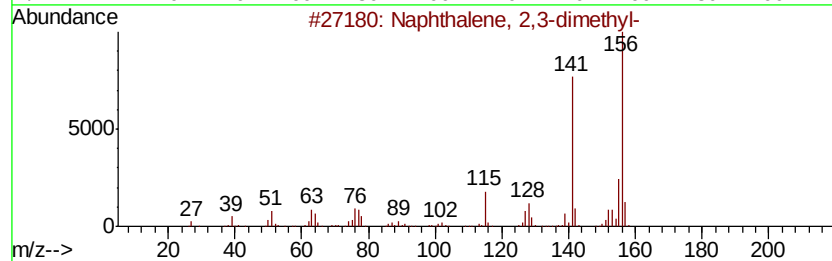
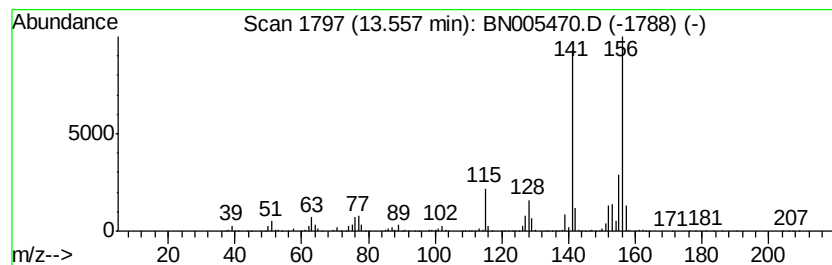
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.56	55.19 ng/ul	31930000	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, 2,3-dimethyl-	156 C12H12	000581-40-8 97
3		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
4		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97



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Sample : K2603-15DL 25X
Misc :
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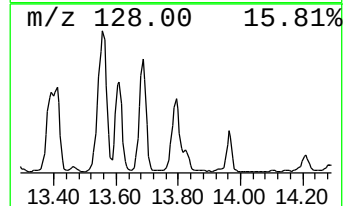
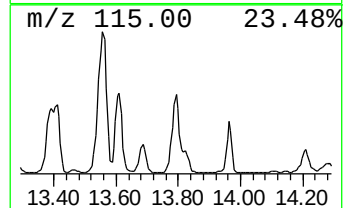
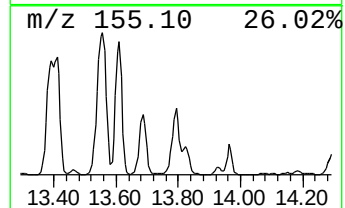
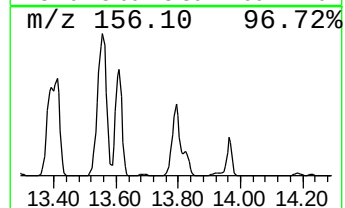
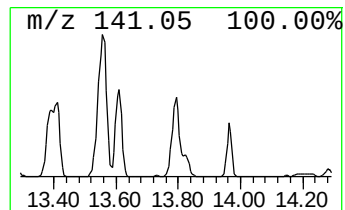
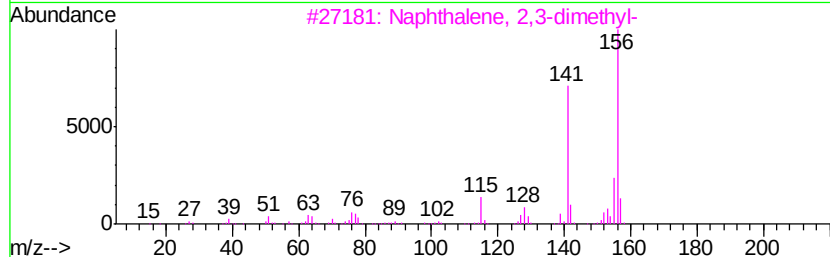
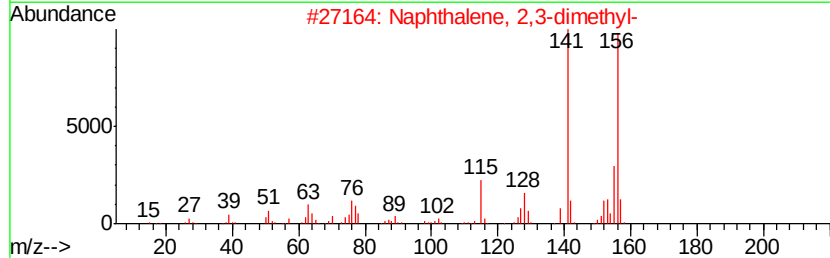
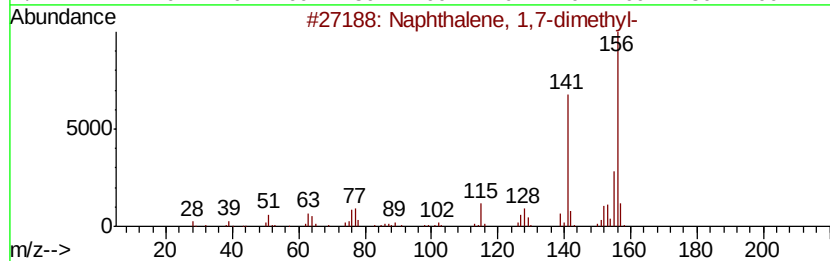
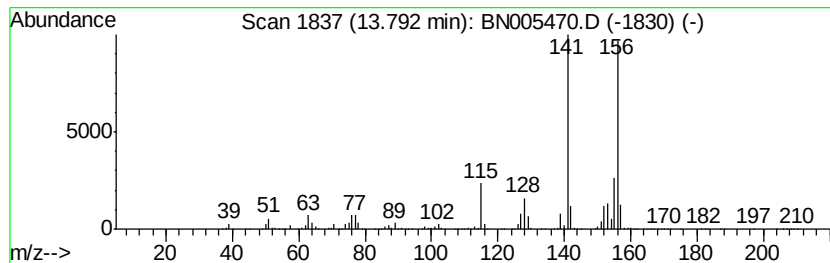
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Naphthalene, 1,7-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	25.36 ng/ul	14673700	Acenaphthene-d10	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
2		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
3		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
4		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
5		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97



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Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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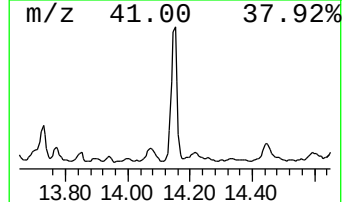
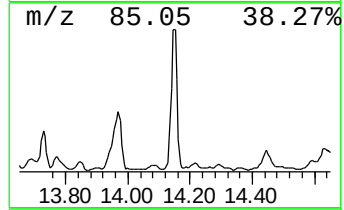
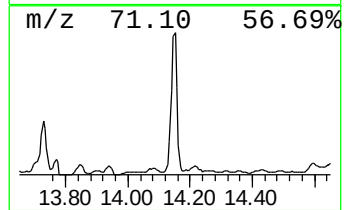
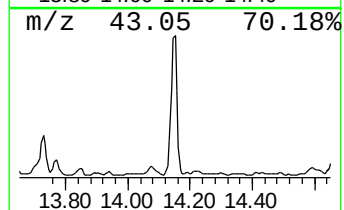
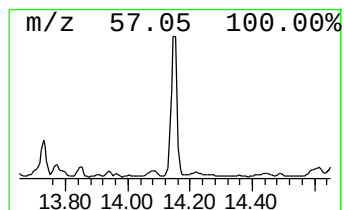
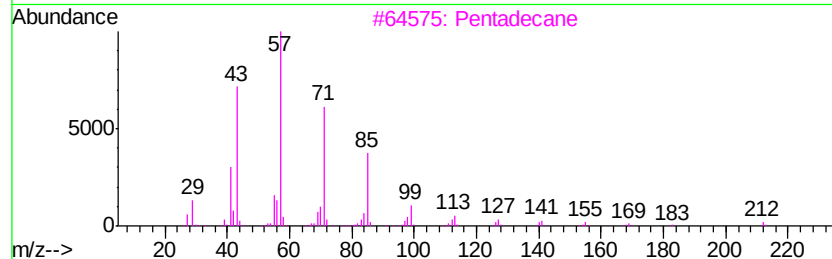
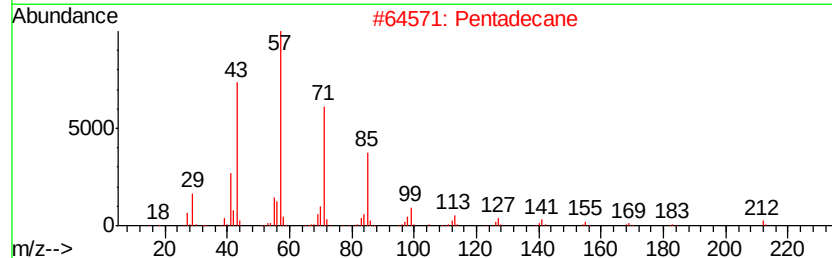
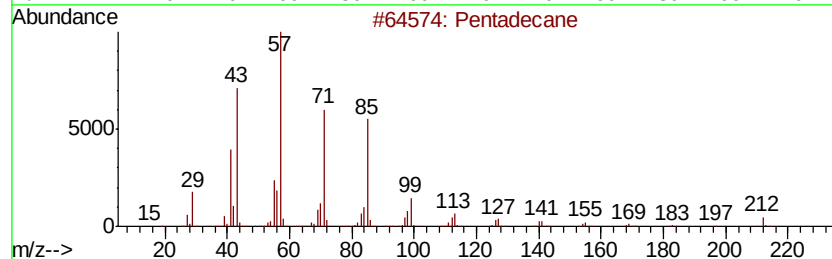
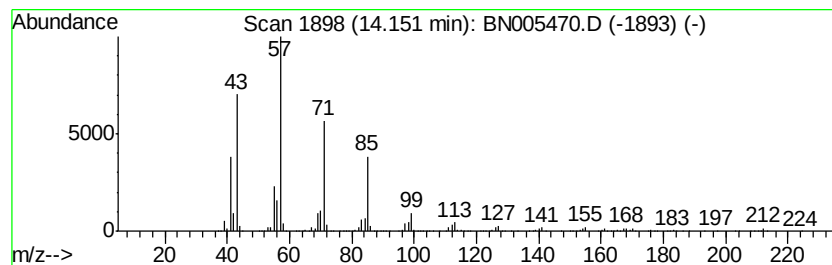
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	7.17 ng/ul	4149350	Acenaphthene-d10	14.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	97
2			Pentadecane	212	C15H32	000629-62-9	96
3			Pentadecane	212	C15H32	000629-62-9	95
4			Eicosane	282	C20H42	000112-95-8	91
5			Undecane, 2-methyl-	170	C12H26	007045-71-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
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Sample : K2603-15DL 25X
Misc :
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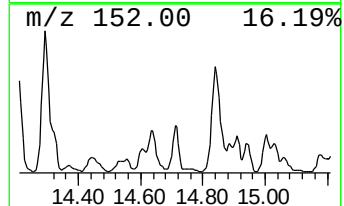
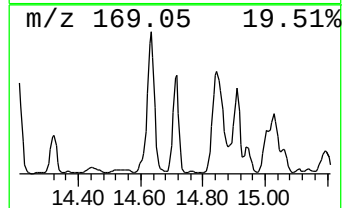
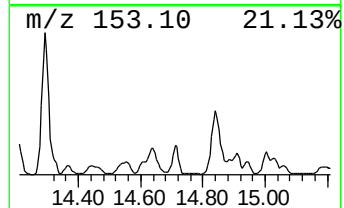
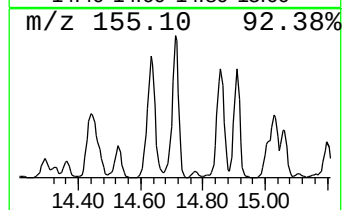
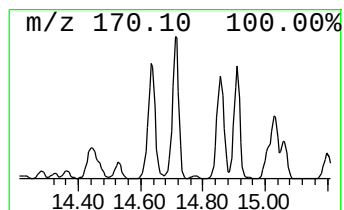
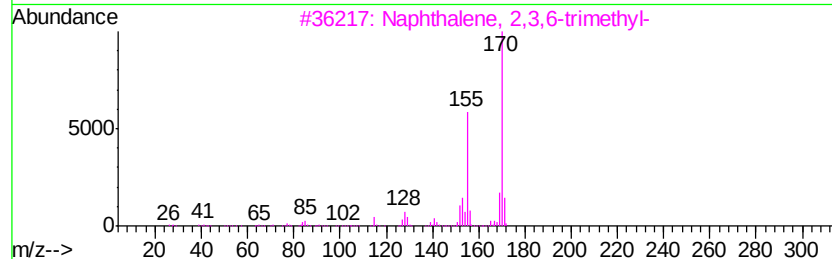
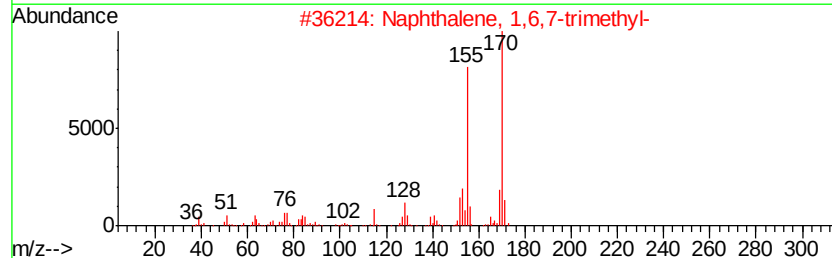
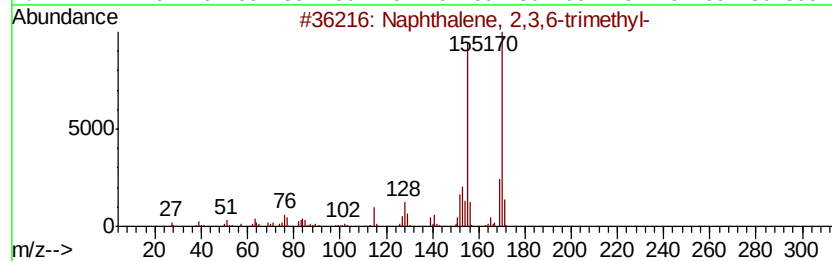
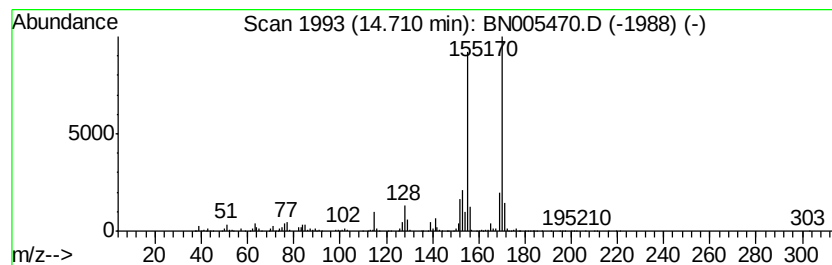
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Naphthalene, 2,3,6-trimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	11.45 ng/ul	6624340	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, 1,4,6-trimethyl-	170 C13H14	002131-42-2 96



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
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Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

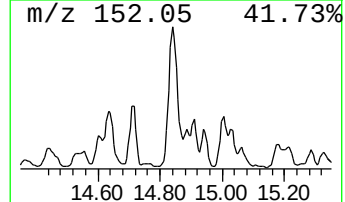
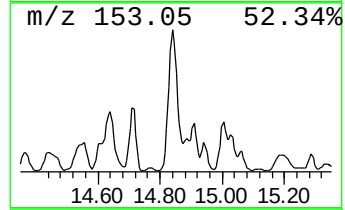
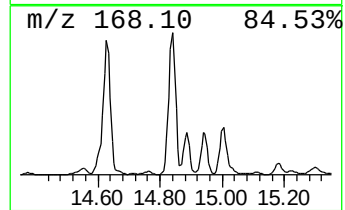
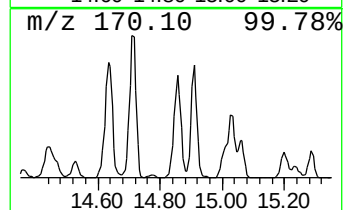
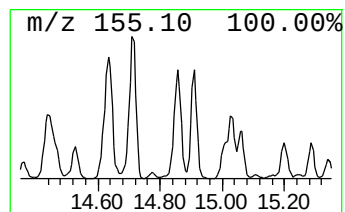
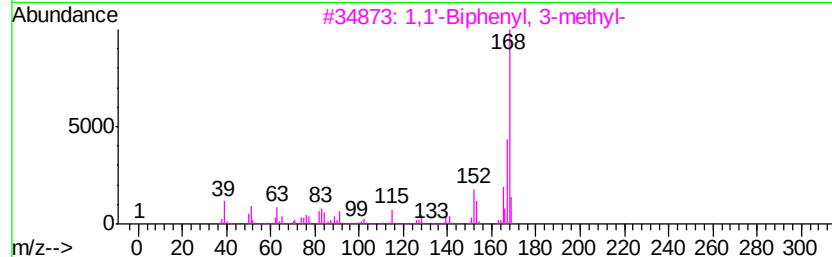
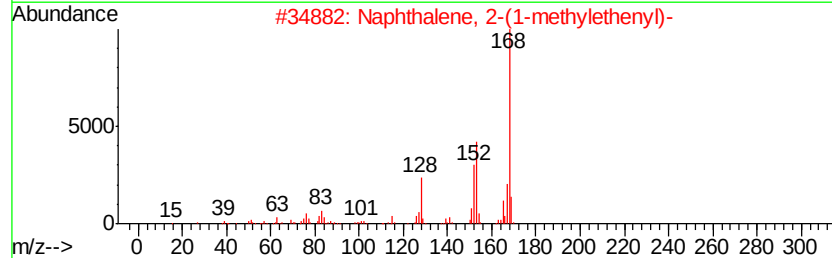
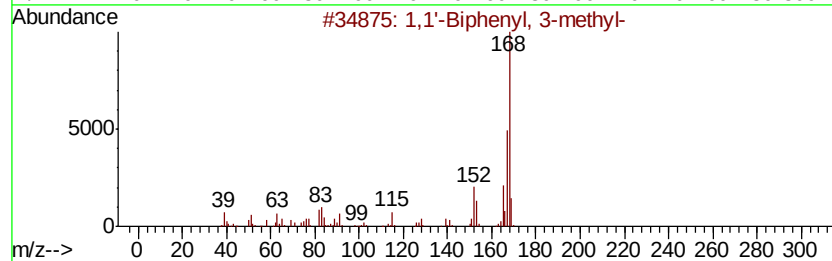
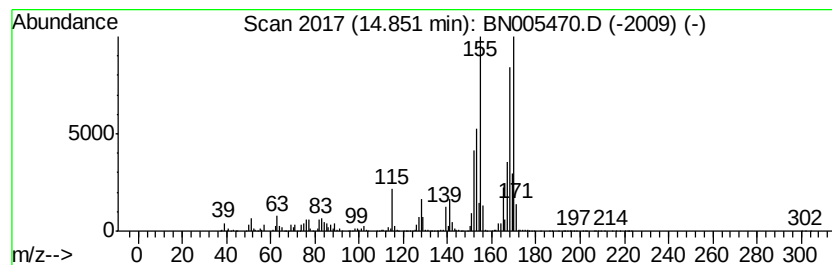
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 1,1'-Biphenyl, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.85	25.35 ng/ul	14666300	Acenaphthene-d10	14.23	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	50
2	Naphthalene, 2-(1-methylethenyl)-	168	C13H12	003710-23-4	47
3	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	43
4	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	38
5	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

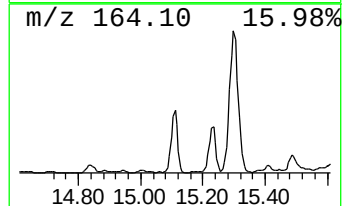
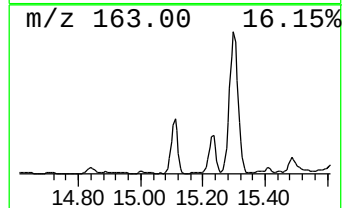
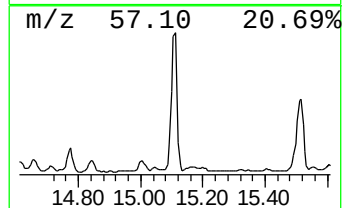
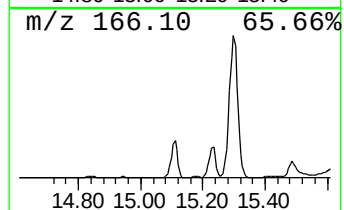
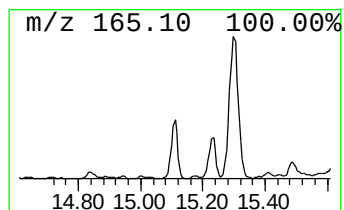
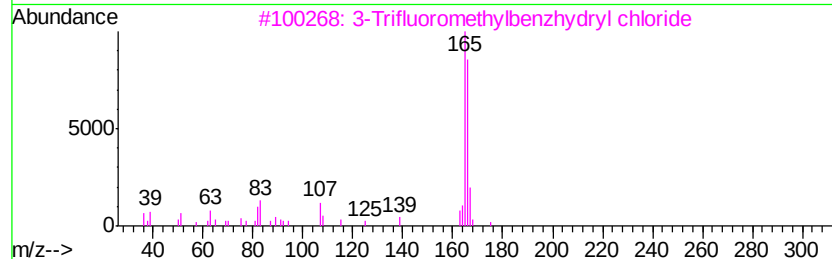
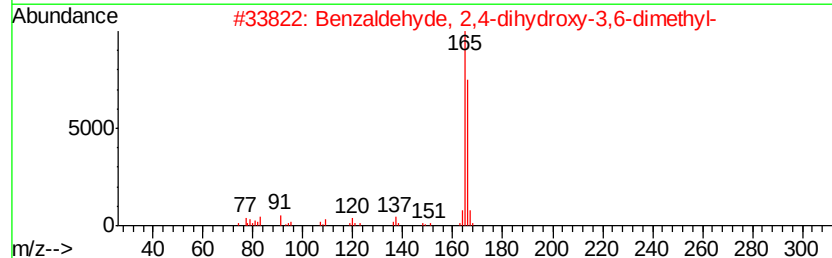
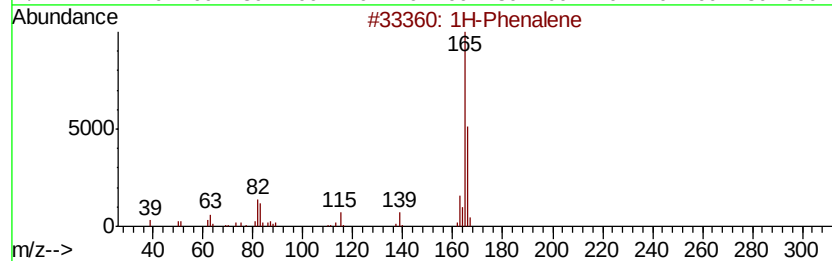
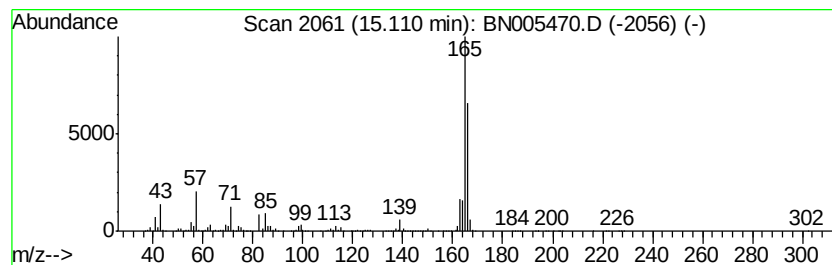
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 22 1H-Phenalene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	23.99 ng/ul	13879100	Acenaphthene-d10	14.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Phenalene	166 C13H10	000203-80-5 64
2		Benzaldehyde, 2,4-dihydroxy-3,6-...	166 C9H10O3	034883-14-2 64
3		3-Trifluoromethylbenzhvdrv1 chlo...	270 C14H10ClF3	067240-79-3 64
4		Fluorene-9-methanol	196 C14H12O	024324-17-2 64
5		9H-Fluorene, 9-(1-methylethyl)-	208 C16H16	003299-99-8 59



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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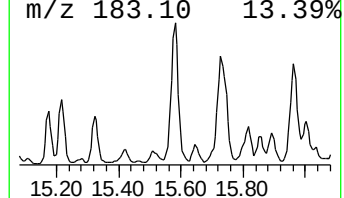
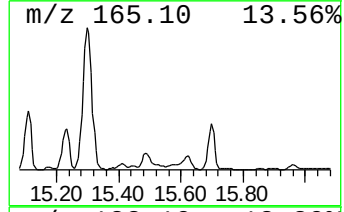
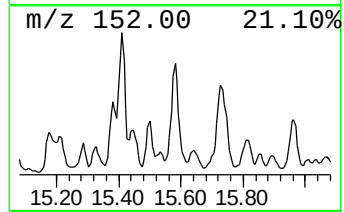
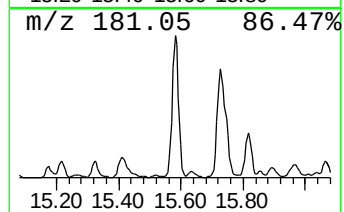
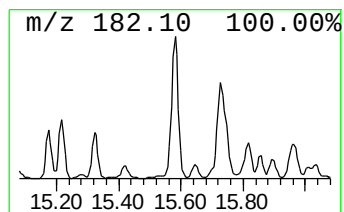
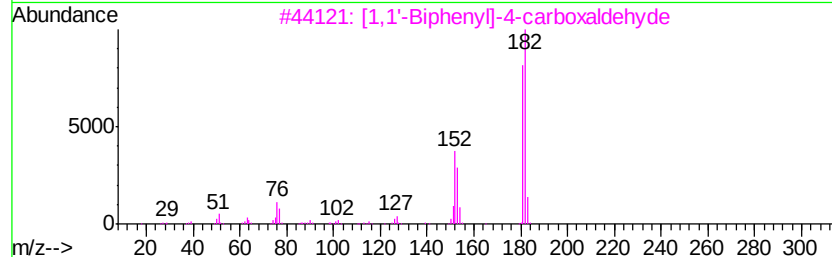
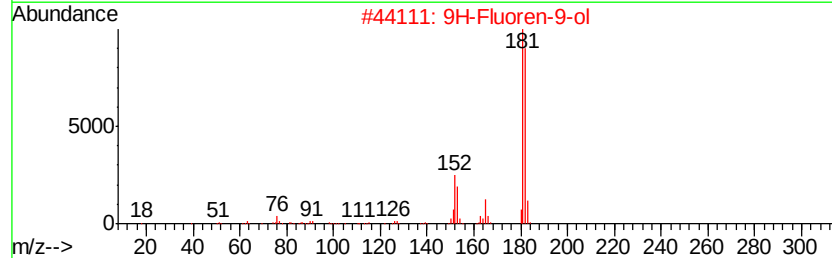
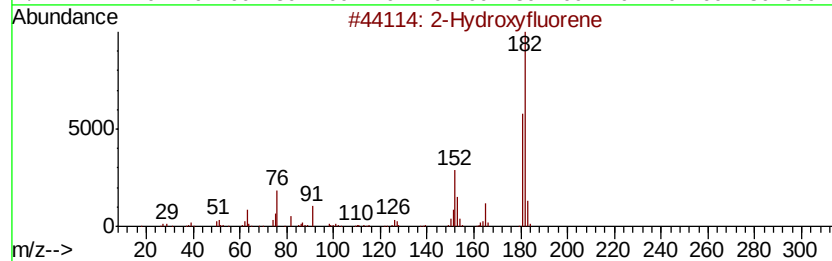
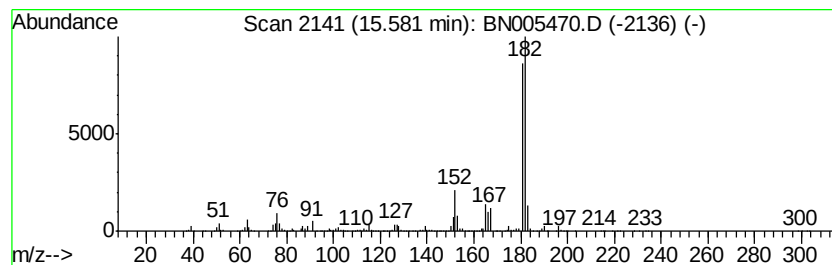
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 2-Hydroxyfluorene Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	10.32 ng/ul	5967760	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hydroxyfluorene	182	C13H10O	002443-58-5	76
2		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	72
3		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	68
4		Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	64
5		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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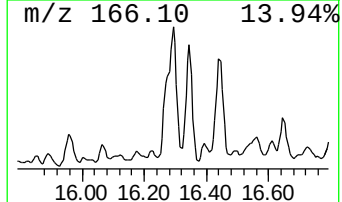
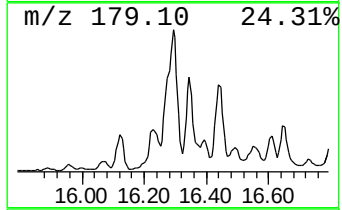
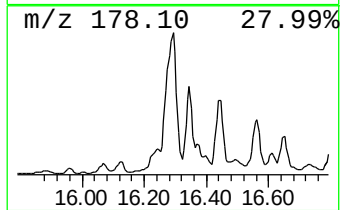
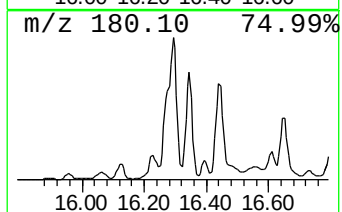
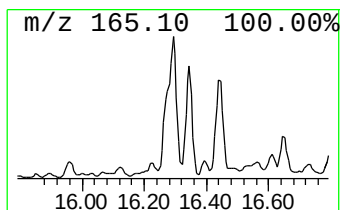
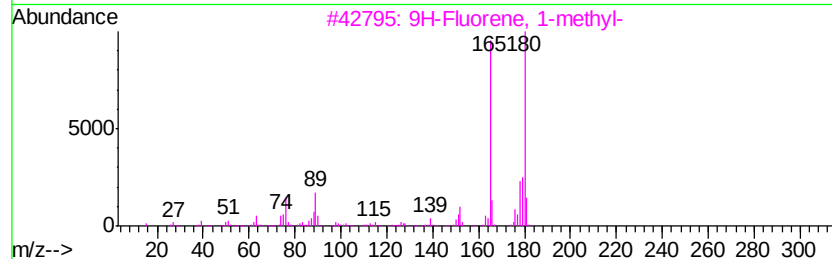
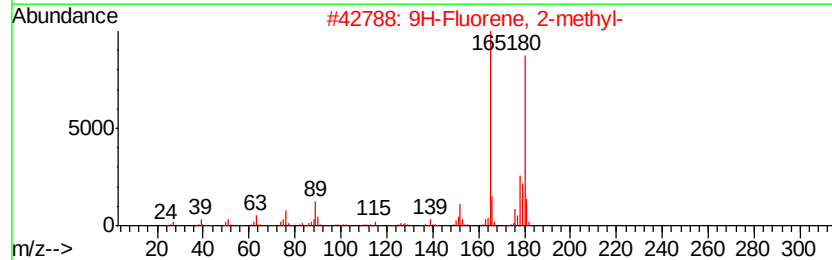
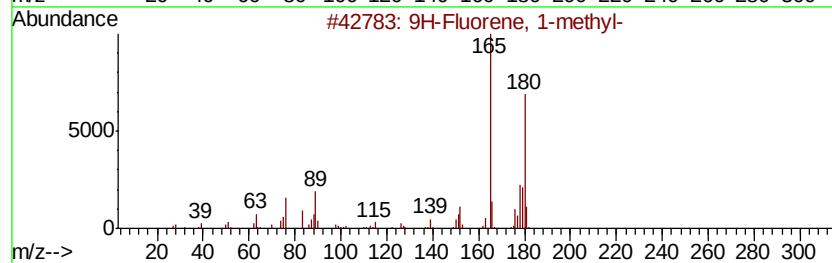
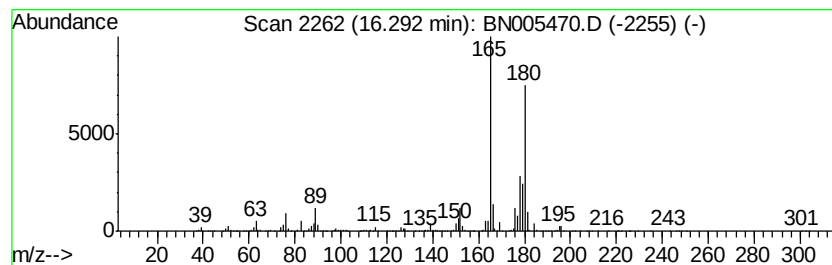
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 9H-Fluorene, 1-methyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.29	2.79 ng/ul	16250900	Phenanthrene-d10	16.99

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, 1-methyl-	180	C14H12	001730-37-6	93
4		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	93
5		4a,9a-Methano-9H-fluorene	180	C14H12	019540-84-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
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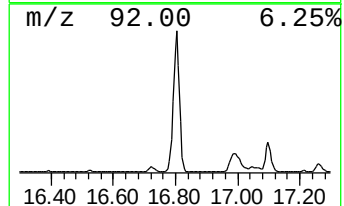
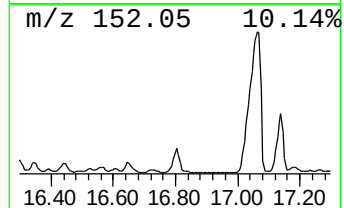
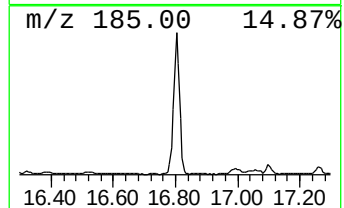
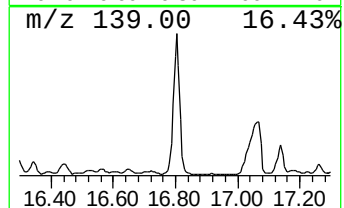
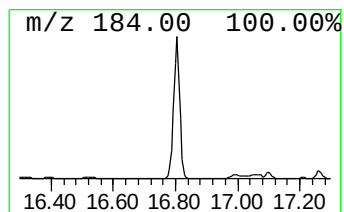
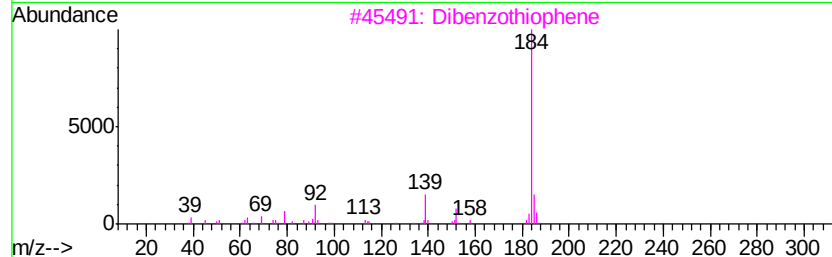
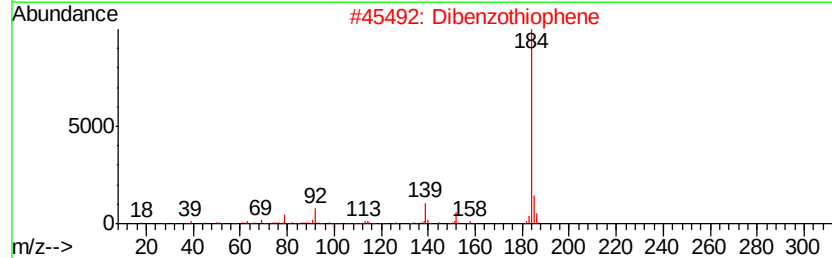
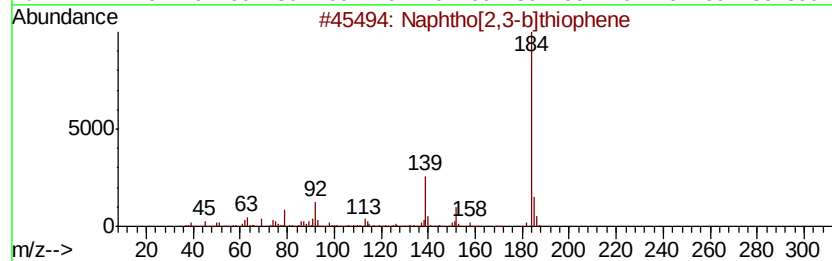
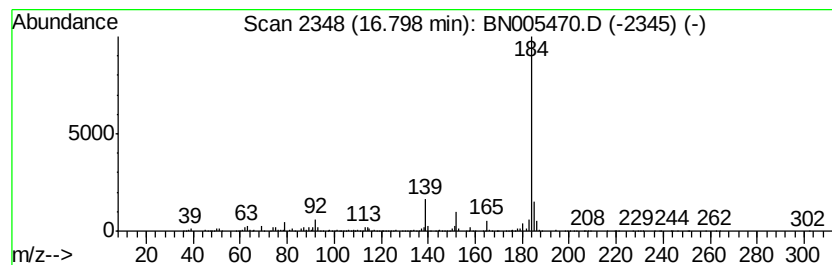
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 Naphtho[2,3-b]thiophene Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.80	2.77 ng/ul	16109700	Phenanthrene-d10	16.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
2		Dibenzothiophene	184	C12H8S	000132-65-0	95
3		Dibenzothiophene	184	C12H8S	000132-65-0	95
4		Dibenzothiophene	184	C12H8S	000132-65-0	94
5		Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
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Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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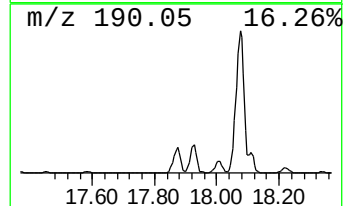
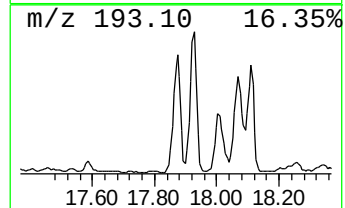
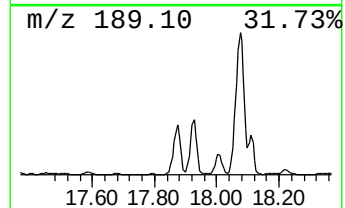
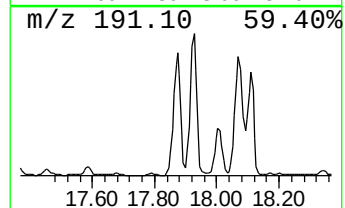
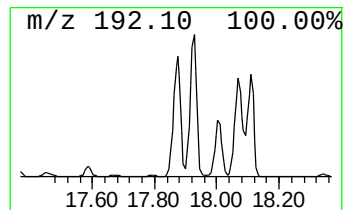
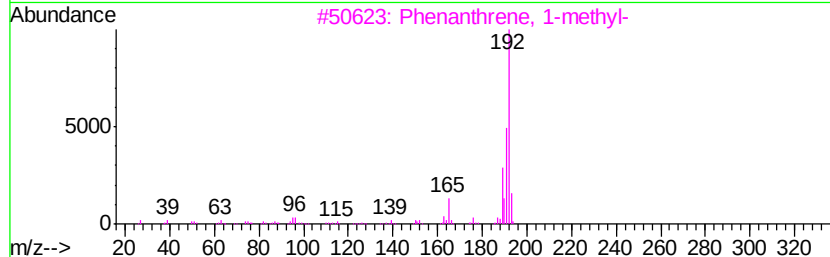
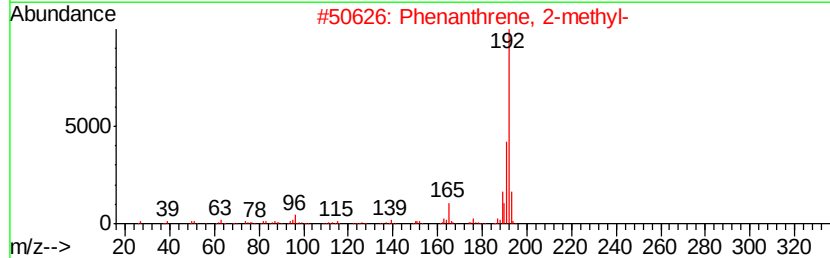
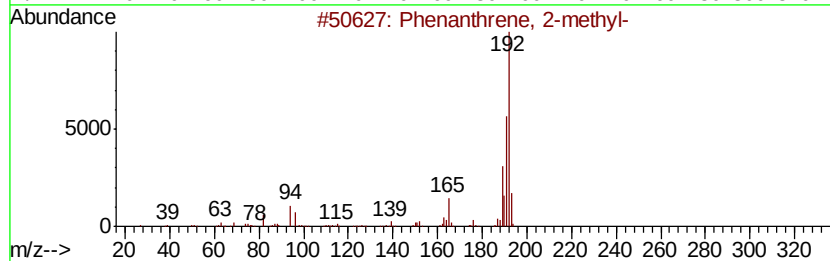
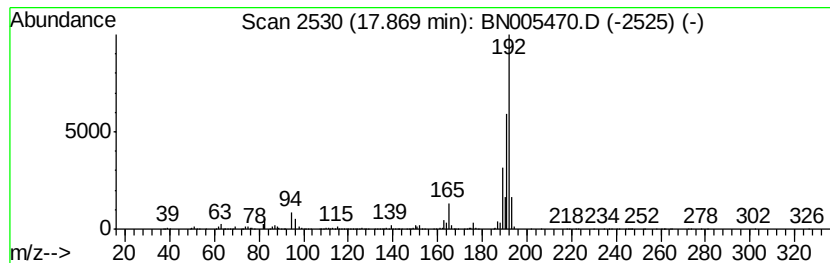
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Phenanthrene, 2-methyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	5.45 ng/ul	31690800	Phenanthrene-d10	16.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
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Misc :
ALS Vial : 22 Sample Multiplier: 1

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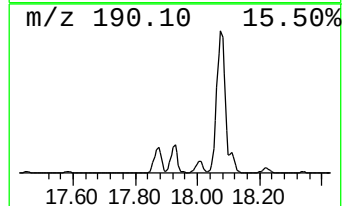
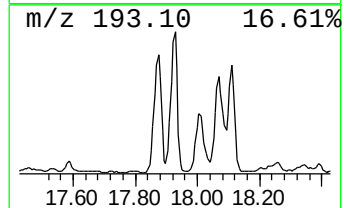
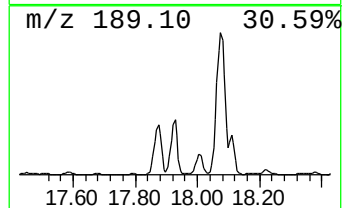
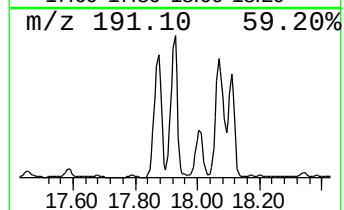
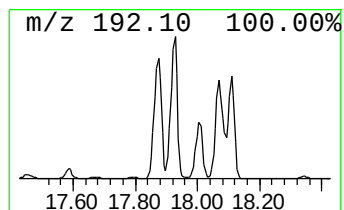
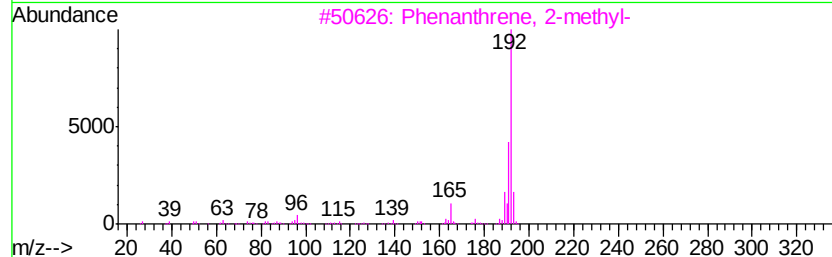
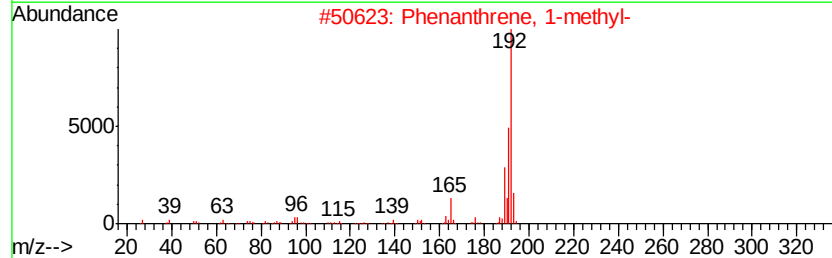
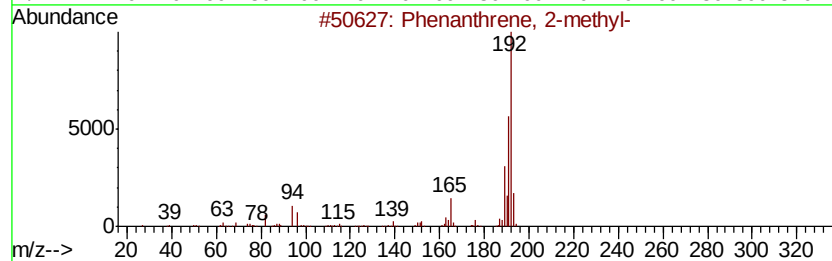
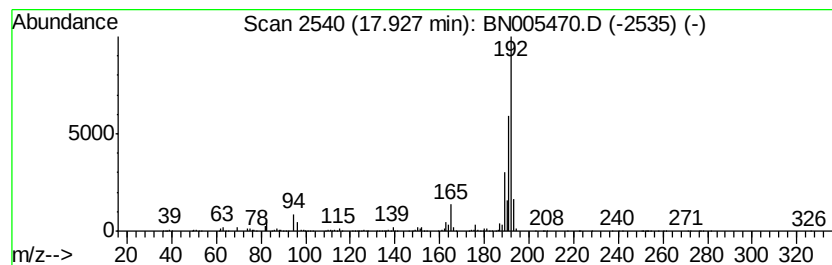
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 Phenanthrene, 1-methyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.93	5.74 ng/ul	33393500	Phenanthrene-d10	16.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
4			1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	96
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GAJ67DL

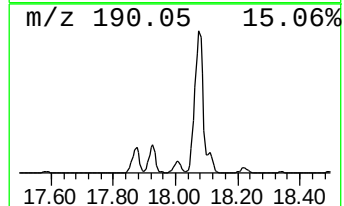
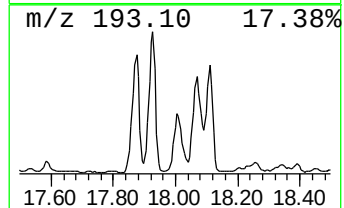
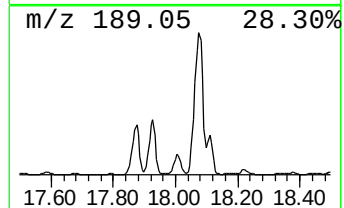
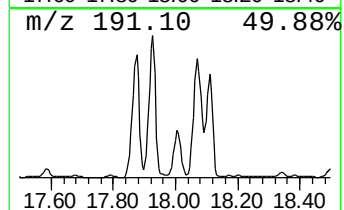
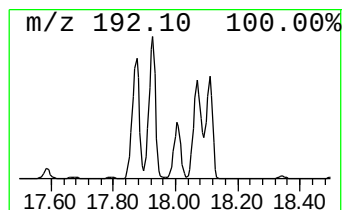
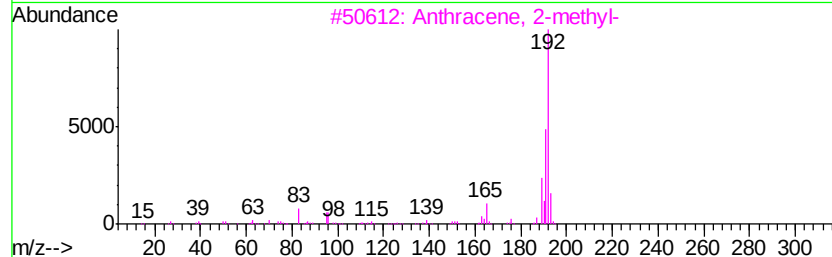
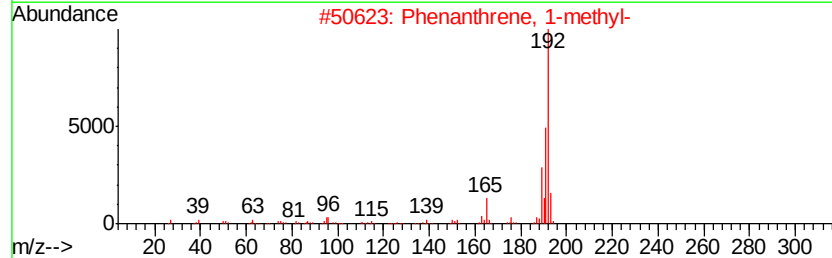
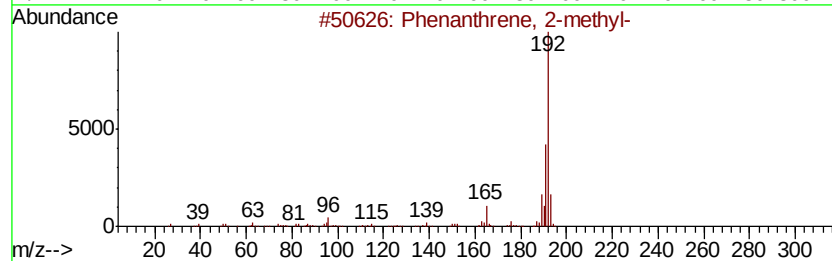
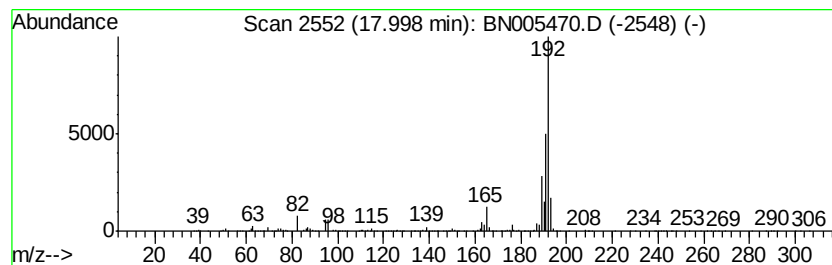
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Anthracene, 2-methyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	2.22 ng/ul	12931900	Phenanthrene-d10	16.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5		1H-Cyclopropa[l]phenanthrene, 1a,...	192	C15H12	000949-41-7	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

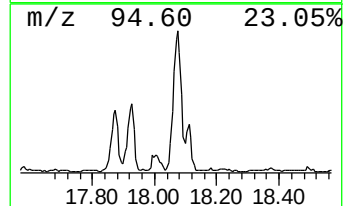
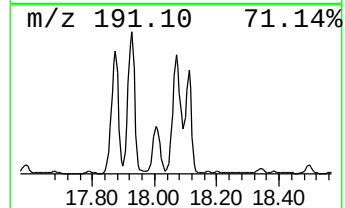
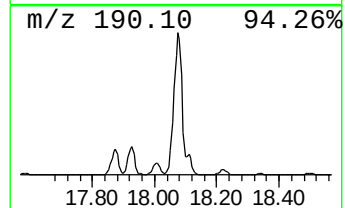
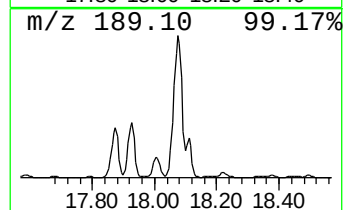
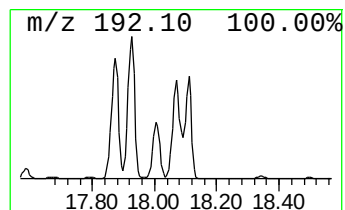
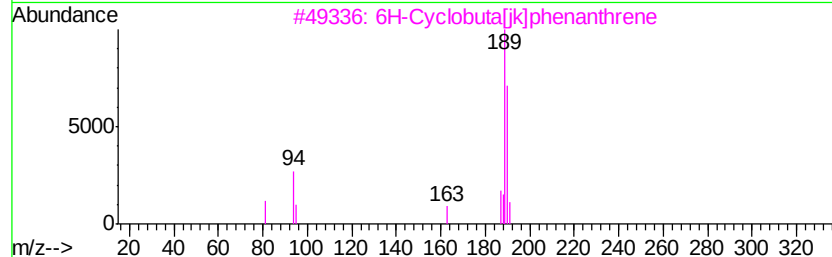
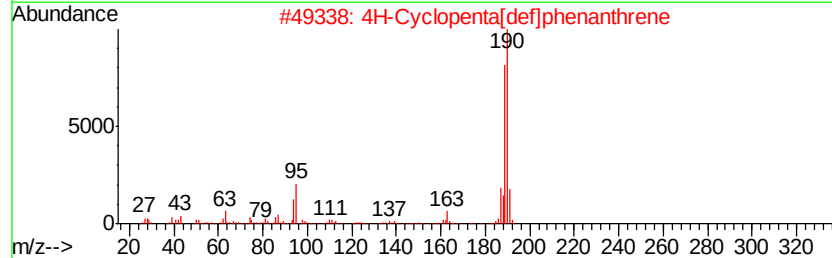
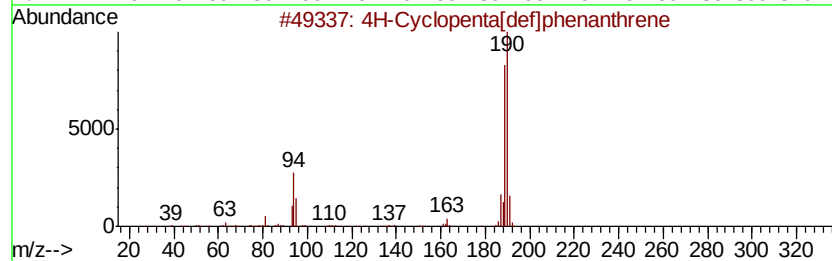
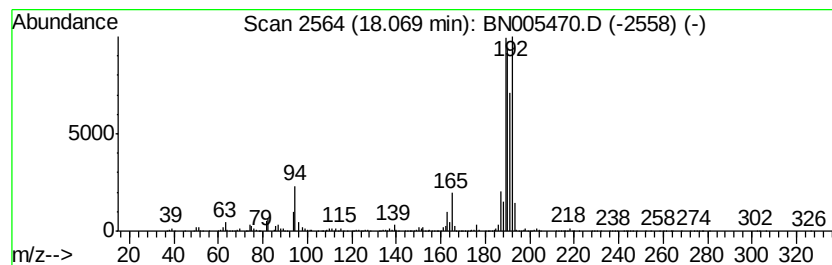
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 4H-Cyclopenta[def]phenanthrene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.07	9.75 ng/ul	56692200	Phenanthrene-d10	16.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
3	6H-Cyclobuta[f]phenanthrene	190	C15H10	083469-43-6	49
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN050319\
Data File : BN005470.D
Acq On : 04 May 2019 02:26
Operator : JU/SJ
Sample : K2603-15DL 25X
Misc :
ALS Vial : 22 Sample Multiplier: 1

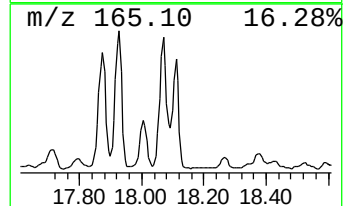
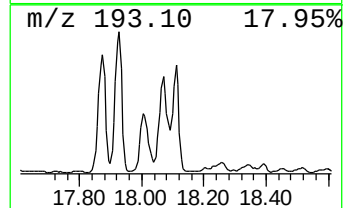
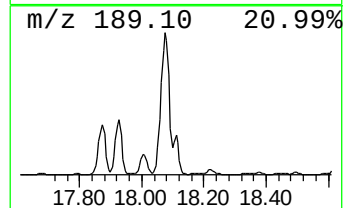
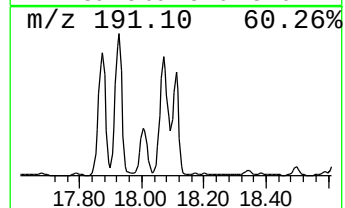
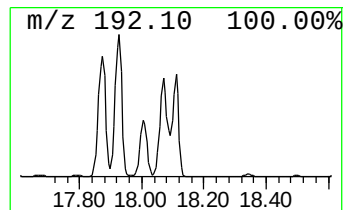
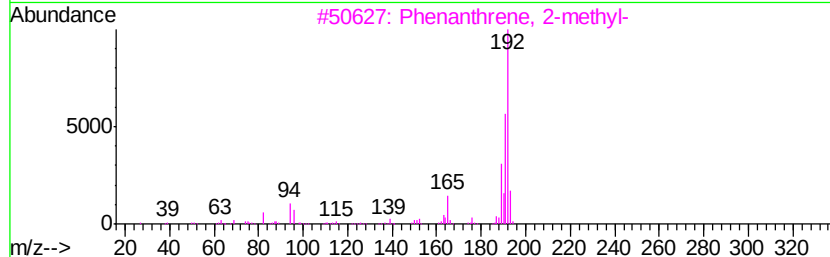
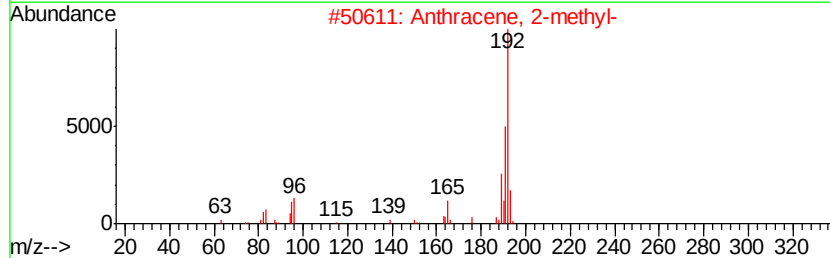
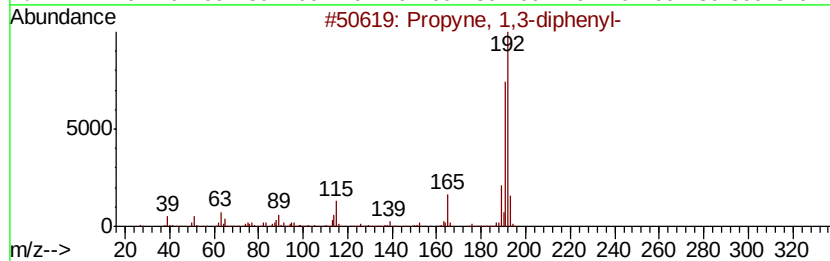
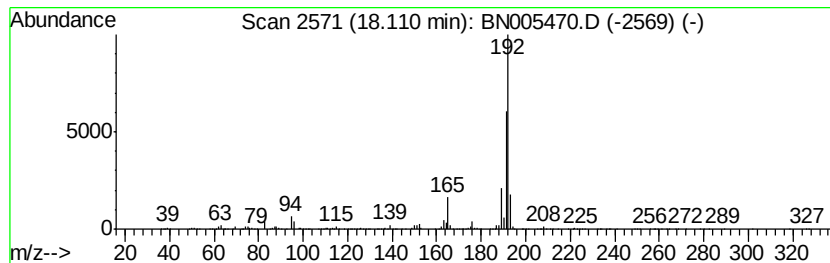
Instrument :
BNA_N
ClientSampleId :
GAJ67DL

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 31 Propyne, 1,3-diphenyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.11	2.93 ng/ul	17022200	Phenanthrene-d10	16.99		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propyne, 1,3-diphenyl-		192	C15H12	004980-70-5	91
2	Anthracene, 2-methyl-		192	C15H12	000613-12-7	91
3	Phenanthrene, 2-methyl-		192	C15H12	002531-84-2	90
4	1H-Indene, 2-phenyl-		192	C15H12	004505-48-0	90
5	1H-Indene, 3-phenyl-		192	C15H12	001961-97-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN050319\
 Data File : BN005470.D
 Acq On : 04 May 2019 02:26
 Operator : JU/SJ
 Sample : K2603-15DL 25X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GAJ67DL

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, 1,3-dime...	5.29	44.7	ng/ul	7899810	1	7.59	3533110	20.0
Bicyclo[4.2.0]oct...	7.25	86.9	ng/ul	15347500	1	7.59	3533110	20.0
Benzene, 1-etheny...	7.33	20.0	ng/ul	3533110	1	7.59	3533110	20.0
Indene	8.12	163.2	ng/ul	28832700	1	7.59	3533110	20.0
Benzene, 1-methyl...	8.85	19.2	ng/ul	3397340	1	7.59	3533110	20.0
2-Methylindene	9.79	67.8	ng/ul	11441100	2	10.36	3374770	20.0
1H-Indene, 1-methyl-	9.86	48.0	ng/ul	8093710	2	10.36	3374770	20.0
1H-Indene, 1,3-di...	11.39	21.7	ng/ul	3658940	2	10.36	3374770	20.0
1H-Inden-1-one, 2...	11.76	22.9	ng/ul	3872570	2	10.36	3374770	20.0
Naphthalene, 1-me...	12.27	354.9	ng/ul	59880600	2	10.36	3374770	20.0
Naphthalene, 1-et...	13.25	13.9	ng/ul	8037520	3	14.23	11570700	20.0
Naphthalene, 1,5-...	13.39	26.6	ng/ul	15418200	3	14.23	11570700	20.0
Naphthalene, 2,3-...	13.56	55.2	ng/ul	31930000	3	14.23	11570700	20.0
Naphthalene, 1,7-...	13.79	25.4	ng/ul	14673700	3	14.23	11570700	20.0
(DEL) Alkane: Str...	14.15	7.2	ng/ul	4149350	3	14.23	11570700	20.0
Naphthalene, 2,3,...	14.71	11.4	ng/ul	6624340	3	14.23	11570700	20.0
1,1'-Biphenyl, 3-...	14.85	25.4	ng/ul	14666300	3	14.23	11570700	20.0
1H-Phenylene	15.11	24.0	ng/ul	13879100	3	14.23	11570700	20.0
2-Hydroxyfluorene	15.58	10.3	ng/ul	5967760	3	14.23	11570700	20.0
9H-Fluorene, 1-me...	16.29	2.8	ng/ul	16250900	4	16.99	116329000	20.0
Naphtho[2,3-b]thi...	16.80	2.8	ng/ul	16109700	4	16.99	116329000	20.0
Phenanthrene, 2-m...	17.87	5.5	ng/ul	31690800	4	16.99	116329000	20.0
Phenanthrene, 1-m...	17.93	5.7	ng/ul	33393500	4	16.99	116329000	20.0
Anthracene, 2-met...	18.00	2.2	ng/ul	12931900	4	16.99	116329000	20.0
4H-Cyclopenta[def...	18.07	9.8	ng/ul	56692200	4	16.99	116329000	20.0
Propyne, 1,3-diph...	18.11	2.9	ng/ul	17022200	4	16.99	116329000	20.0