

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023445.D
 Acq On : 29 Oct 2019 16:32
 Operator : JU
 Sample : K5423-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE4H8

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_M\METHODS\SOM-EPA-BM102319MA.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	3.152	25	28	32	rBV	112134	127626	1.14%	0.102%
2	4.970	332	337	344	rVB	375105	548948	4.89%	0.439%
3	5.105	357	360	366	rVB3	26173	41132	0.37%	0.033%
4	5.317	388	396	401	rVB	208653	450255	4.01%	0.360%
5	5.658	451	454	461	rVB2	43110	74836	0.67%	0.060%
6	6.134	528	535	546	rVB	590680	923223	8.23%	0.738%
7	6.316	561	566	575	rVB4	80576	137809	1.23%	0.110%
8	6.616	613	617	621	rBV	33186	48168	0.43%	0.039%
9	6.805	642	649	664	rVB2	548439	1093933	9.75%	0.875%
10	6.963	670	676	683	rBV	1360158	2142246	19.09%	1.713%
11	7.022	683	686	690	rVB3	36010	50239	0.45%	0.040%
12	7.163	703	710	718	rBV	1628236	2695883	24.02%	2.156%
13	7.281	725	730	736	rVB2	77437	126868	1.13%	0.101%
14	7.346	737	741	746	rBV4	36322	55159	0.49%	0.044%
15	7.505	764	768	779	rVB9	23421	57820	0.52%	0.046%
16	7.622	782	788	798	rBV	2050291	3339685	29.76%	2.671%
17	7.705	798	802	805	rVV	100841	148815	1.33%	0.119%
18	7.740	805	808	813	rVB2	35923	52032	0.46%	0.042%
19	7.840	821	825	835	rVB3	86603	148100	1.32%	0.118%
20	7.975	843	848	857	rVB2	104410	252962	2.25%	0.202%
21	8.334	903	909	920	rVB	1403951	2225729	19.83%	1.780%
22	8.610	952	956	967	rVB6	39820	82887	0.74%	0.066%
23	8.728	967	976	977	rBV2	72857	100034	0.89%	0.080%
24	8.769	977	983	995	rVV	1740740	2901614	25.85%	2.321%
25	8.952	1009	1014	1024	rVB	141864	230726	2.06%	0.185%
26	9.240	1058	1063	1070	rVB4	41585	74442	0.66%	0.060%
27	9.310	1070	1075	1081	rVB3	21839	37736	0.34%	0.030%
28	9.422	1083	1094	1097	rBV3	81162	155947	1.39%	0.125%
29	9.487	1097	1105	1115	rVV	1403182	2497410	22.25%	1.997%
30	9.663	1131	1135	1140	rBV4	43521	72705	0.65%	0.058%
31	9.787	1151	1156	1158	rBV	41269	66851	0.60%	0.053%
32	9.822	1158	1162	1170	rVB2	125745	220947	1.97%	0.177%
33	10.028	1189	1197	1205	rBV	2120028	3602762	32.10%	2.881%
34	10.346	1246	1251	1254	rVV3	79582	131915	1.18%	0.105%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023445.D
 Acq On : 29 Oct 2019 16:32
 Operator : JU
 Sample : K5423-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE4H8

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_M\METHODS\SOM-EPA-BM102319MA.M
 Title : SVOA CALIBRATION

35	10.399	1254	1260	1268	rVV	2683549	4579113	40.80%	3.662%
36	10.463	1268	1271	1274	rVV5	22780	35834	0.32%	0.029%
37	10.540	1278	1284	1285	rVV5	24347	40414	0.36%	0.032%
38	10.575	1287	1290	1296	rVV5	29815	51305	0.46%	0.041%
39	10.810	1319	1330	1339	rVB5	50522	174374	1.55%	0.139%
40	11.010	1357	1364	1371	rVV2	61505	128310	1.14%	0.103%
41	11.763	1482	1492	1501	rBV	163934	304809	2.72%	0.244%
42	11.998	1524	1532	1537	rBV2	39575	70961	0.63%	0.057%
43	12.081	1541	1546	1551	rVB5	21611	36243	0.32%	0.029%
44	12.275	1572	1579	1587	rVB6	33676	86305	0.77%	0.069%
45	12.528	1617	1622	1633	rVB6	28778	63839	0.57%	0.051%
46	12.763	1656	1662	1668	rVB3	42425	79445	0.71%	0.064%
47	12.828	1668	1673	1679	rBV5	29303	65839	0.59%	0.053%
48	13.004	1695	1703	1710	rBV5	19632	54065	0.48%	0.043%
49	13.110	1715	1721	1725	rVV7	17606	39764	0.35%	0.032%
50	13.187	1729	1734	1741	rVV	75310	128642	1.15%	0.103%
51	13.569	1792	1799	1809	rVB6	34145	73281	0.65%	0.059%
52	13.687	1812	1819	1824	rVV	3940312	5665722	50.48%	4.531%
53	13.734	1824	1827	1833	rVV	251162	378783	3.38%	0.303%
54	13.798	1833	1838	1842	rVV8	24376	53630	0.48%	0.043%
55	13.963	1856	1866	1877	rVV	4510054	6812567	60.70%	5.448%
56	14.051	1877	1881	1885	rVV	53281	89086	0.79%	0.071%
57	14.269	1912	1918	1924	rBV	4041176	6099087	54.34%	4.878%
58	14.316	1924	1926	1934	rVB3	82342	128915	1.15%	0.103%
59	14.487	1948	1955	1971	rBV	397273	804837	7.17%	0.644%
60	14.710	1987	1993	1997	rVB8	17569	36560	0.33%	0.029%
61	14.975	2034	2038	2043	rVB4	34704	50157	0.45%	0.040%
62	15.034	2043	2048	2052	rBV	48977	67748	0.60%	0.054%
63	15.110	2057	2061	2065	rVB	41845	49735	0.44%	0.040%
64	15.269	2081	2088	2099	rBV	5950976	8584126	76.49%	6.865%
65	15.392	2103	2109	2122	rVV	1523891	2260095	20.14%	1.807%
66	15.510	2124	2129	2137	rVV2	93433	186767	1.66%	0.149%
67	15.786	2171	2176	2181	rBV2	61142	97357	0.87%	0.078%
68	15.951	2199	2204	2209	rBV	129980	210429	1.87%	0.168%
69	16.051	2219	2221	2228	rVB7	30821	43725	0.39%	0.035%
70	16.292	2258	2262	2266	rBV3	30925	58923	0.53%	0.047%
71	16.410	2279	2282	2286	rBV2	25467	35256	0.31%	0.028%

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023445.D
 Acq On : 29 Oct 2019 16:32
 Operator : JU
 Sample : K5423-12
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE4H8

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_M\METHODS\SOM-EPA-BM102319MA.M
 Title : SVOA CALIBRATION

72	16.698	2326	2331	2337	rBV6	40245	77923	0.69%	0.062%
73	16.798	2346	2348	2354	rVB2	22621	35964	0.32%	0.029%
74	17.016	2379	2385	2395	rBV2	4738921	6970000	62.10%	5.574%
75	17.116	2395	2402	2412	rBV	6563771	9101058	81.09%	7.278%
76	17.275	2426	2429	2436	rVB7	27682	51691	0.46%	0.041%
77	17.369	2441	2445	2452	rVB8	33207	62448	0.56%	0.050%
78	17.633	2486	2490	2495	rVB3	47965	62473	0.56%	0.050%
79	17.704	2498	2502	2507	rVV6	39590	86664	0.77%	0.069%
80	17.751	2508	2510	2517	rVB3	33444	38897	0.35%	0.031%
81	17.939	2537	2542	2552	rBV2	111082	210347	1.87%	0.168%
82	18.063	2558	2563	2569	rVB	186844	257437	2.29%	0.206%
83	18.163	2575	2580	2582	rBV	139235	195808	1.74%	0.157%
84	18.210	2582	2588	2600	rVB	298554	572635	5.10%	0.458%
85	18.757	2677	2681	2684	rBV3	44908	68263	0.61%	0.055%
86	19.051	2727	2731	2735	rBV2	84451	111147	0.99%	0.089%
87	19.227	2758	2761	2766	rBV6	36766	53001	0.47%	0.042%
88	19.304	2770	2774	2778	rVV	67878	98910	0.88%	0.079%
89	19.416	2787	2793	2799	rBV	8236999	10654200	94.93%	8.520%
90	19.469	2800	2802	2809	rVB2	50531	62918	0.56%	0.050%
91	20.951	3050	3054	3061	rBV	977725	1298025	11.57%	1.038%
92	21.216	3094	3099	3108	rBV	6224834	7879033	70.20%	6.301%
93	21.398	3127	3130	3134	rBV2	269764	298305	2.66%	0.239%
94	21.827	3200	3203	3209	rBV	441223	535259	4.77%	0.428%
95	22.280	3277	3280	3285	rBV	591273	756832	6.74%	0.605%
96	22.780	3362	3365	3371	rVB	719087	1000327	8.91%	0.800%
97	23.274	3442	3449	3455	rBV	6365292	11223192	100.00%	8.976%
98	23.345	3457	3461	3465	rVV	782252	1216801	10.84%	0.973%
99	23.409	3466	3472	3483	rVB	4537618	8309704	74.04%	6.646%
100	23.980	3566	3569	3577	rVB	662379	1109263	9.88%	0.887%

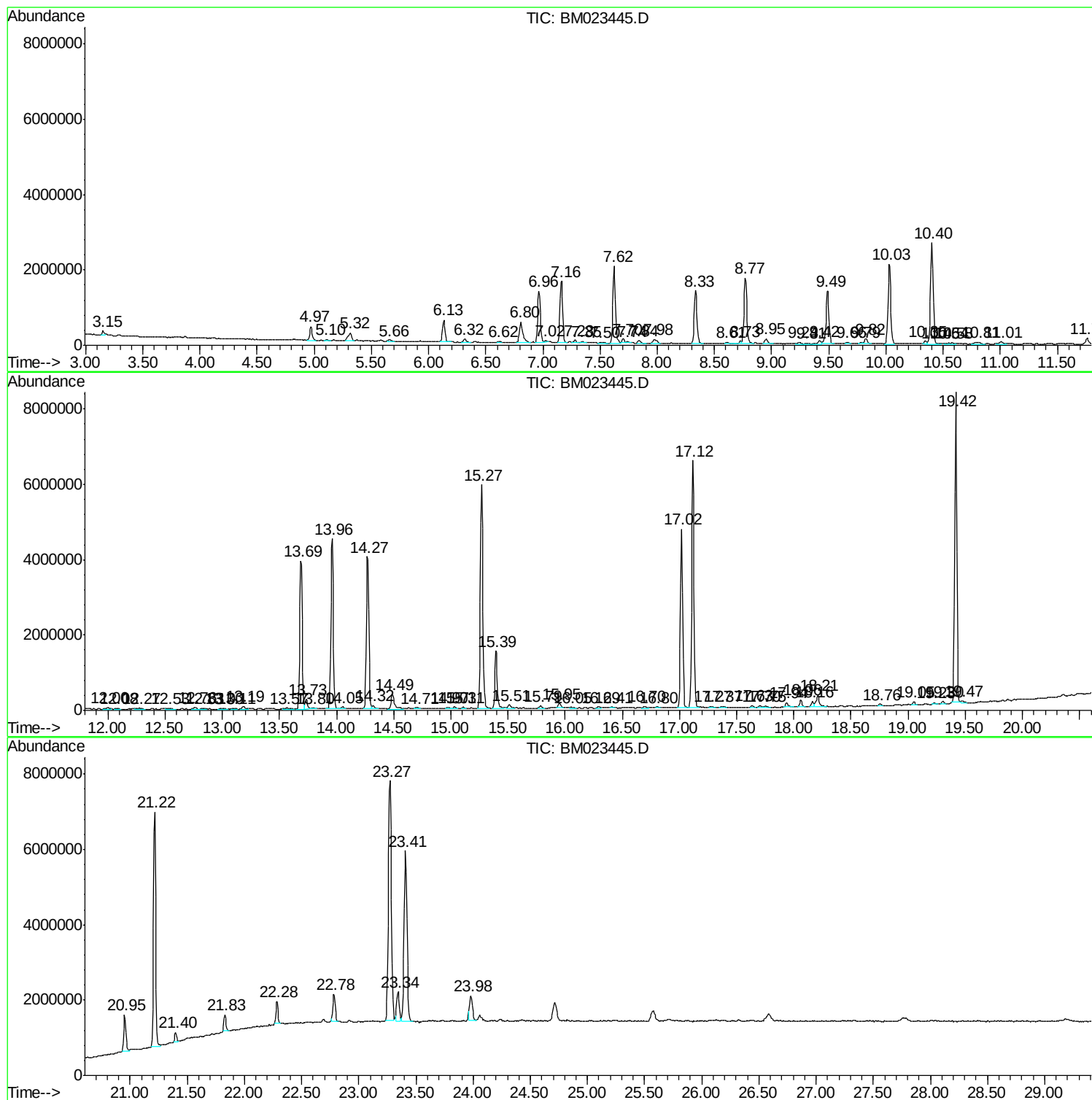
Sum of corrected areas: 125042017

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

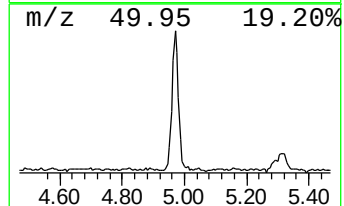
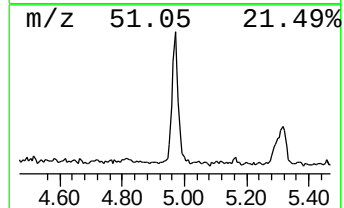
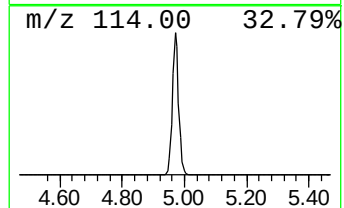
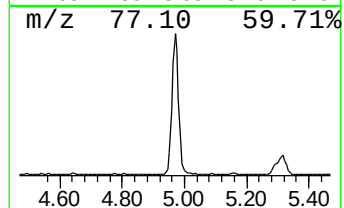
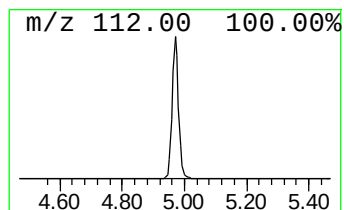
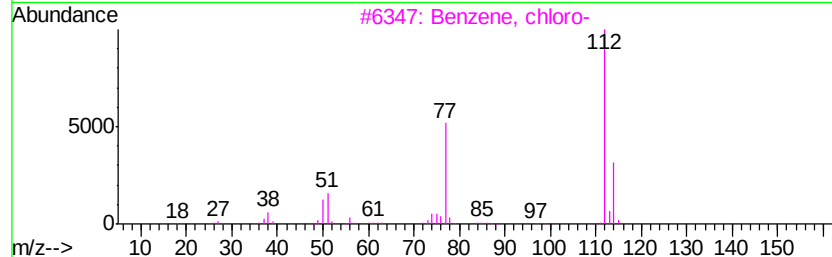
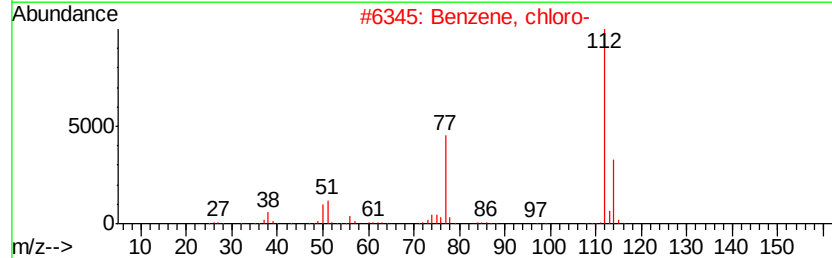
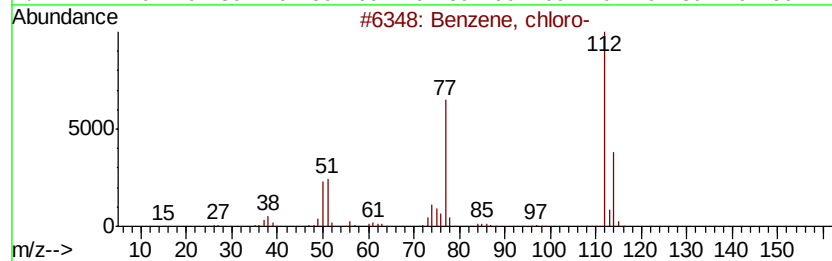
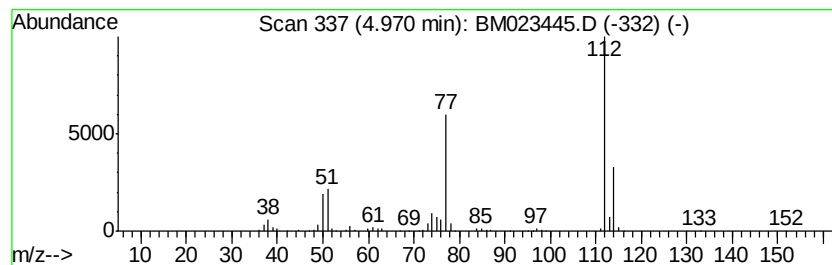
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	3.29 ng/ul	548948	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	91
5		Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
BE4H8

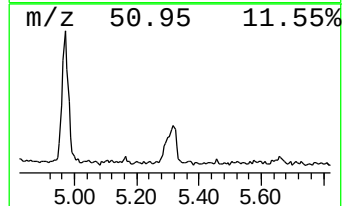
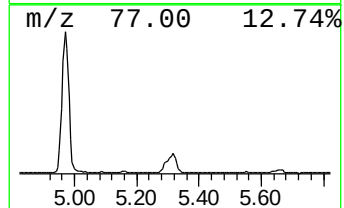
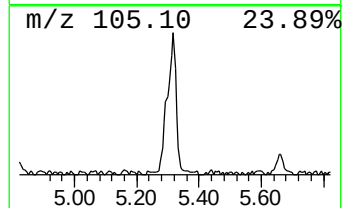
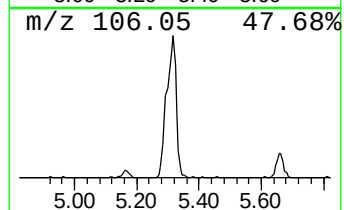
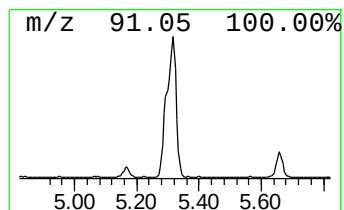
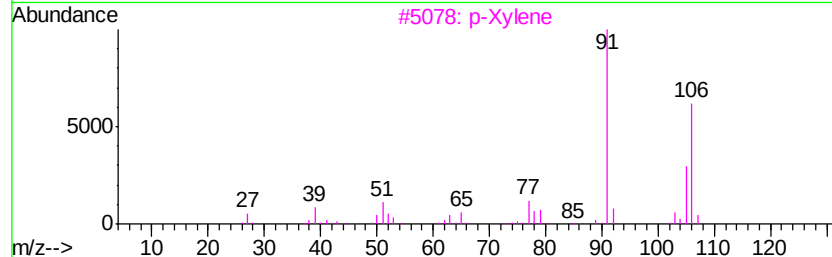
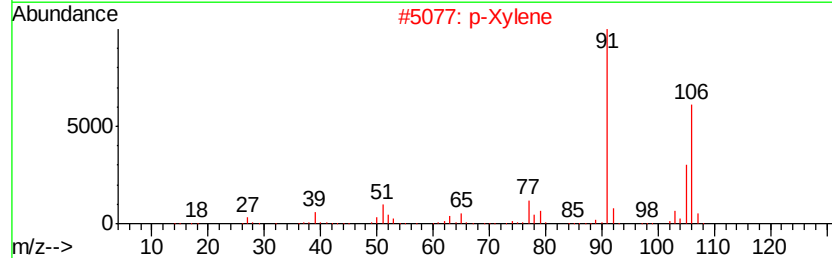
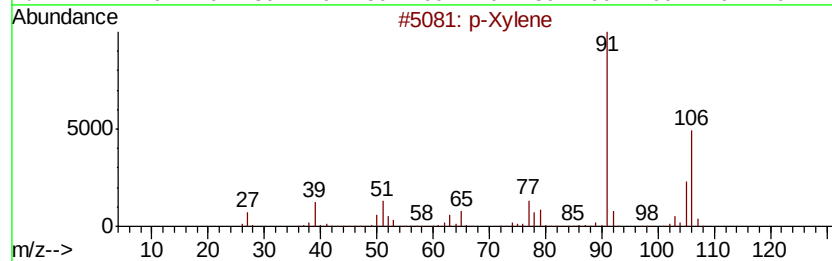
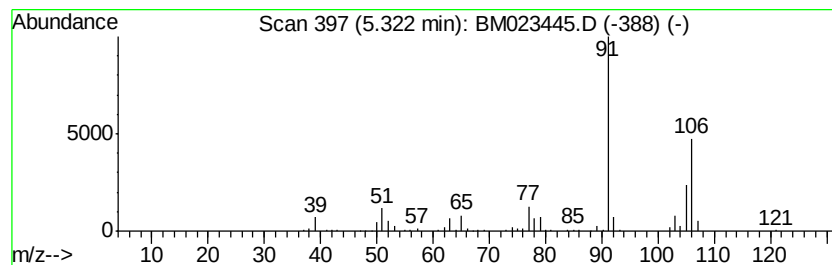
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 p-Xylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	2.70 ng/ul	450255	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Xylene	106	C8H10	000106-42-3	97
2		p-Xylene	106	C8H10	000106-42-3	97
3		p-Xylene	106	C8H10	000106-42-3	97
4		Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
5		p-Xylene	106	C8H10	000106-42-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

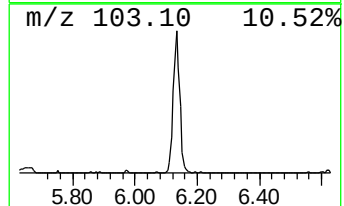
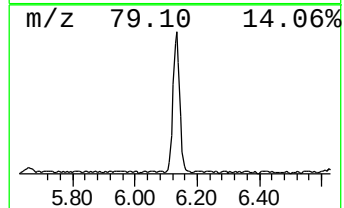
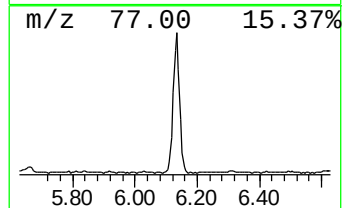
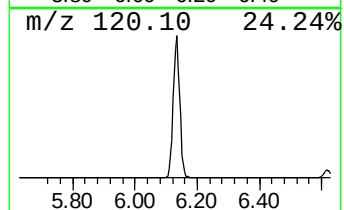
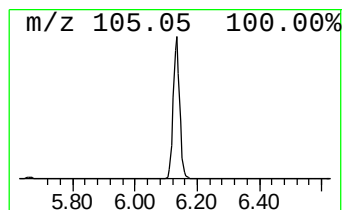
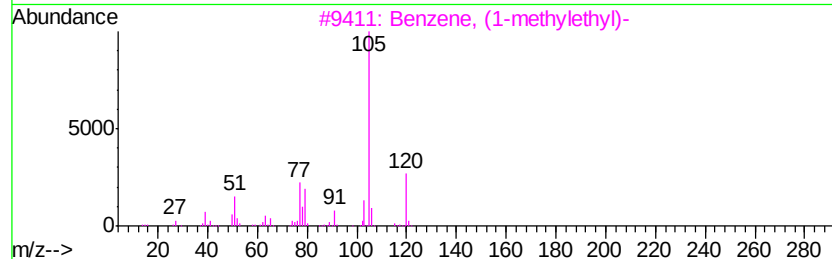
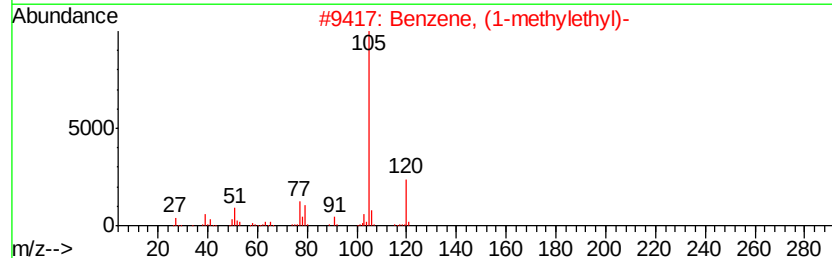
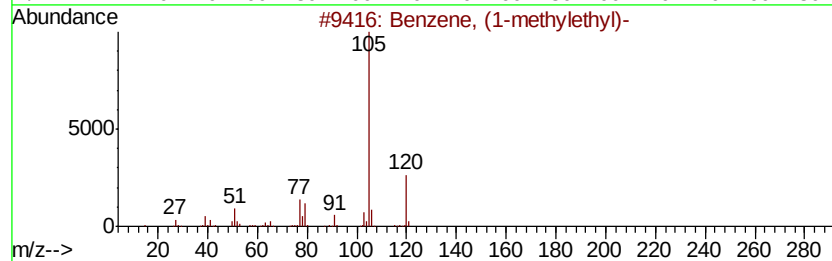
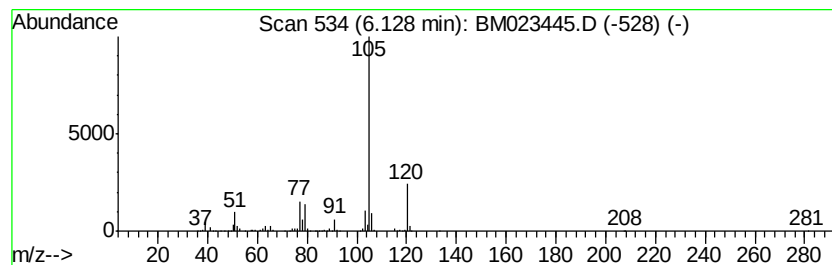
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.13	5.53 ng/ul	923223	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
3		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

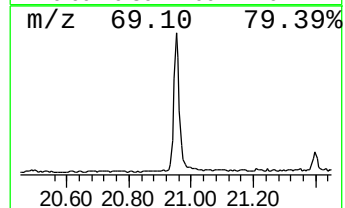
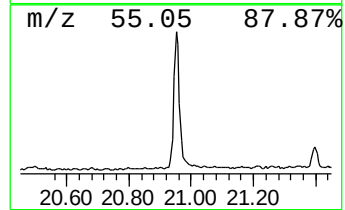
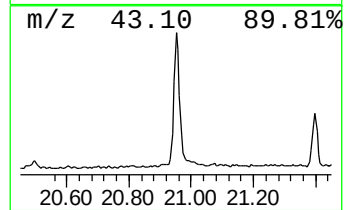
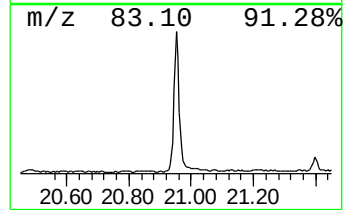
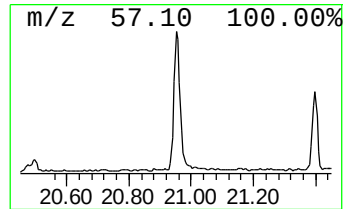
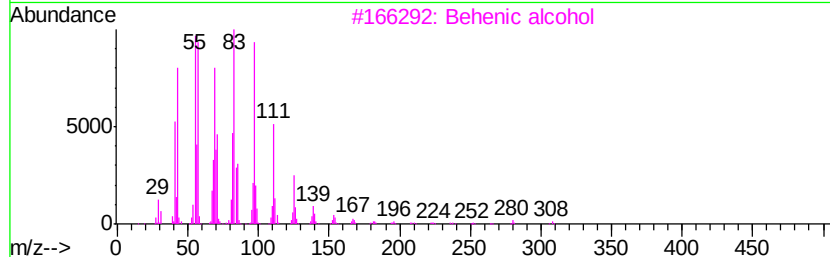
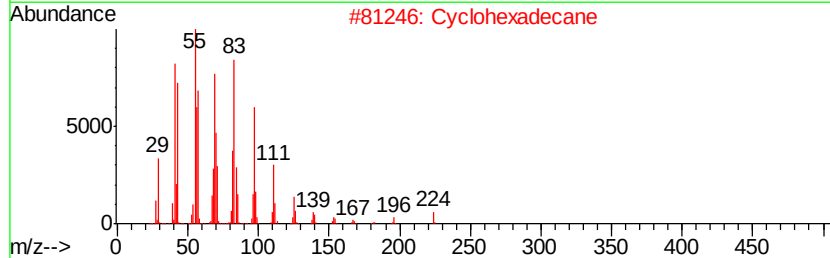
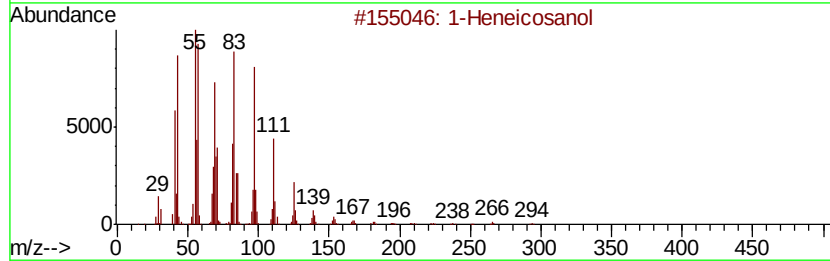
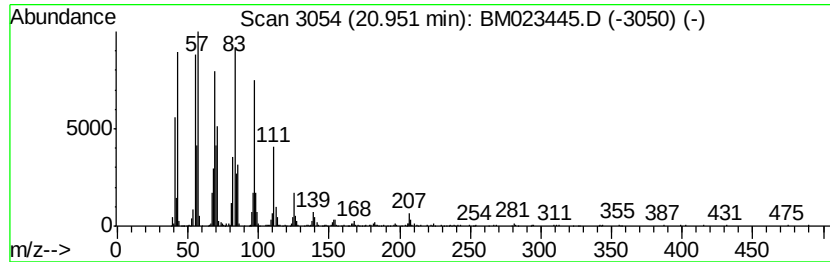
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	3.29 ng/ul	1298030	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	95
2		Cyclohexadecane	224	C16H32	000295-65-8	94
3		Behenic alcohol	326	C22H46O	000661-19-8	94
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

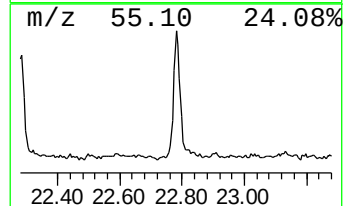
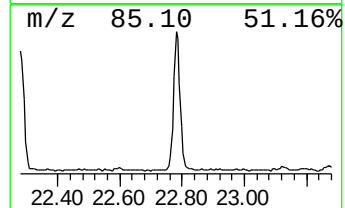
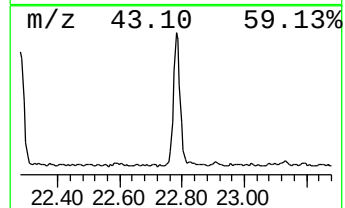
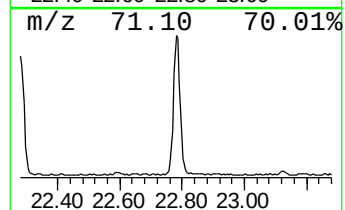
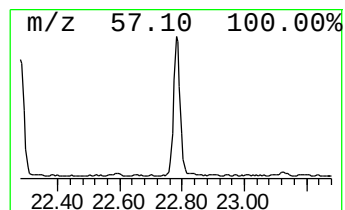
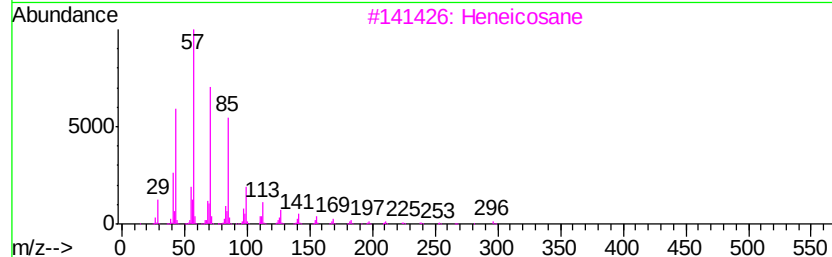
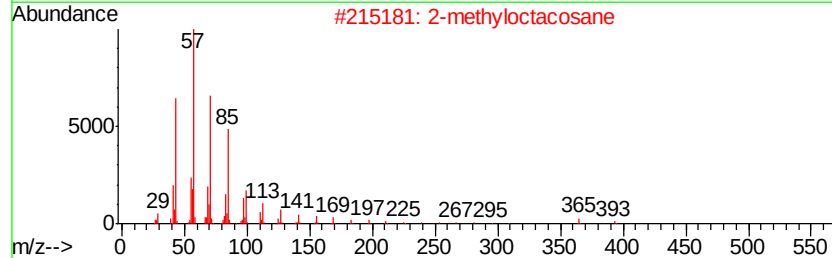
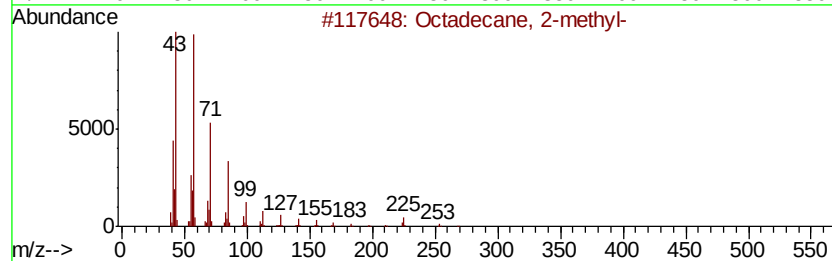
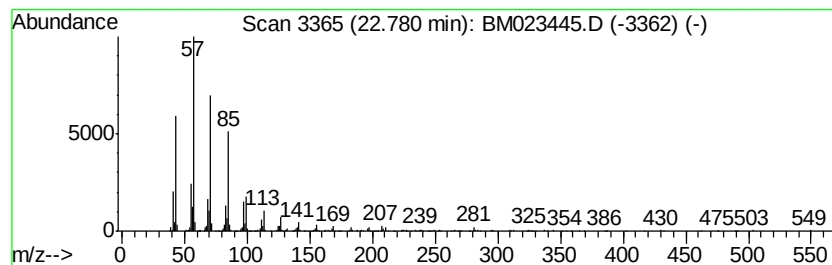
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.78	2.41 ng/ul	1000330	Perylene-d12	23.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Tetratetracontane	619	C44H90	007098-22-8	90
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

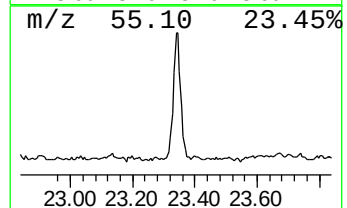
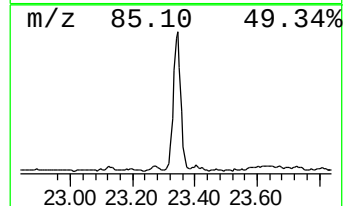
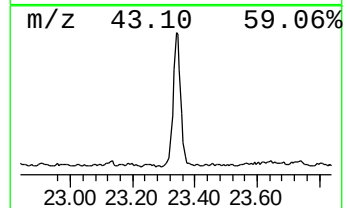
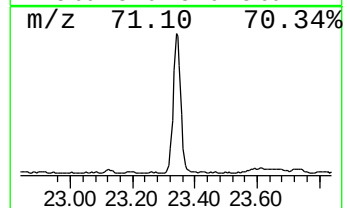
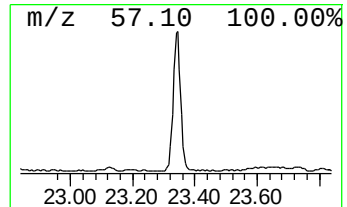
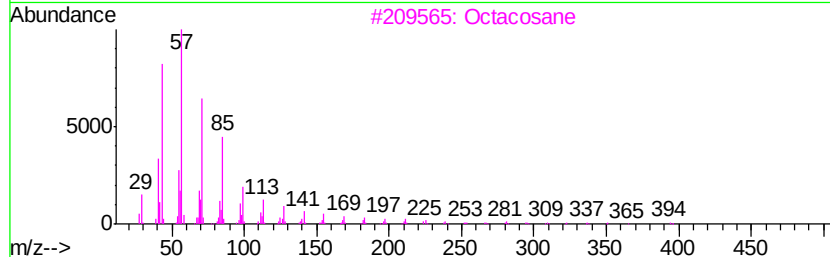
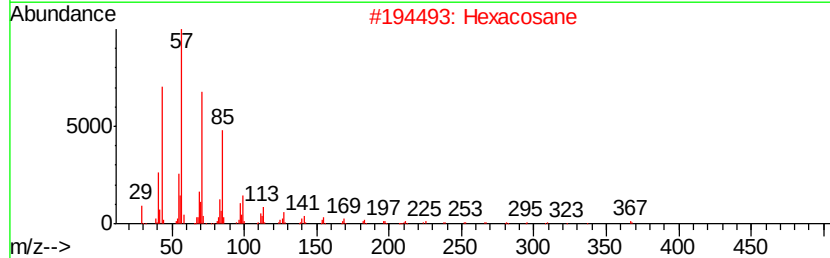
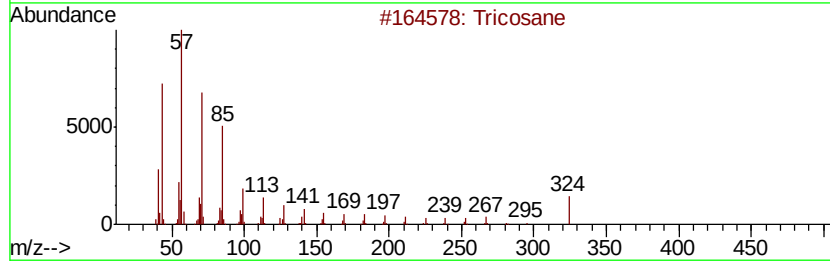
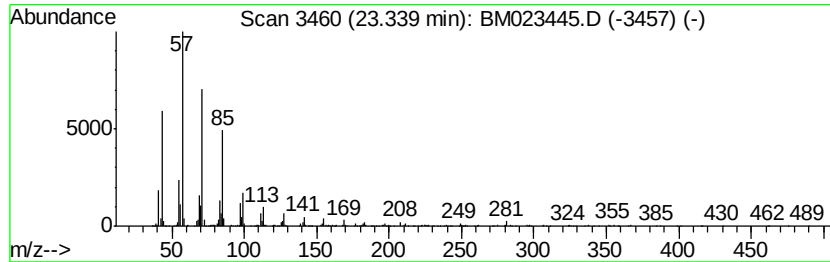
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.34	2.93 ng/ul	1216800	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	93
2		Hexacosane	366	C26H54	000630-01-3	93
3		Octacosane	394	C28H58	000630-02-4	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

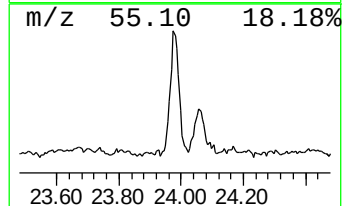
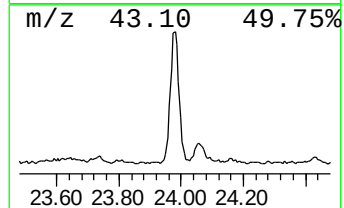
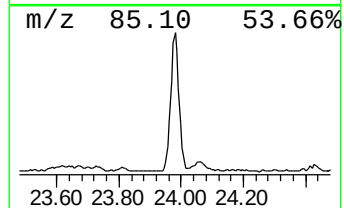
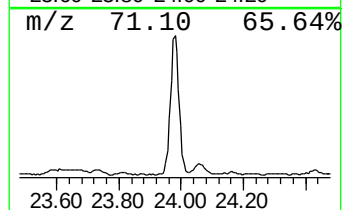
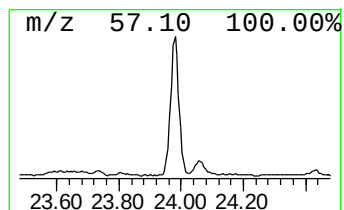
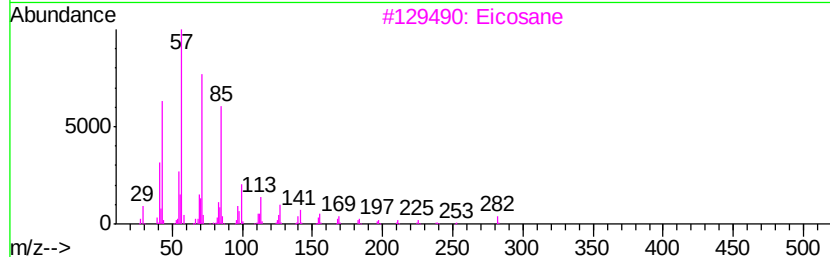
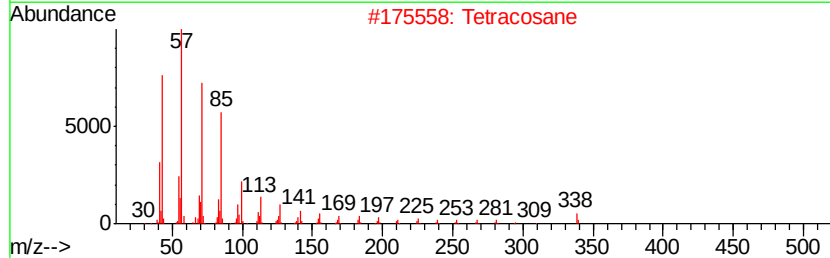
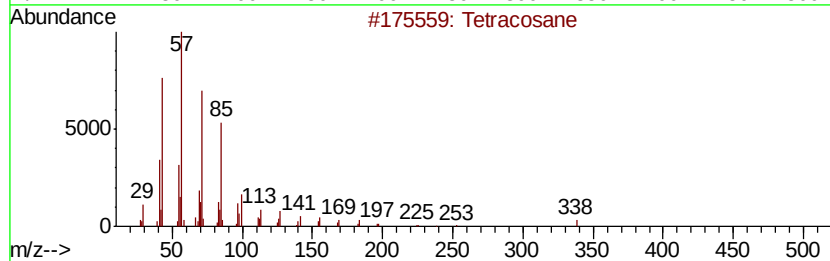
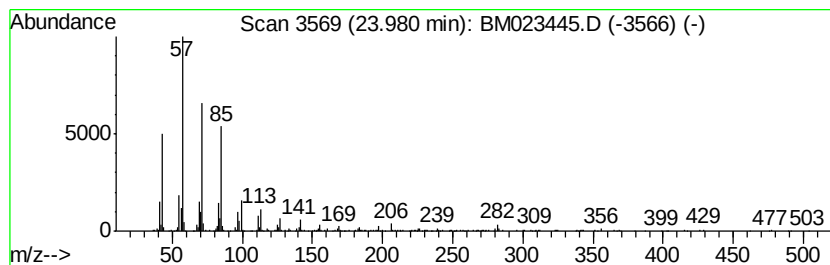
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.98	2.67 ng/ul	1109260	Perylene-d12	23.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	93
2		Tetracosane	338	C24H50	000646-31-1	93
3		Eicosane	282	C20H42	000112-95-8	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102919\
Data File : BM023445.D
Acq On : 29 Oct 2019 16:32
Operator : JU
Sample : K5423-12
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE4H8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.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, chloro-	4.97	3.3	ng/ul	548948	1	7.62	3339690	20.0
p-Xylene	5.32	2.7	ng/ul	450255	1	7.62	3339690	20.0
Benzene, (1-methy...	6.13	5.5	ng/ul	923223	1	7.62	3339690	20.0
1-Heneicosanol	20.95	3.3	ng/ul	1298030	5	21.22	7879030	20.0
(DEL) Alkane: Str...	22.78	2.4	ng/ul	1000330	6	23.41	8309700	20.0
(DEL) Alkane: Str...	23.34	2.9	ng/ul	1216800	6	23.41	8309700	20.0
(DEL) Alkane: Str...	23.98	2.7	ng/ul	1109260	6	23.41	8309700	20.0