

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122718\
 Data File : BN004247.D
 Acq On : 27 Dec 2018 18:50
 Operator : JU/SJ
 Sample : J6441-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BE802

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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-BN121218MA.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	3.058	10	12	18	rVB	8832	11216	1.16%	0.088%
2	4.940	324	332	337	rBB2	6832	10835	1.12%	0.085%
3	5.328	393	398	404	rBB	35892	51835	5.38%	0.405%
4	6.299	558	563	568	rBB	18763	25002	2.59%	0.195%
5	6.787	641	646	652	rVB	73420	104622	10.85%	0.817%
6	6.893	658	664	670	rBV	264638	397292	41.22%	3.103%
7	6.958	670	675	678	rVV	128691	207261	21.50%	1.619%
8	6.987	678	680	683	rVV	85088	115810	12.02%	0.905%
9	7.028	683	687	694	rVB	194672	299289	31.05%	2.338%
10	7.158	703	709	712	rBV	76693	118116	12.25%	0.923%
11	7.199	712	716	722	rVB	94690	137088	14.22%	1.071%
12	7.352	736	742	748	rBV	84687	127901	13.27%	0.999%
13	7.458	754	760	767	rVB	370889	553222	57.40%	4.321%
14	7.822	816	822	832	rBB	134934	222947	23.13%	1.741%
15	7.922	833	839	844	rBV	86474	129393	13.42%	1.011%
16	8.293	897	902	907	rBV	72233	103552	10.74%	0.809%
17	8.352	907	912	917	rVB	31162	43894	4.55%	0.343%
18	8.440	921	927	933	rBV	94107	141745	14.71%	1.107%
19	8.528	936	942	951	rVB	106478	170836	17.72%	1.334%
20	8.605	951	955	960	rBV	7010	9773	1.01%	0.076%
21	8.752	975	980	985	rBV	22925	31439	3.26%	0.246%
22	8.805	985	989	995	rVB	25379	35252	3.66%	0.275%
23	8.905	1000	1006	1012	rBB	103296	148457	15.40%	1.160%
24	8.993	1013	1021	1029	rBB	95769	159941	16.59%	1.249%
25	9.705	1137	1142	1149	rBB	56132	88631	9.20%	0.692%
26	10.240	1225	1233	1244	rBB2	104650	184822	19.18%	1.444%
27	10.616	1290	1297	1303	rBV	225244	365659	37.94%	2.856%
28	10.663	1303	1305	1313	rVB2	7272	9842	1.02%	0.077%
29	10.763	1316	1322	1339	rBB	120110	203994	21.16%	1.593%
30	10.987	1355	1360	1371	rBB	94631	139095	14.43%	1.086%
31	11.816	1492	1501	1511	rBV	63452	135323	14.04%	1.057%
32	13.863	1844	1849	1854	rBV	219052	295174	30.62%	2.306%
33	13.910	1854	1857	1862	rVB	33813	41103	4.26%	0.321%
34	14.157	1892	1899	1910	rBB	240243	347238	36.03%	2.712%

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 Title : SVOA CALIBRATION

35	14.463	1945	1951	1958	rBB2	290133	440412	45.69%	3.440%
36	14.963	2031	2036	2042	rBB	31259	38776	4.02%	0.303%
37	15.451	2113	2119	2130	rBB	329033	462268	47.96%	3.611%
38	16.181	2239	2243	2248	rVB	11043	13231	1.37%	0.103%
39	16.275	2254	2259	2262	rBV	32685	43422	4.51%	0.339%
40	16.328	2262	2268	2275	rVB2	41139	74032	7.68%	0.578%
41	16.392	2275	2279	2283	rBV2	37474	48500	5.03%	0.379%
42	16.434	2283	2286	2291	rVB	21362	25049	2.60%	0.196%
43	16.486	2291	2295	2297	rBV	9357	12032	1.25%	0.094%
44	16.592	2307	2313	2320	rBV4	22205	38348	3.98%	0.300%
45	16.657	2320	2324	2328	rVV	37498	52978	5.50%	0.414%
46	16.692	2328	2330	2334	rVV	13020	19497	2.02%	0.152%
47	16.734	2334	2337	2342	rVB2	16823	21827	2.26%	0.170%
48	17.204	2410	2417	2423	rBV2	384078	551158	57.18%	4.305%
49	17.304	2428	2434	2446	rVB2	342239	476944	49.48%	3.725%
50	18.122	2568	2573	2577	rVB5	6282	10968	1.14%	0.086%
51	18.281	2591	2600	2603	rBV4	4584	11631	1.21%	0.091%
52	19.263	2764	2767	2774	rVB	24407	31955	3.32%	0.250%
53	19.604	2816	2825	2838	rVB	556290	765421	79.41%	5.979%
54	19.916	2874	2878	2881	rBV	8426	10176	1.06%	0.079%
55	20.104	2902	2910	2912	rBV	13954	21878	2.27%	0.171%
56	20.127	2912	2914	2919	rVB2	20745	24089	2.50%	0.188%
57	20.427	2961	2965	2968	rBV2	6956	10147	1.05%	0.079%
58	20.604	2991	2995	2999	rBV2	21423	22173	2.30%	0.173%
59	21.069	3071	3074	3080	rBV2	43798	62876	6.52%	0.491%
60	21.251	3101	3105	3107	rBV	304900	395668	41.05%	3.091%
61	21.269	3107	3108	3114	rVB	320772	279684	29.02%	2.185%
62	21.398	3123	3130	3134	rBV	762378	963853	100.00%	7.529%
63	21.433	3134	3136	3144	rVB	40457	47190	4.90%	0.369%
64	21.504	3144	3148	3155	rBV	57066	85486	8.87%	0.668%
65	21.645	3168	3172	3176	rVB2	26968	30664	3.18%	0.240%
66	21.951	3220	3224	3229	rBV2	59782	75530	7.84%	0.590%
67	22.421	3300	3304	3309	rBV	75024	96323	9.99%	0.752%
68	22.557	3323	3327	3332	rVB2	13643	22055	2.29%	0.172%
69	22.939	3387	3392	3399	rVB	83697	123391	12.80%	0.964%
70	23.016	3400	3405	3410	rBV	29750	65562	6.80%	0.512%
71	23.521	3485	3491	3494	rBV2	102201	183428	19.03%	1.433%

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72	23.568	3494	3499	3505	rVB	385375	709746	73.64%	5.544%
73	23.721	3517	3525	3539	rVB2	464827	930246	96.51%	7.266%
74	24.186	3598	3604	3612	rVB	70883	132295	13.73%	1.033%
75	24.951	3728	3734	3744	rVB	58877	127009	13.18%	0.992%
76	25.845	3880	3886	3896	rVB2	33037	76763	7.96%	0.600%

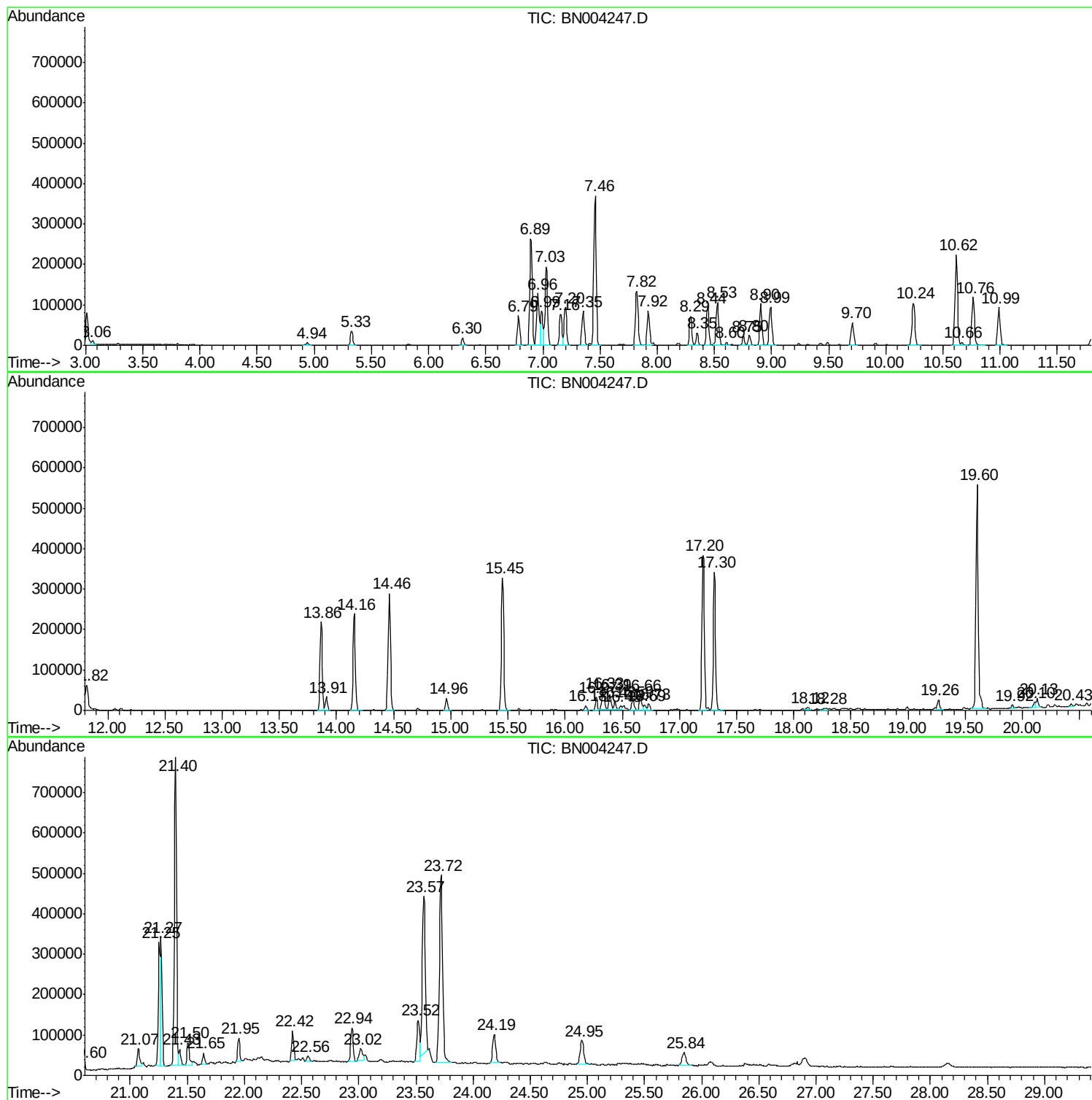
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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



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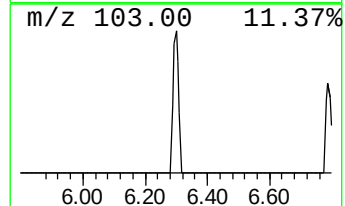
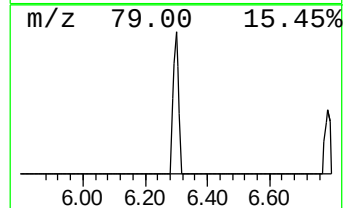
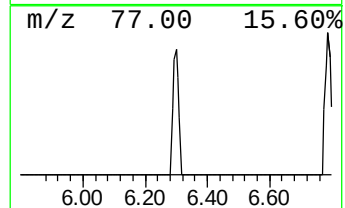
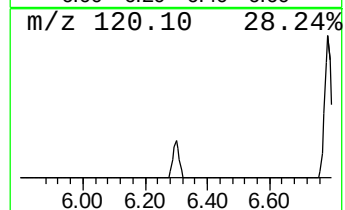
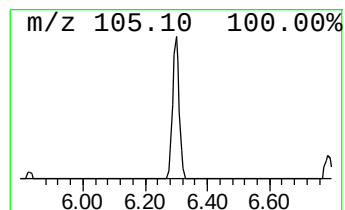
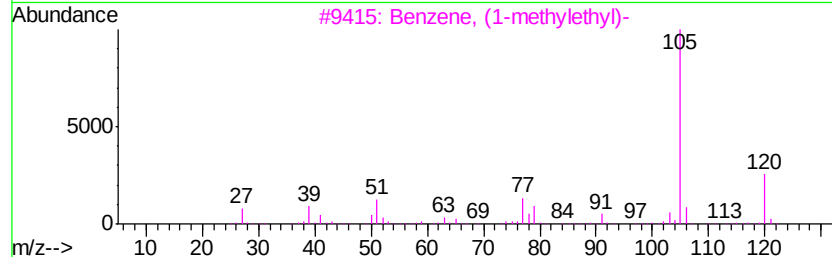
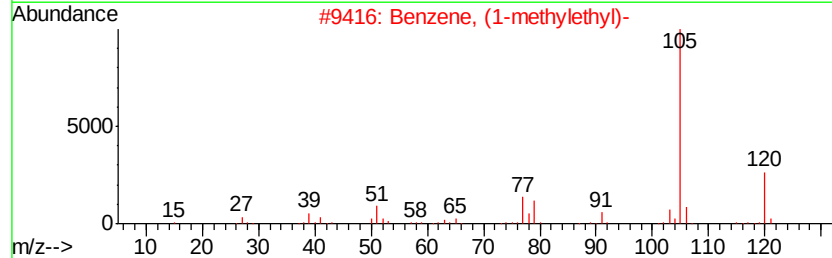
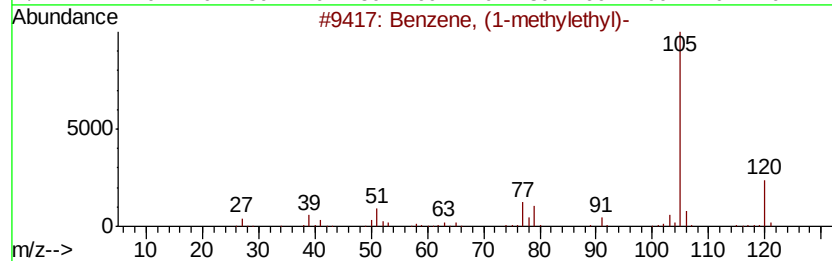
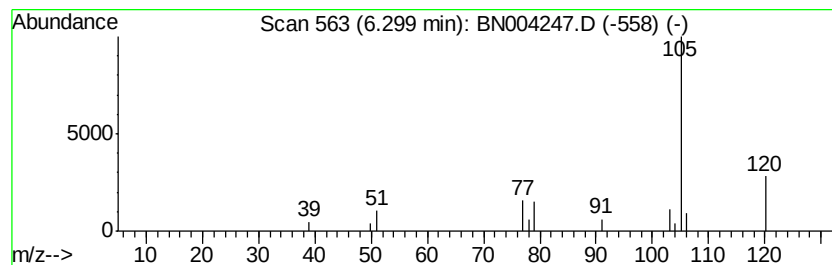
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, (1-methylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.30	2.24 ng/ul	25002	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
2		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
3		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
4		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	83
5		2,4-Nonadiyne	120	C9H12	063621-15-8	78



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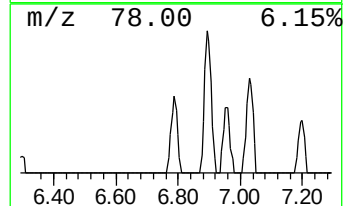
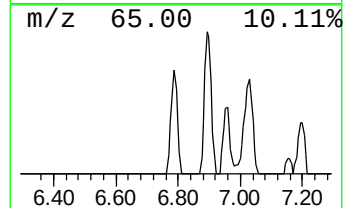
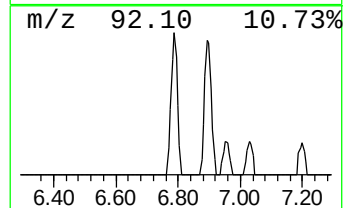
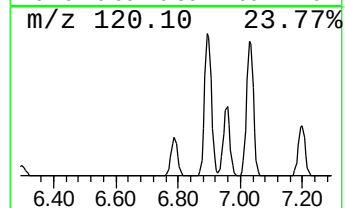
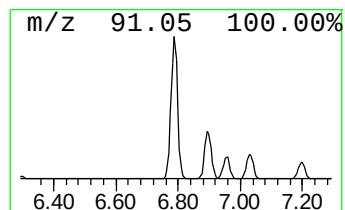
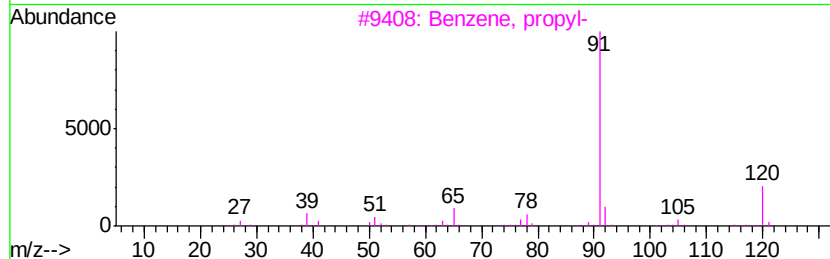
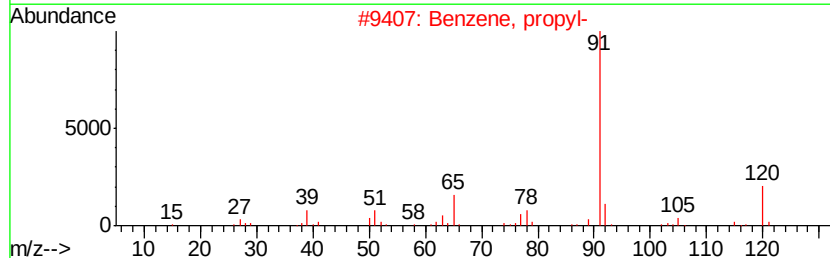
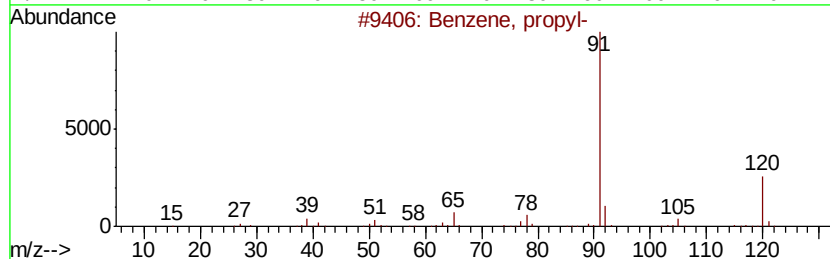
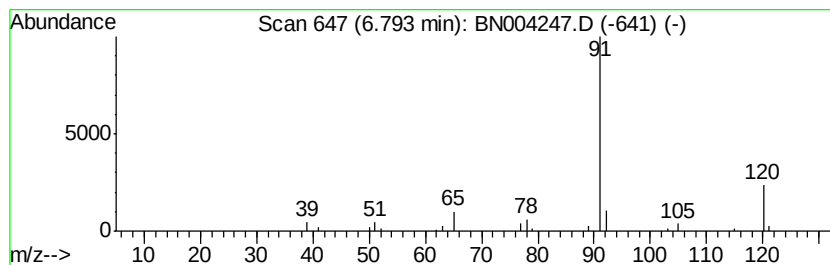
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	9.39 ng/ul	104622	1,4-Dichlorobenzene-d4	7.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, propyl-	120 C9H12	000103-65-1	91
2	Benzene, propyl-	120 C9H12	000103-65-1	91
3	Benzene, propyl-	120 C9H12	000103-65-1	91
4	N-Benzyl-2-phenethylamine	211 C15H17N	003647-71-0	78
5	Ethanol, 2-[(phenylmethyl)amino]-	151 C9H13NO	000104-63-2	72



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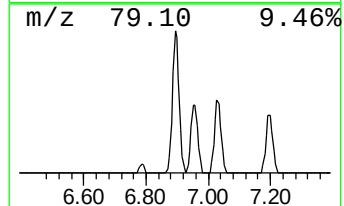
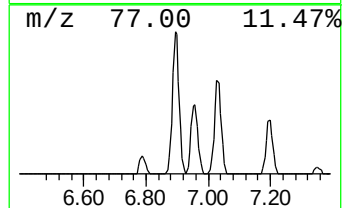
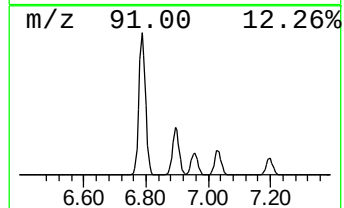
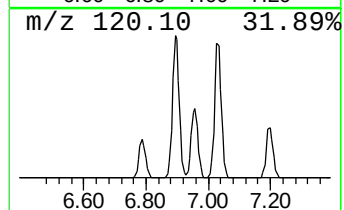
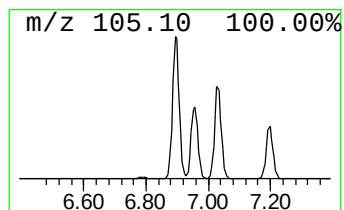
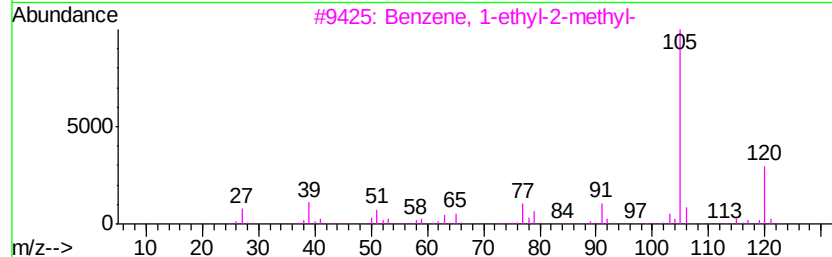
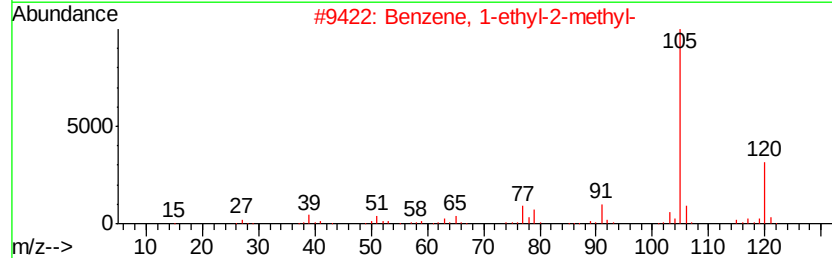
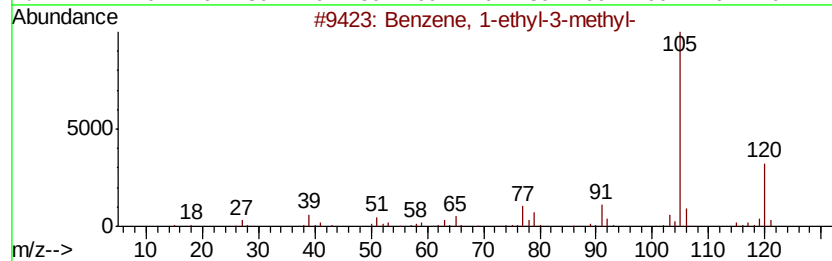
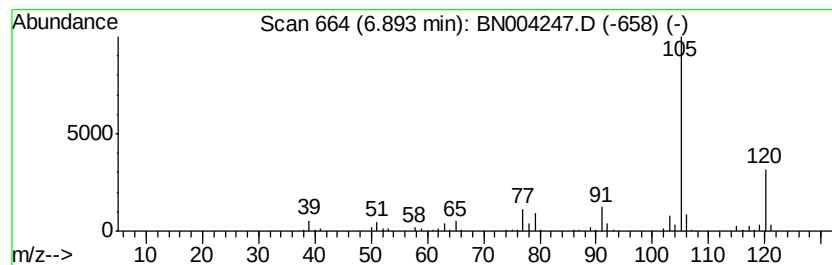
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.89	35.64 ng/ul	397292	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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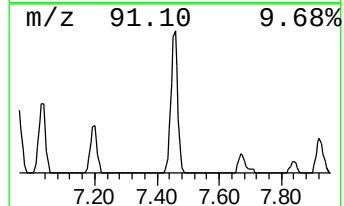
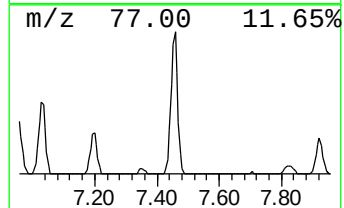
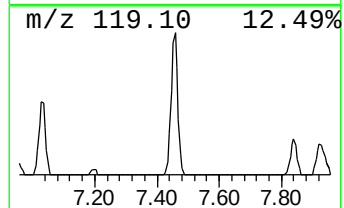
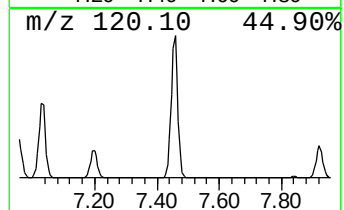
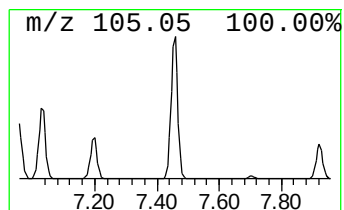
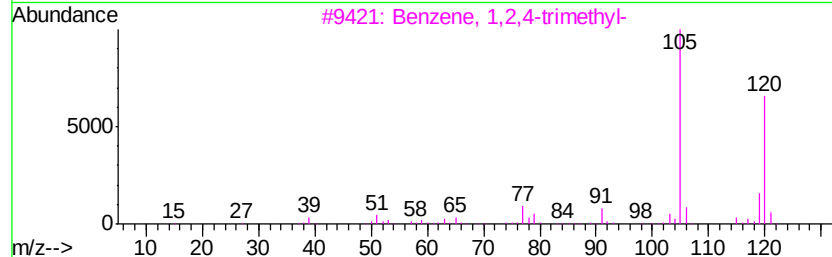
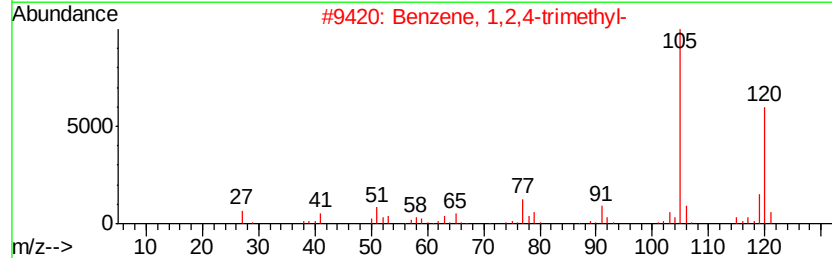
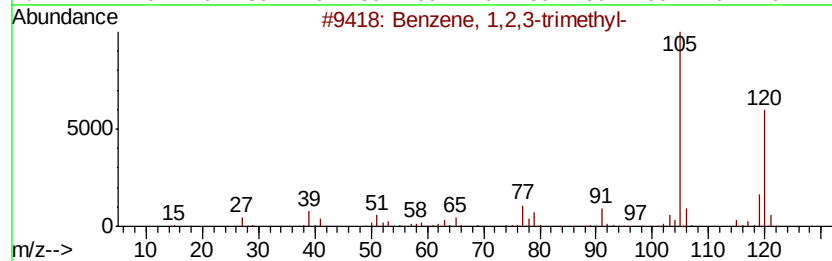
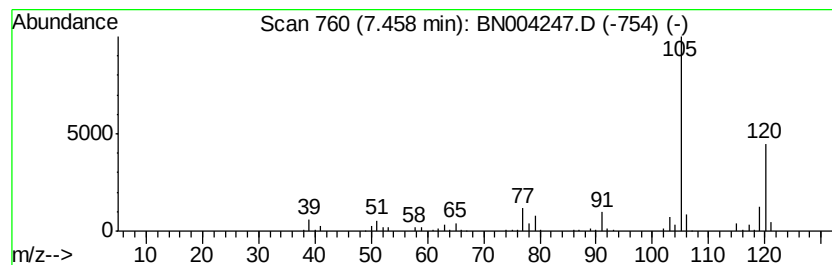
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.46	49.63 ng/ul	553222	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
5		Mesitylene	120	C9H12	000108-67-8	94



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Acq On : 27 Dec 2018 18:50
Operator : JU/SJ
Sample : J6441-09
Misc :
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Instrument :
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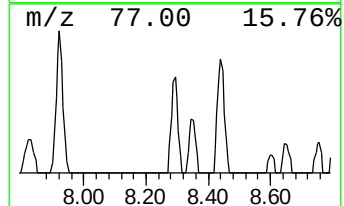
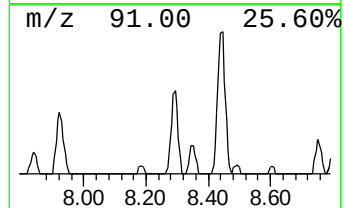
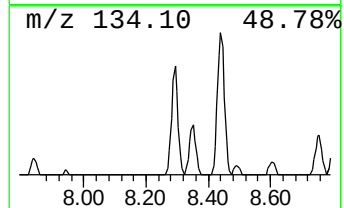
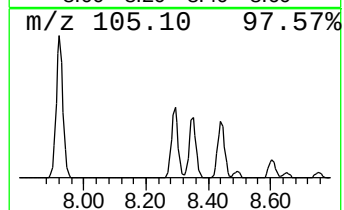
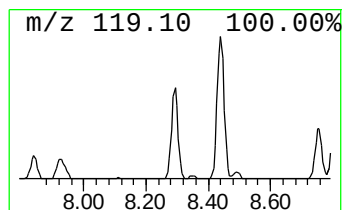
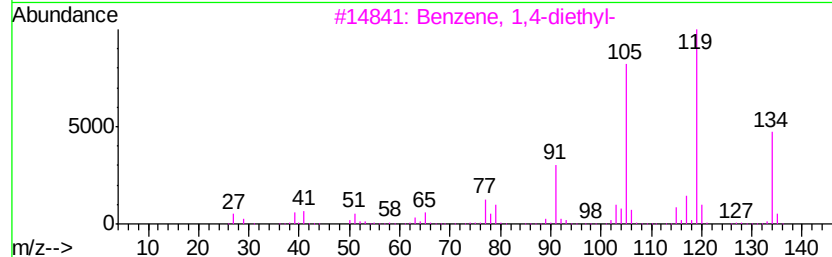
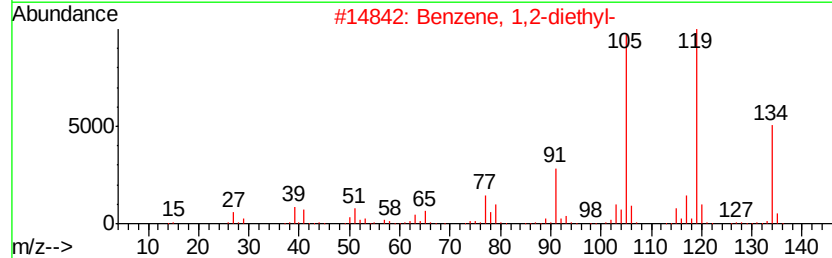
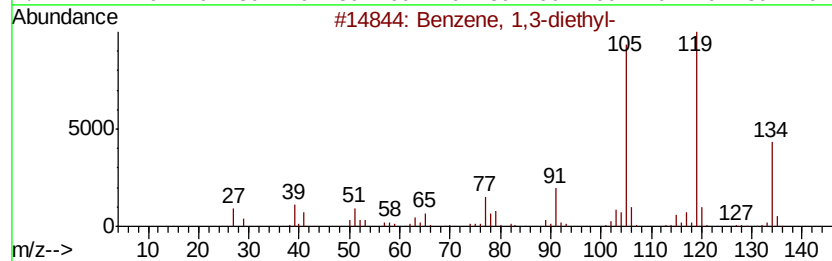
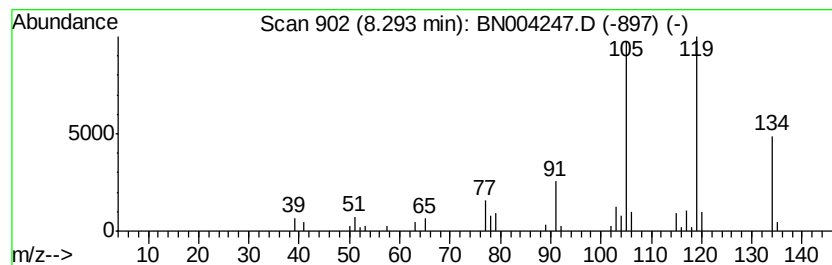
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,3-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.29	9.29 ng/ul	103552	1,4-Dichlorobenzene-d4	7.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
5			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94



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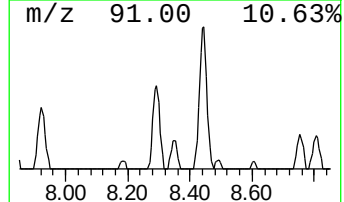
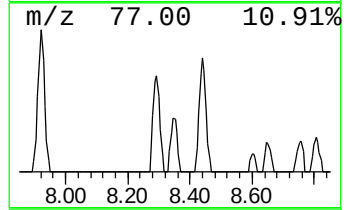
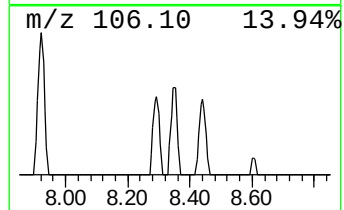
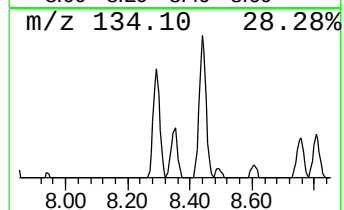
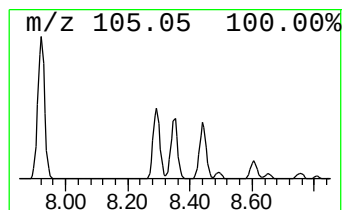
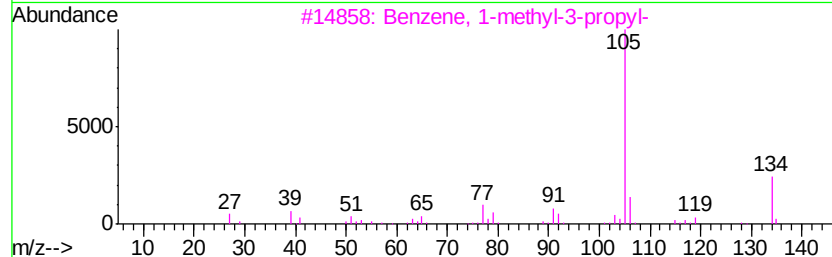
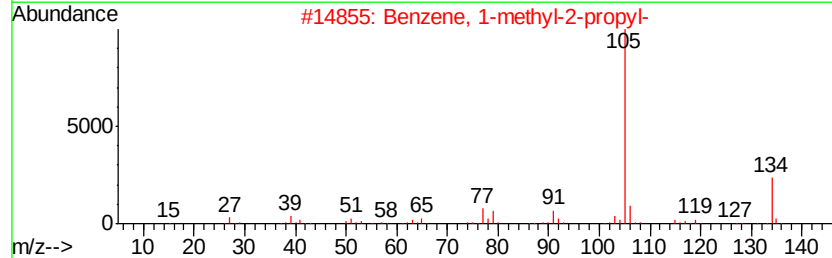
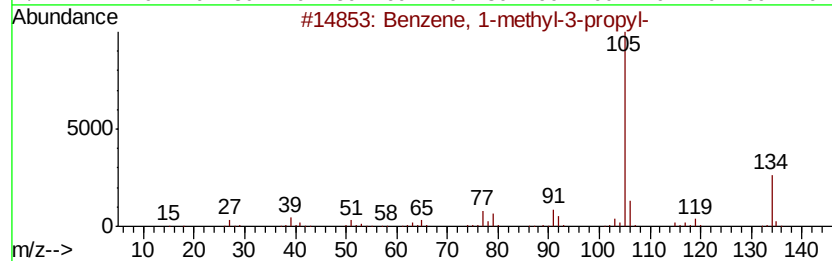
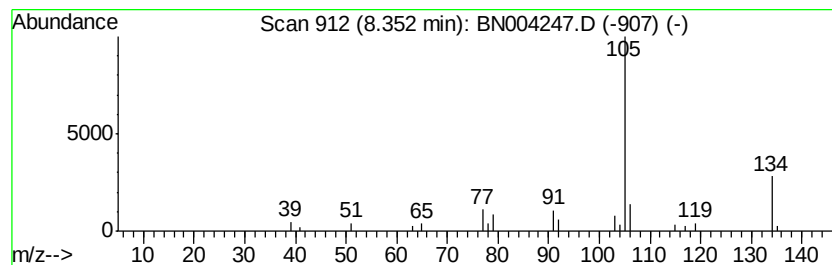
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1-methyl-3-propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.35	3.94 ng/ul	43894	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	94
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
4		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91



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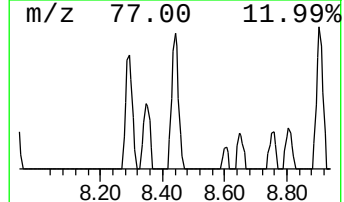
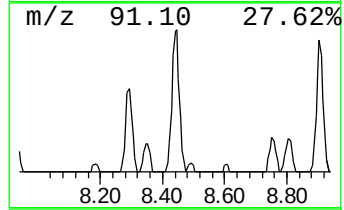
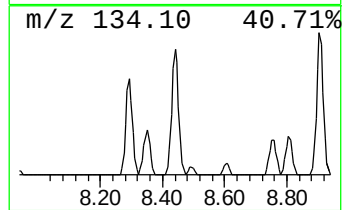
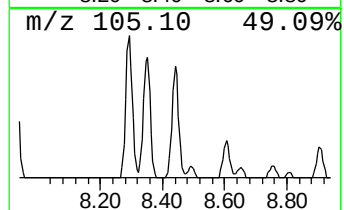
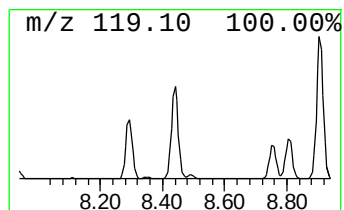
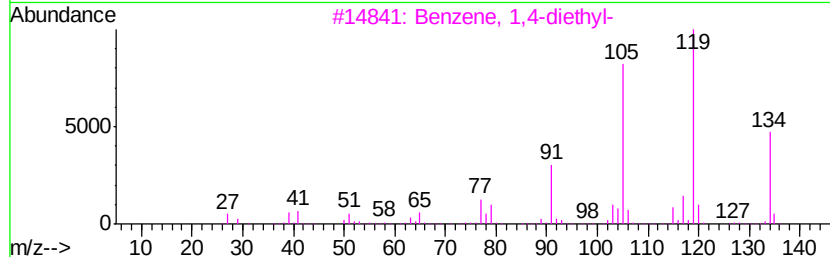
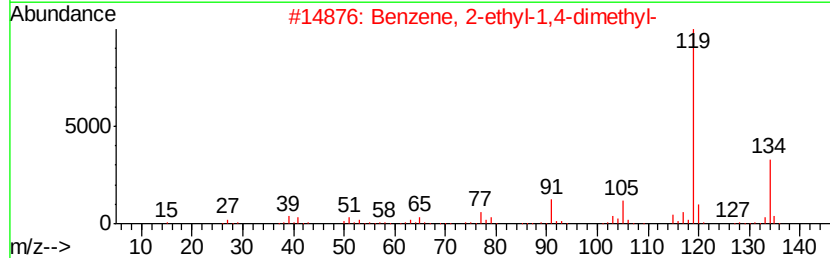
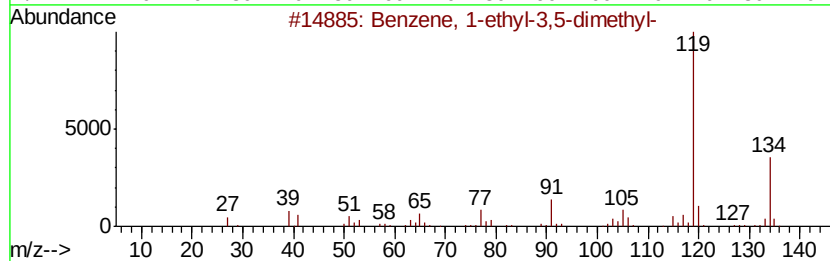
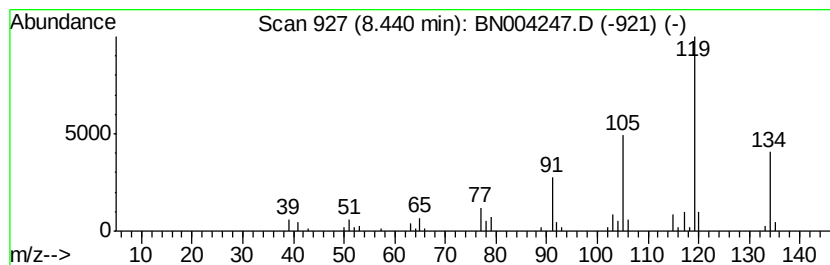
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.44	12.72 ng/ul	141745	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4		o-Cymene	134	C10H14	000527-84-4	93
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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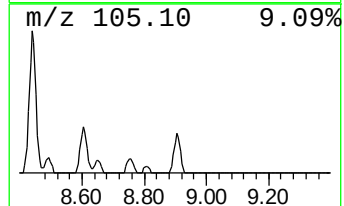
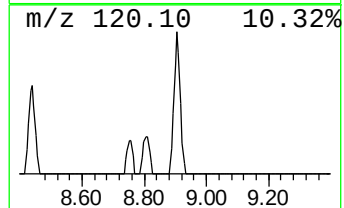
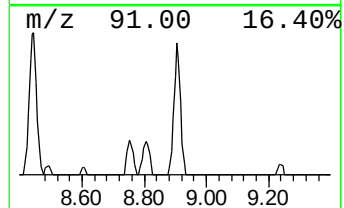
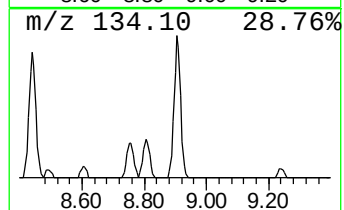
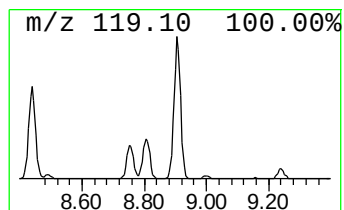
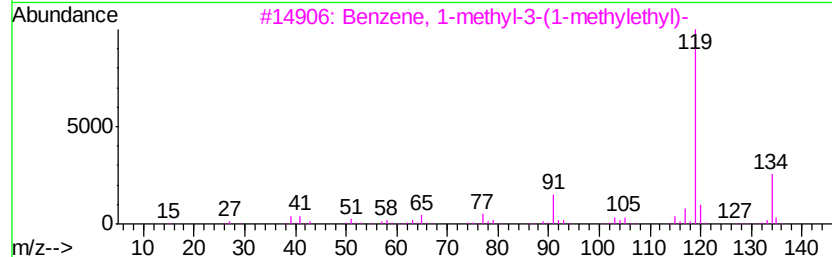
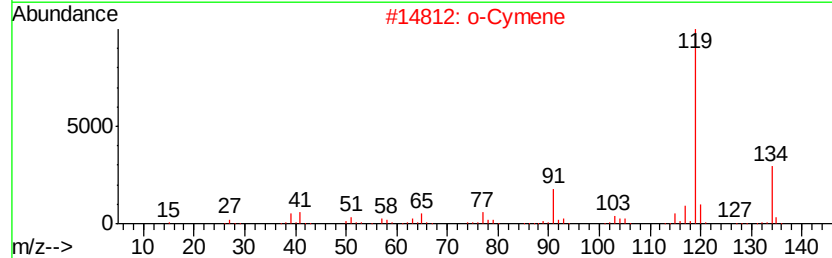
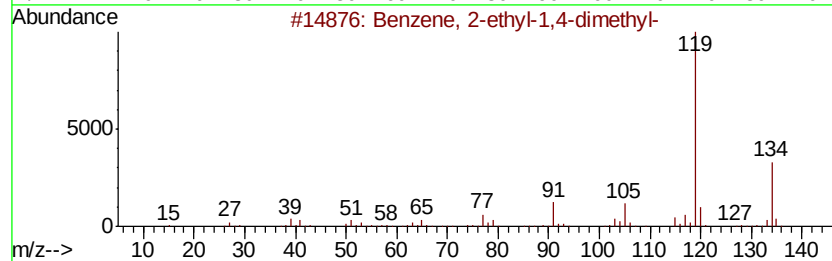
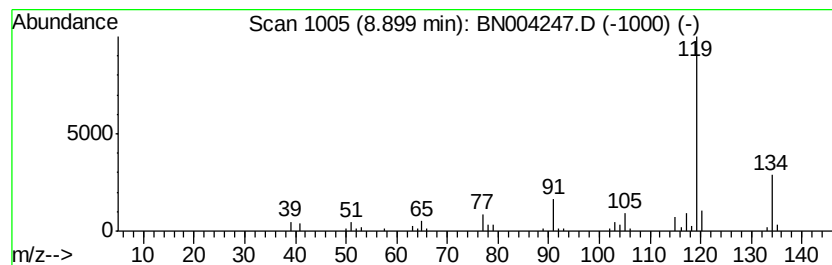
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.90	13.32 ng/ul	148457	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		o-Cymene	134	C10H14	000527-84-4	97
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95



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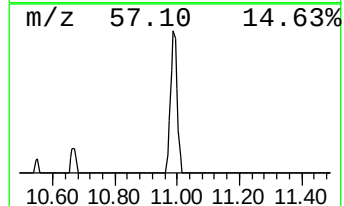
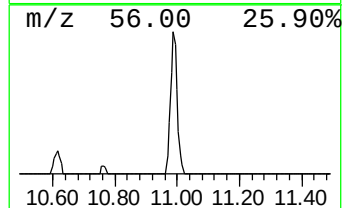
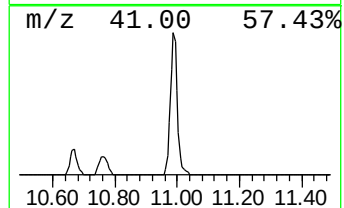
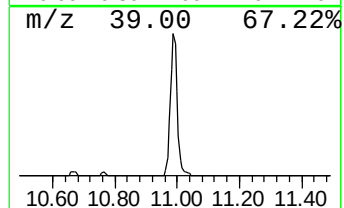
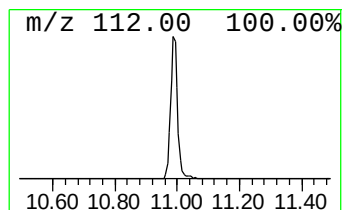
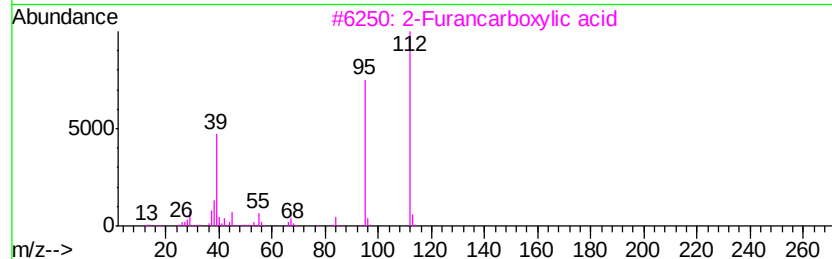
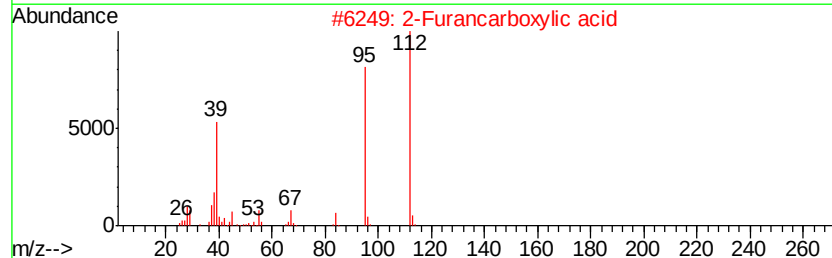
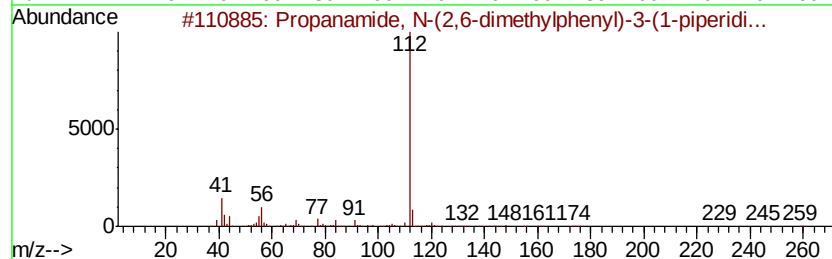
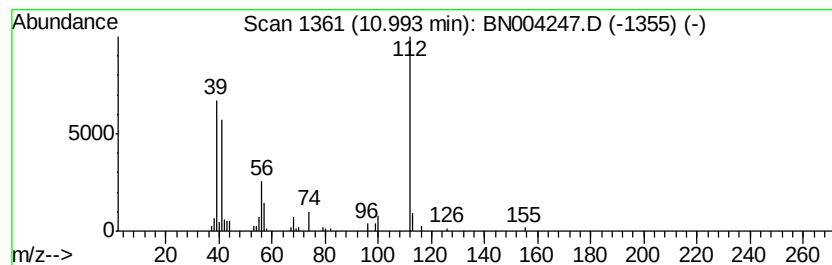
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	7.61 ng/ul	139095	Naphthalene-d8	10.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Propanamide, N-(2,6-dimethylphen...	260 C16H24N2O	004933-38-4 37
2		2-Furancarboxylic acid	112 C5H4O3	000088-14-2 36
3		2-Furancarboxylic acid	112 C5H4O3	000088-14-2 36
4		Benzene, chloro-	112 C6H5Cl	000108-90-7 32
5		N-Methyl-tetrahydro-solasodine	431 C28H49NO2	1000256-11-3 28



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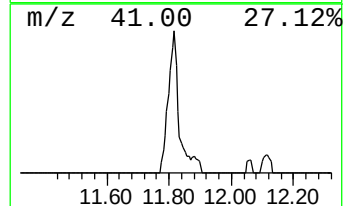
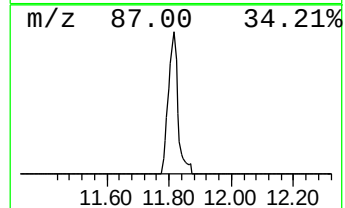
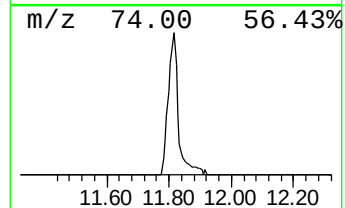
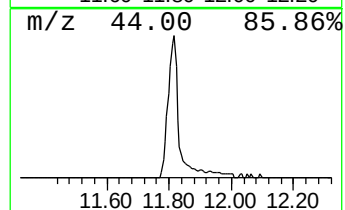
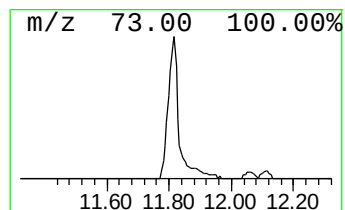
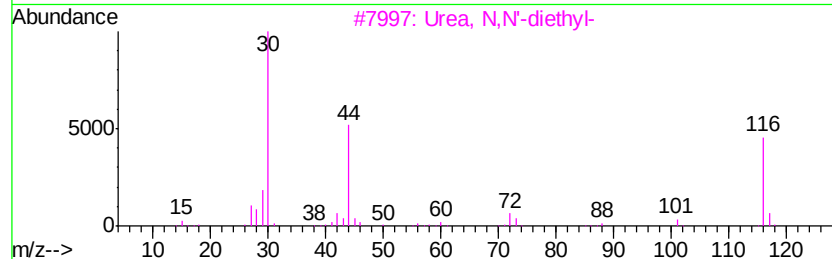
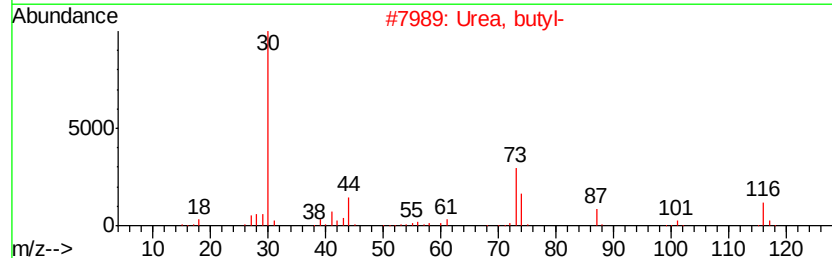
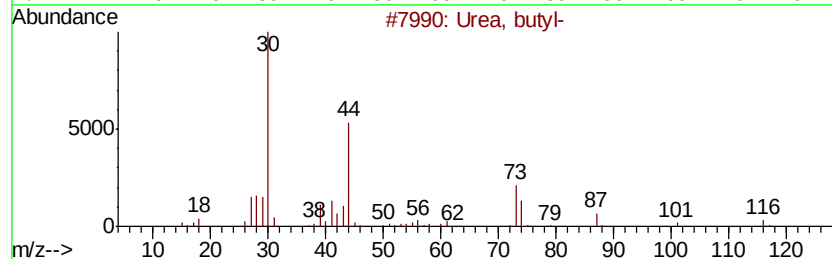
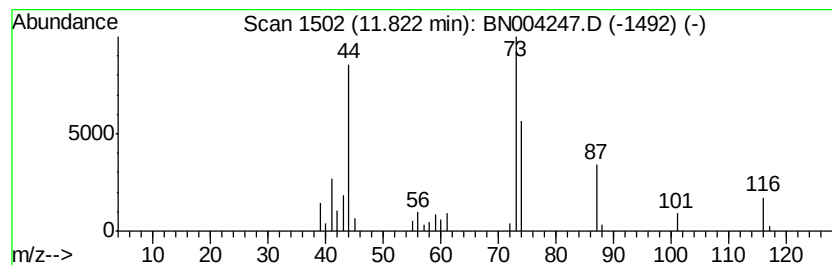
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Urea, butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	7.40 ng/ul	135323	Napthalene-d8	10.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Urea, butyl-	116 C5H12N2O	000592-31-4 91
2		Urea, butyl-	116 C5H12N2O	000592-31-4 50
3		Urea, N,N'-diethyl-	116 C5H12N2O	000623-76-7 46
4		Guanidine, N,N-dimethyl-	87 C3H9N3	006145-42-2 32
5		Guanidine, N,N-dimethyl-	87 C3H9N3	006145-42-2 32



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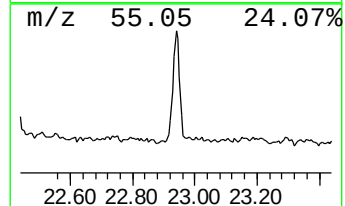
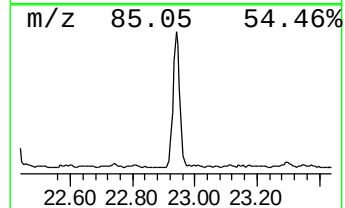
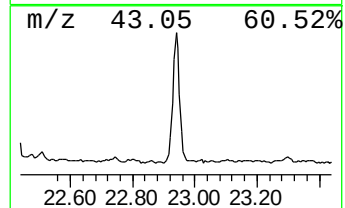
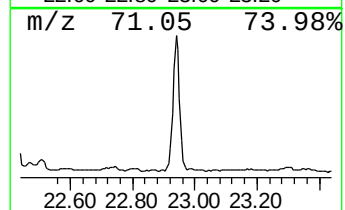
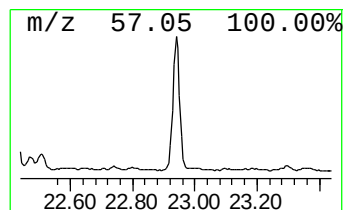
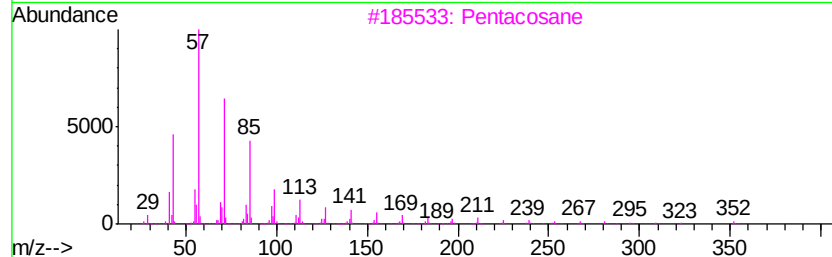
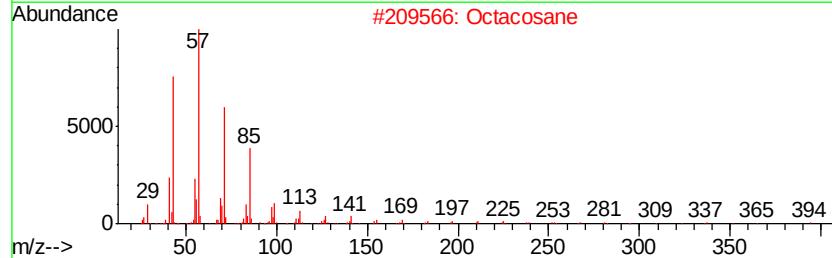
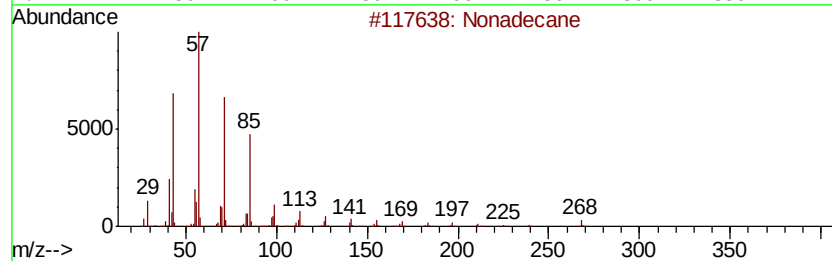
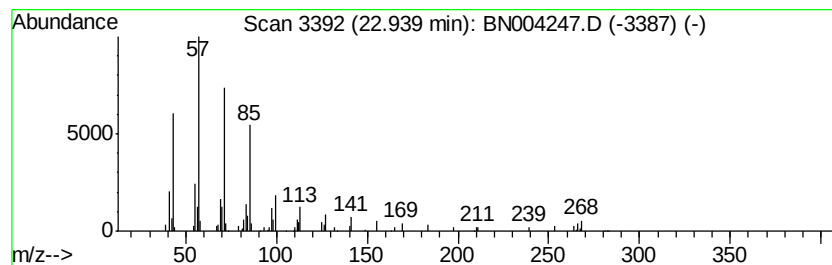
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.94	2.65 ng/ul	123391	Perylene-d12	23.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	96
2		Octacosane	394	C28H58	000630-02-4	90
3		Pentacosane	352	C25H52	000629-99-2	90
4		triacontane	422	C30H62	000638-68-6	90
5		Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122718\
Data File : BN004247.D
Acq On : 27 Dec 2018 18:50
Operator : JU/SJ
Sample : J6441-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BE802

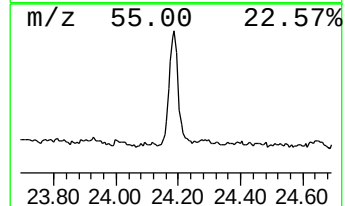
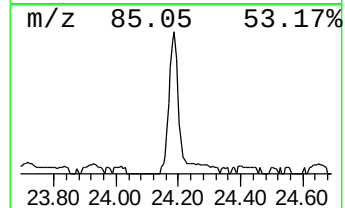
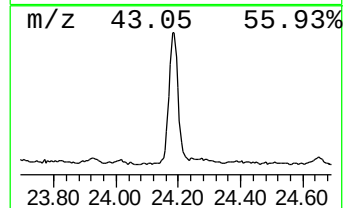
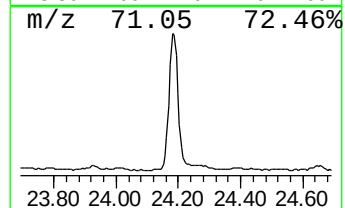
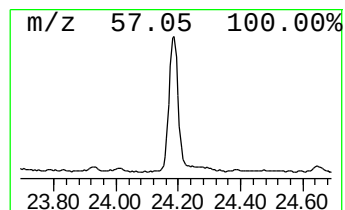
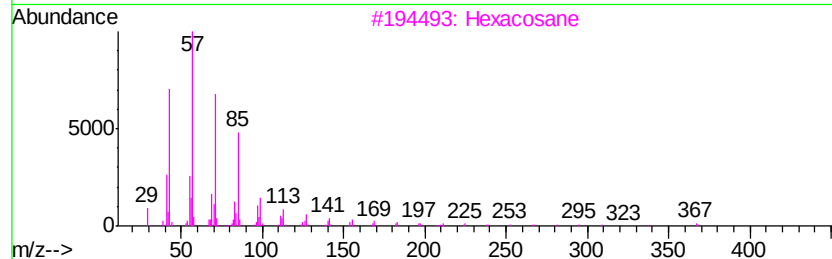
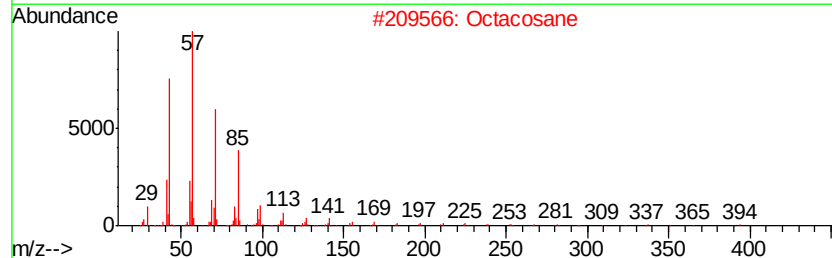
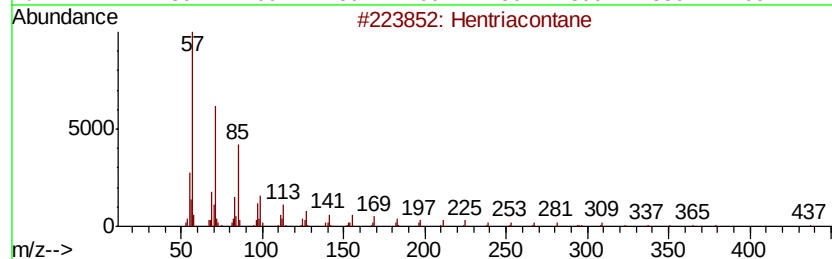
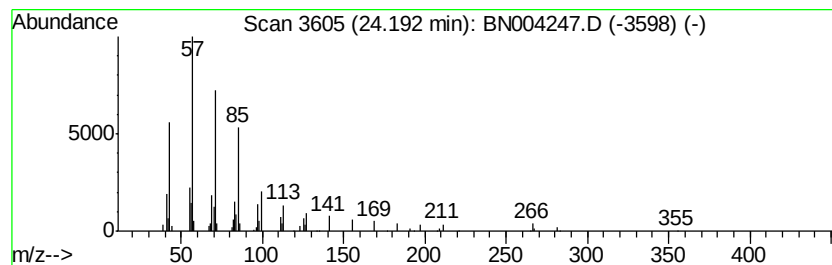
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.19	2.84 ng/ul	132295	Perylene-d12	23.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	94
2			Octacosane	394	C28H58	000630-02-4	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122718\
Data File : BN004247.D
Acq On : 27 Dec 2018 18:50
Operator : JU/SJ
Sample : J6441-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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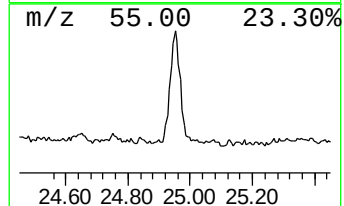
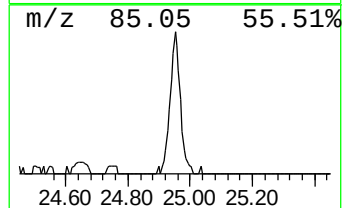
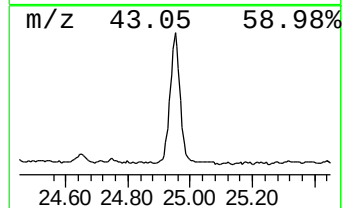
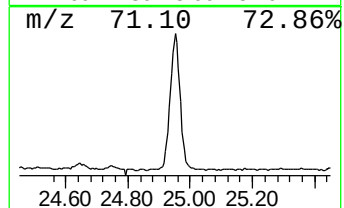
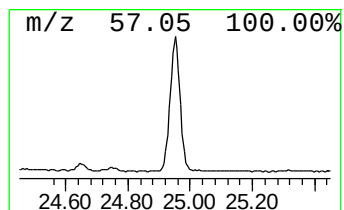
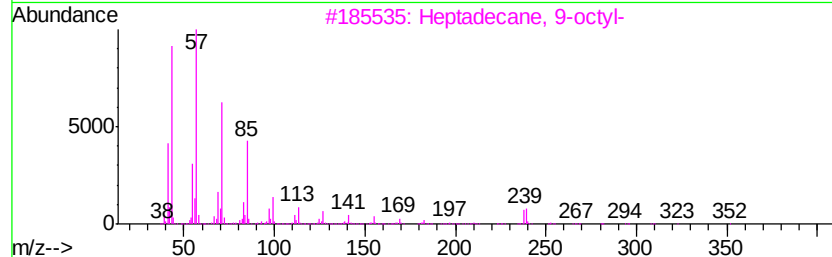
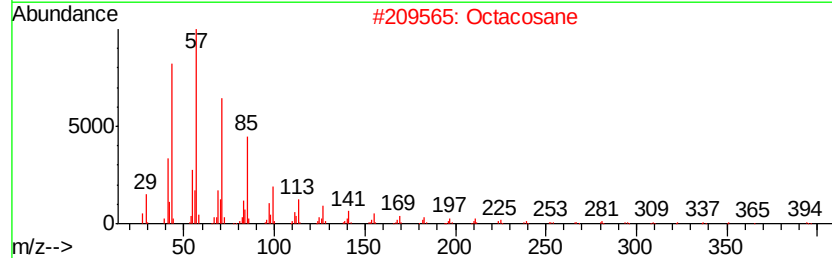
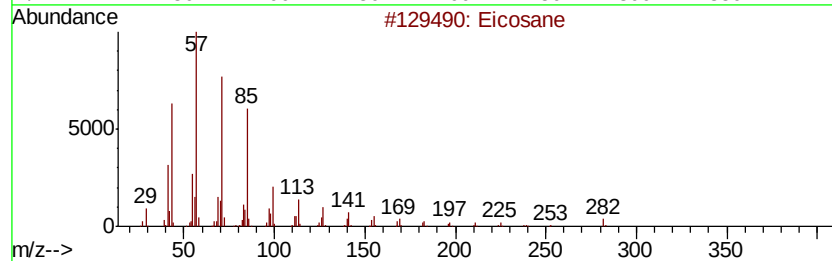
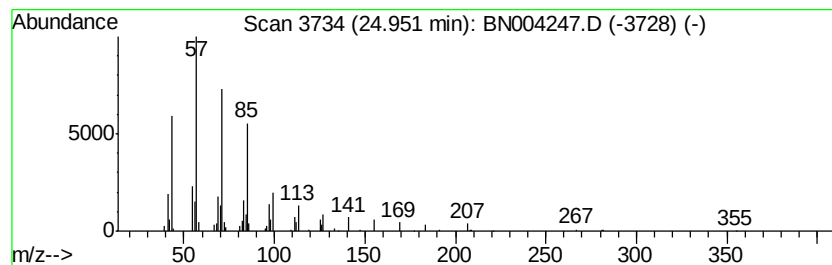
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.95	2.73 ng/ul	127009	Perylene-d12	23.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN122718\
Data File : BN004247.D
Acq On : 27 Dec 2018 18:50
Operator : JU/SJ
Sample : J6441-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BE802

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, (1-methy...	6.30	2.2	ng/ul	25002	1	7.82	222947	20.0
Benzene, propyl-	6.79	9.4	ng/ul	104622	1	7.82	222947	20.0
Benzene, 1-ethyl-...	6.89	35.6	ng/ul	397292	1	7.82	222947	20.0
Benzene, 1,2,3-tr...	7.46	49.6	ng/ul	553222	1	7.82	222947	20.0
Benzene, 1,3-diet...	8.29	9.3	ng/ul	103552	1	7.82	222947	20.0
Benzene, 1-methyl...	8.35	3.9	ng/ul	43894	1	7.82	222947	20.0
Benzene, 1-ethyl-...	8.44	12.7	ng/ul	141745	1	7.82	222947	20.0
Benzene, 2-ethyl-...	8.90	13.3	ng/ul	148457	1	7.82	222947	20.0
unknown-01	10.99	7.6	ng/ul	139095	2	10.62	365659	20.0
Urea, butyl-	11.82	7.4	ng/ul	135323	2	10.62	365659	20.0
(DEL) Alkane: Str...	22.94	2.6	ng/ul	123391	6	23.72	930246	20.0
(DEL) Alkane: Str...	24.19	2.8	ng/ul	132295	6	23.72	930246	20.0
(DEL) Alkane: Str...	24.95	2.7	ng/ul	127009	6	23.72	930246	20.0