

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
 Data File : BM021275.D
 Acq On : 29 Jun 2019 23:05
 Operator : HP/JU
 Sample : K3593-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C5BC5

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-BM061419MA.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	3.016	3	5	24	rVB	1870427	2240423	49.66%	3.212%
2	4.404	237	241	248	rVB	148832	194386	4.31%	0.279%
3	6.857	653	658	663	rBV	97317	139031	3.08%	0.199%
4	6.934	664	671	681	rVB	729331	1212046	26.87%	1.738%
5	7.098	694	699	707	rBV	542597	852107	18.89%	1.222%
6	7.292	726	732	742	rVB	775886	1214719	26.93%	1.742%
7	7.757	805	811	818	rBV	1001856	1579867	35.02%	2.265%
8	7.816	818	821	826	rVB3	11550	19065	0.42%	0.027%
9	7.887	829	833	836	rBV2	11280	13590	0.30%	0.019%
10	8.010	850	854	860	rBV	27841	50852	1.13%	0.073%
11	8.463	924	931	943	rVV	897150	1492474	33.08%	2.140%
12	8.586	947	952	957	rVB3	28504	45067	1.00%	0.065%
13	8.922	1002	1009	1017	rBV	744881	1231938	27.31%	1.766%
14	9.075	1029	1035	1042	rVB3	32065	53734	1.19%	0.077%
15	9.245	1056	1064	1066	rBV7	6363	11534	0.26%	0.017%
16	9.286	1066	1071	1077	rVB	68463	104601	2.32%	0.150%
17	9.639	1123	1131	1140	rBV	631415	1085367	24.06%	1.556%
18	9.822	1156	1162	1166	rBV5	6630	12039	0.27%	0.017%
19	9.957	1180	1185	1189	rBV4	9313	16660	0.37%	0.024%
20	10.028	1194	1197	1201	rVB3	5795	7946	0.18%	0.011%
21	10.069	1201	1204	1211	rVB7	7470	11624	0.26%	0.017%
22	10.169	1213	1221	1233	rBV	1096611	1823465	40.42%	2.615%
23	10.545	1277	1285	1297	rBV	1521989	2514843	55.75%	3.606%
24	10.633	1297	1300	1303	rBV5	6385	7689	0.17%	0.011%
25	10.692	1303	1310	1327	rBV	863549	1571078	34.83%	2.253%
26	11.339	1415	1420	1422	rVV5	7191	12801	0.28%	0.018%
27	11.492	1442	1446	1450	rVB5	8190	12857	0.28%	0.018%
28	11.727	1482	1486	1489	rBV3	8817	12360	0.27%	0.018%
29	11.775	1489	1494	1503	rVB	64660	97446	2.16%	0.140%
30	12.051	1536	1541	1545	rBV2	13842	22703	0.50%	0.033%
31	12.104	1545	1550	1552	rVV5	5457	9508	0.21%	0.014%
32	12.139	1552	1556	1562	rVB	20577	33070	0.73%	0.047%
33	12.210	1565	1568	1571	rBV4	6093	7923	0.18%	0.011%
34	12.292	1577	1582	1587	rVB2	18299	30409	0.67%	0.044%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
 Title : SVOA CALIBRATION

35	12.745	1651	1659	1669	rVV	41728	82527	1.83%	0.118%
36	12.857	1673	1678	1682	rVV4	11667	21177	0.47%	0.030%
37	12.910	1682	1687	1692	rVB5	5857	13430	0.30%	0.019%
38	12.998	1696	1702	1708	rBV	69531	113752	2.52%	0.163%
39	13.204	1732	1737	1740	rBV5	7305	10345	0.23%	0.015%
40	13.263	1742	1747	1750	rVV3	9539	14920	0.33%	0.021%
41	13.357	1757	1763	1768	rVB6	10590	17070	0.38%	0.024%
42	13.816	1830	1841	1845	rBV	2086654	2929549	64.94%	4.201%
43	13.863	1845	1849	1857	rVB	217512	319051	7.07%	0.457%
44	14.045	1873	1880	1882	rBV4	20312	32525	0.72%	0.047%
45	14.092	1882	1888	1896	rVB	2380649	3449714	76.47%	4.946%
46	14.263	1914	1917	1918	rBV	10402	10018	0.22%	0.014%
47	14.398	1934	1940	1948	rBV2	2329635	3586189	79.49%	5.142%
48	14.621	1971	1978	1988	rBV	576719	1018684	22.58%	1.461%
49	14.710	1989	1993	1999	rVV	50566	97720	2.17%	0.140%
50	14.780	1999	2005	2009	rVV9	11478	29628	0.66%	0.042%
51	14.833	2011	2014	2019	rVV5	15093	26105	0.58%	0.037%
52	14.904	2023	2026	2034	rVB6	4631	10903	0.24%	0.016%
53	15.045	2046	2050	2052	rBV4	9781	13229	0.29%	0.019%
54	15.192	2070	2075	2078	rBV	55417	69619	1.54%	0.100%
55	15.227	2078	2081	2088	rVB3	29504	64755	1.44%	0.093%
56	15.345	2097	2101	2104	rBV	28529	39232	0.87%	0.056%
57	15.398	2104	2110	2117	rVB	3027910	4293383	95.17%	6.156%
58	15.527	2126	2132	2143	rBV	541782	792572	17.57%	1.136%
59	15.633	2145	2150	2157	rVV	47697	75682	1.68%	0.109%
60	15.774	2170	2174	2176	rBV4	11797	13146	0.29%	0.019%
61	15.957	2199	2205	2213	rBV6	29963	56653	1.26%	0.081%
62	16.033	2213	2218	2219	rBV5	8507	10998	0.24%	0.016%
63	16.115	2227	2232	2237	rBV2	98788	145412	3.22%	0.208%
64	16.180	2239	2243	2247	rVB5	10884	17117	0.38%	0.025%
65	16.221	2247	2250	2254	rBV4	6969	9908	0.22%	0.014%
66	16.274	2254	2259	2267	rBV10	10059	29516	0.65%	0.042%
67	16.474	2289	2293	2298	rVB4	21741	31868	0.71%	0.046%
68	16.539	2301	2304	2310	rVB7	10315	17699	0.39%	0.025%
69	16.592	2310	2313	2317	rBV	22310	30593	0.68%	0.044%
70	16.692	2326	2330	2334	rBV2	27461	40380	0.90%	0.058%
71	16.762	2340	2342	2349	rBV7	17737	31027	0.69%	0.044%

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72	16.898	2359	2365	2368	rBV5	19181	31705	0.70%	0.045%
73	16.933	2369	2371	2381	rVB8	15093	36009	0.80%	0.052%
74	17.062	2391	2393	2399	rVB6	15663	19106	0.42%	0.027%
75	17.151	2402	2408	2413	rVV2	2737412	3864420	85.66%	5.541%
76	17.192	2413	2415	2418	rVV	76977	84905	1.88%	0.122%
77	17.251	2418	2425	2434	rVB	2948148	4222873	93.61%	6.055%
78	17.433	2453	2456	2459	rVB	19472	18985	0.42%	0.027%
79	17.533	2469	2473	2481	rVB4	42923	67707	1.50%	0.097%
80	17.639	2488	2491	2492	rBV3	9950	10357	0.23%	0.015%
81	17.662	2492	2495	2500	rVB4	27347	36941	0.82%	0.053%
82	17.809	2518	2520	2524	rBV4	21114	22412	0.50%	0.032%
83	18.033	2554	2558	2563	rBV5	51459	97597	2.16%	0.140%
84	18.115	2568	2572	2576	rVV	44118	59405	1.32%	0.085%
85	19.180	2749	2753	2754	rBV	43337	53236	1.18%	0.076%
86	19.209	2754	2758	2763	rVB	188830	259414	5.75%	0.372%
87	19.545	2810	2815	2829	rVB	3134201	4511272	100.00%	6.468%
88	20.074	2903	2905	2909	rVB2	70379	77556	1.72%	0.111%
89	20.886	3040	3043	3047	rBV	208800	243591	5.40%	0.349%
90	21.039	3065	3069	3077	rBV	506917	687027	15.23%	0.985%
91	21.333	3114	3119	3124	rBV	2585603	3280323	72.71%	4.703%
92	21.915	3215	3218	3219	rBV	244872	272050	6.03%	0.390%
93	21.939	3219	3222	3229	rVB	535958	777267	17.23%	1.114%
94	22.615	3333	3337	3344	rVB	741732	1001979	22.21%	1.437%
95	22.891	3380	3384	3389	rBV	1041140	1528331	33.88%	2.191%
96	22.950	3389	3394	3404	rVB2	1520356	2836790	62.88%	4.068%
97	23.462	3475	3481	3491	rVB	2441219	4294766	95.20%	6.158%
98	23.603	3500	3505	3513	rVB	1843385	3410954	75.61%	4.891%
99	23.785	3533	3536	3543	rVB2	446679	725822	16.09%	1.041%
100	24.121	3588	3593	3602	rVB	1006326	1898597	42.09%	2.722%

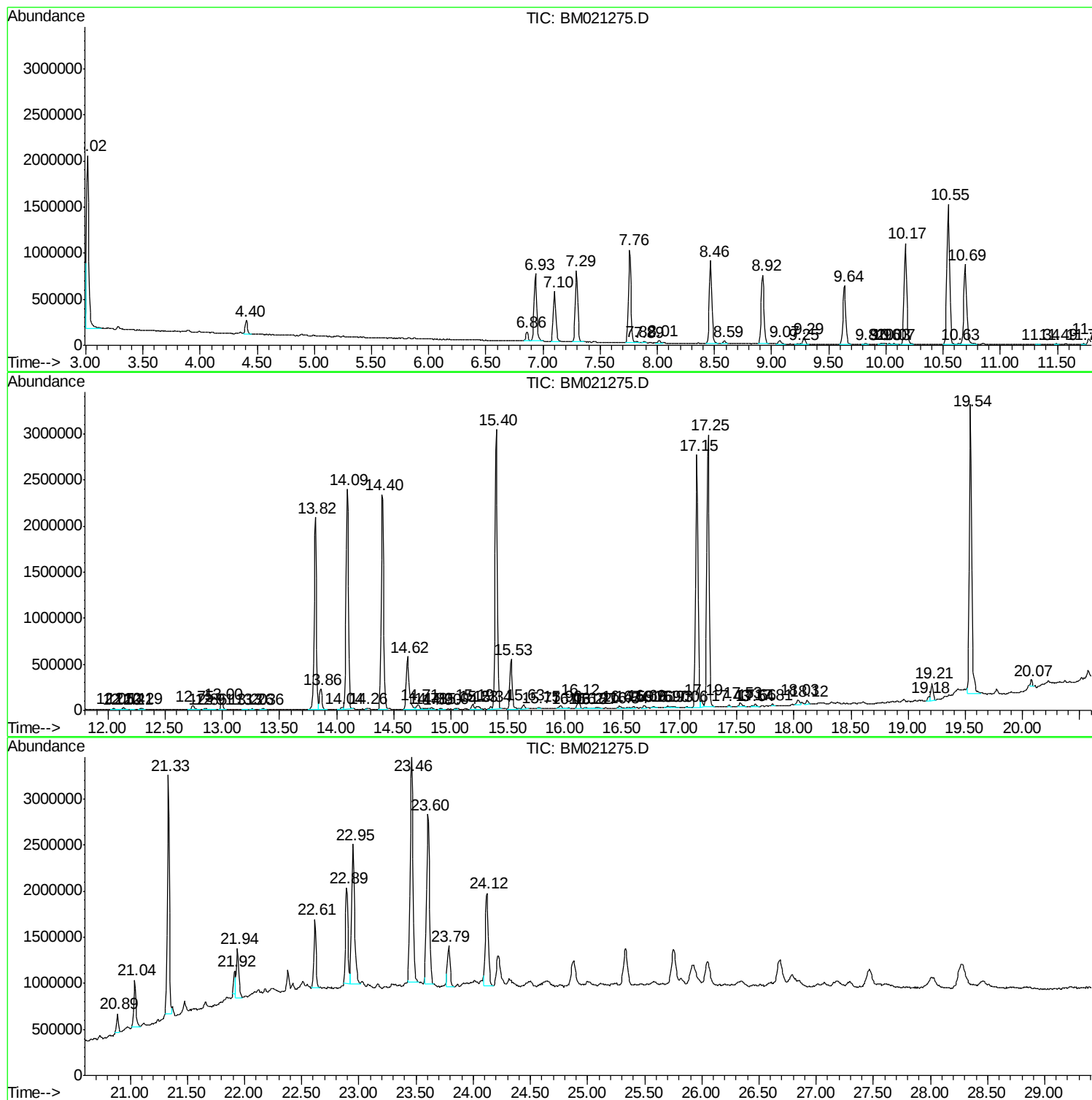
Sum of corrected areas: 69742715

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Instrument :
BNA_M
ClientSampled :
C5BC5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_M
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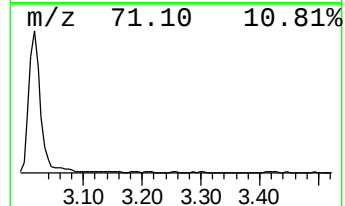
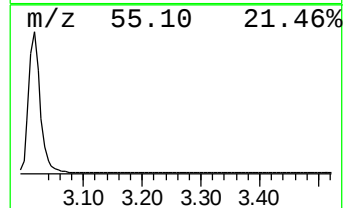
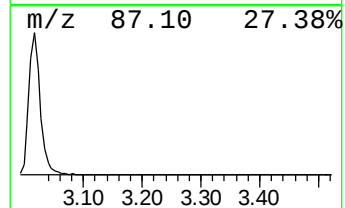
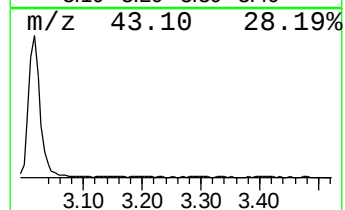
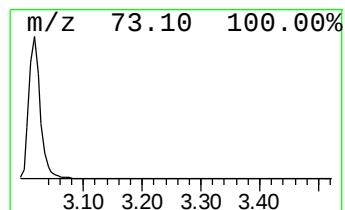
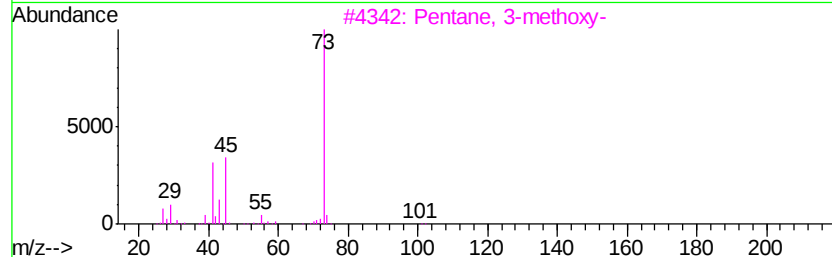
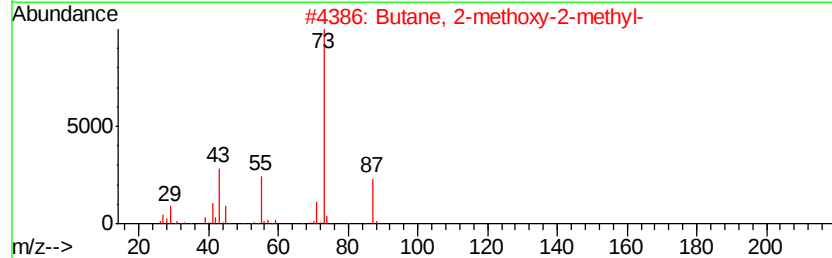
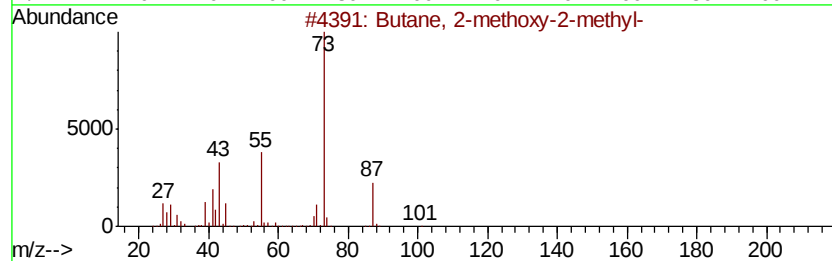
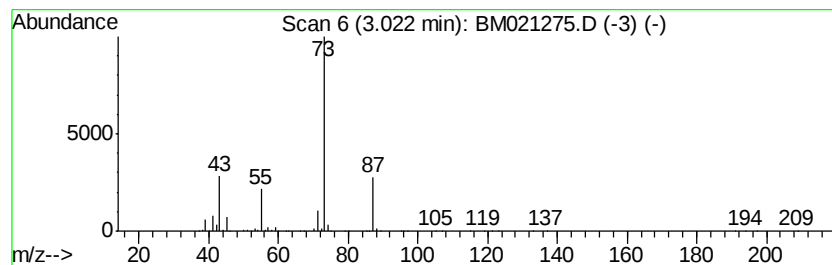
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	28.36 ng/ul	2240420	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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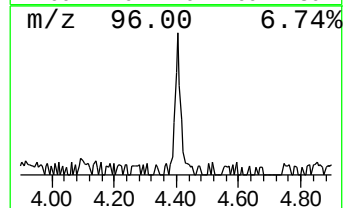
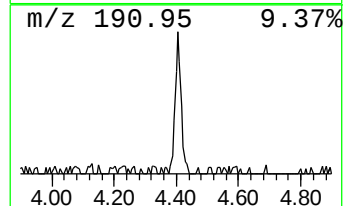
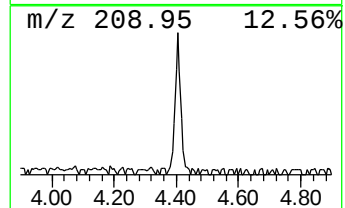
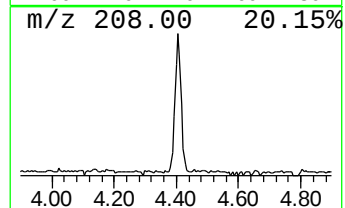
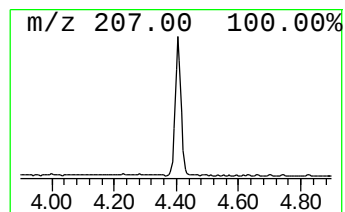
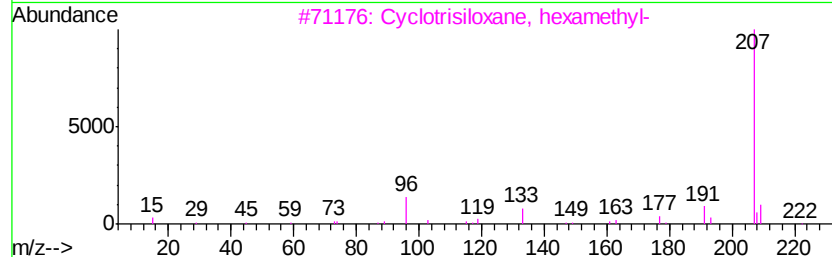
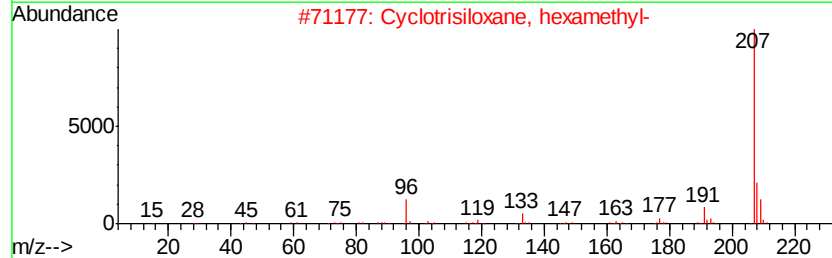
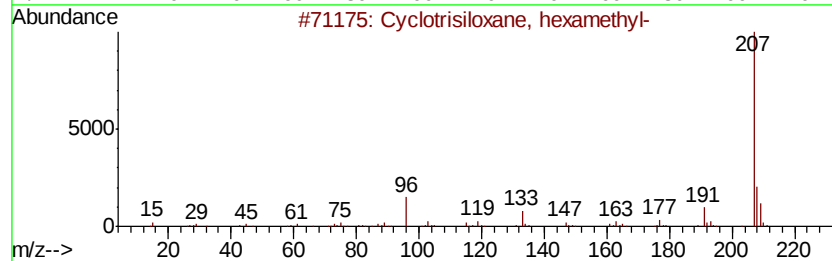
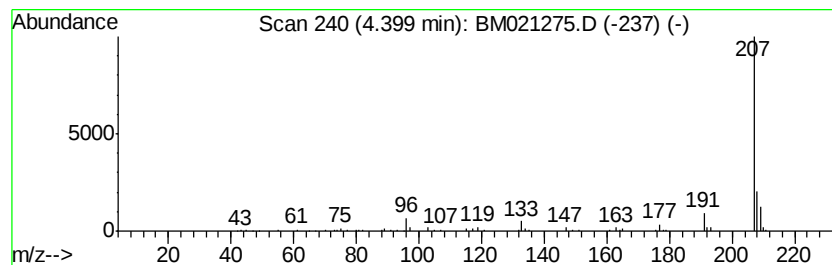
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclotrisiloxane, hexamethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	2.46 ng/ul	194386	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	90
2		Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	86
3		Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	64
4		5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	39
5		1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	39



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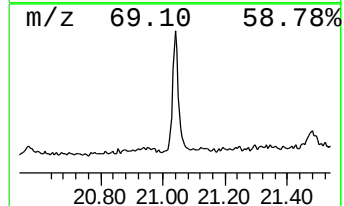
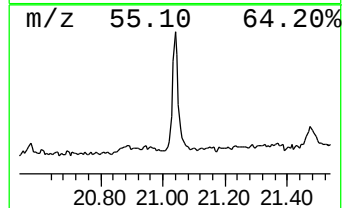
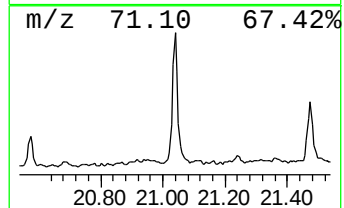
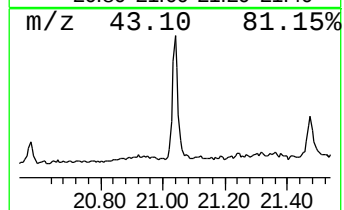
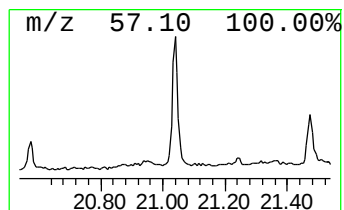
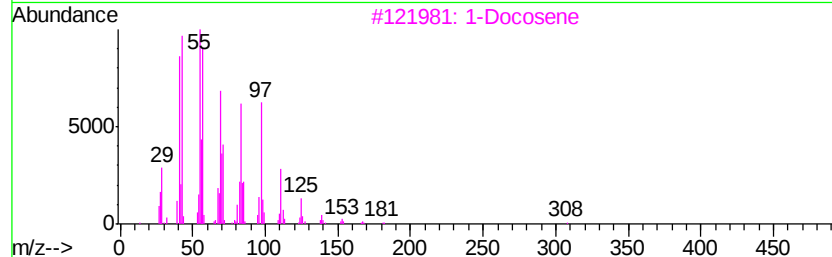
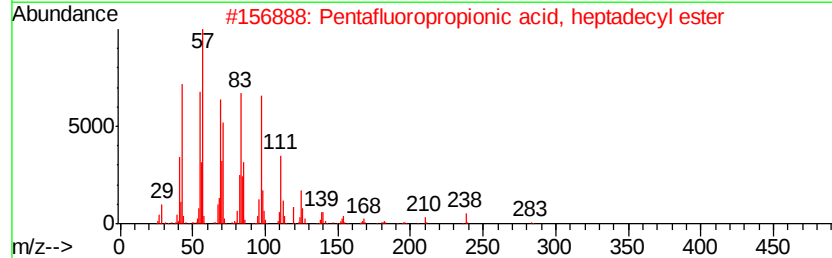
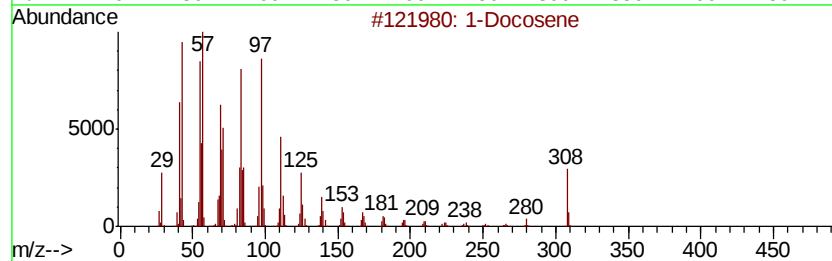
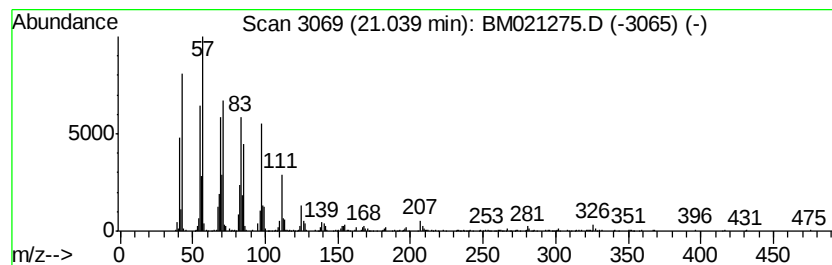
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic21.04 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	4.19 ng/ul	687027	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	96
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87
3		1-Docosene	308	C22H44	001599-67-3	86
4		1-Hentetracontanol	593	C41H84O	040710-42-7	83
5		1-Heptacosanol	396	C27H56O	002004-39-9	74



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Data File : BM021275.D
Acq On : 29 Jun 2019 23:05
Operator : HP/JU
Sample : K3593-05
Misc :
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Instrument :
BNA_M
ClientSampleId :
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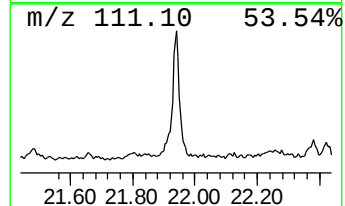
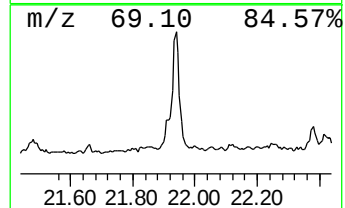
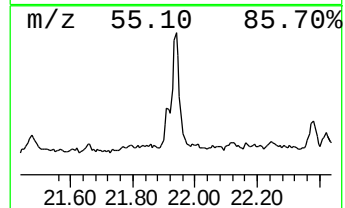
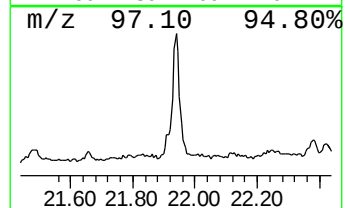
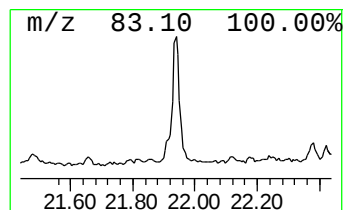
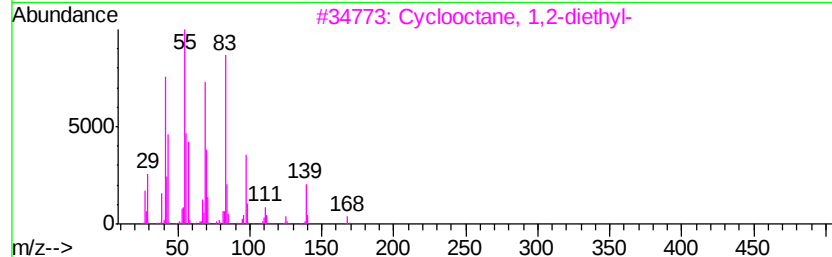
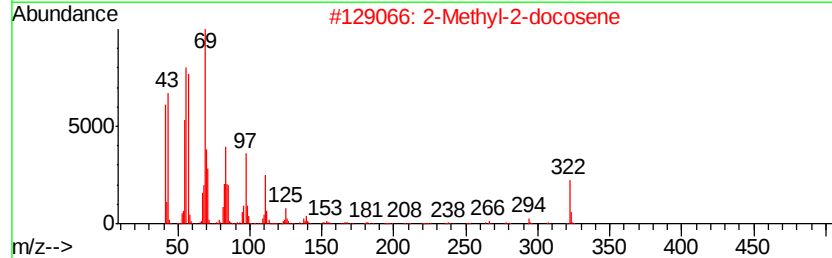
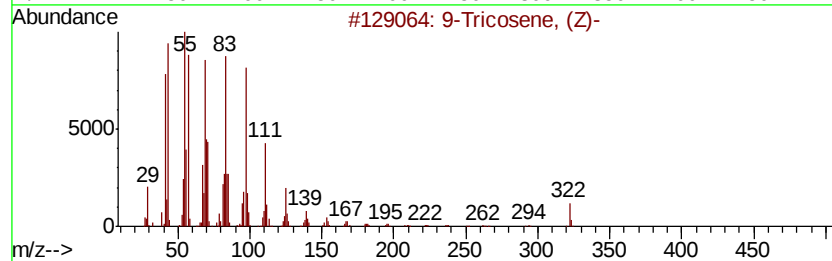
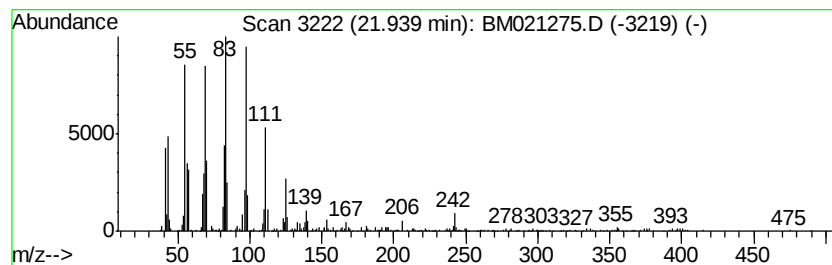
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic21.94 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.94	4.74 ng/ul	777267	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	90
2		2-Methyl-2-docosene	322	C23H46	1000131-16-9	64
3		Cyclooctane, 1,2-diethyl-	168	C12H24	023609-46-3	60
4		Cyclohexane, 1,1',1''-(1-propanyl-...	290	C21H38	055682-89-8	58
5		9-Tricosene, (Z)-	322	C23H46	027519-02-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021275.D
Acq On : 29 Jun 2019 23:05
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Sample : K3593-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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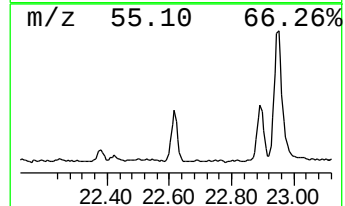
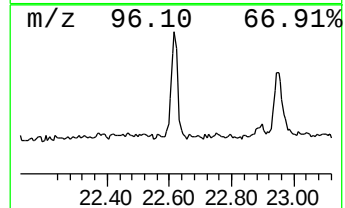
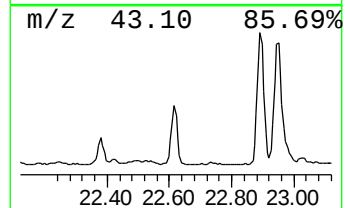
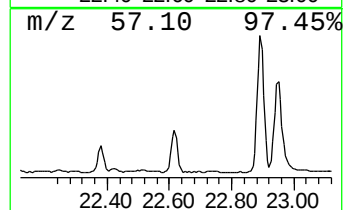
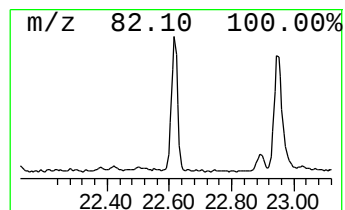
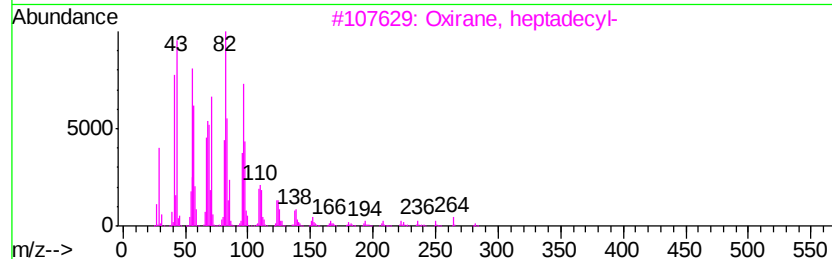
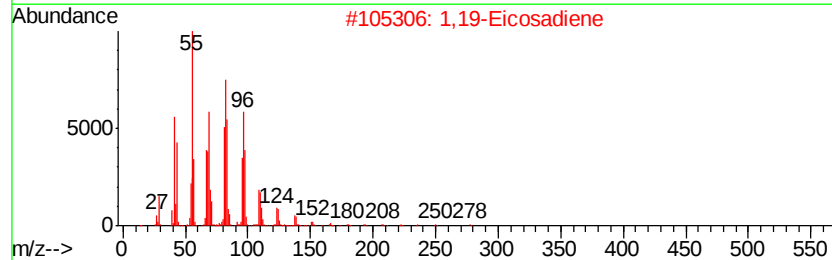
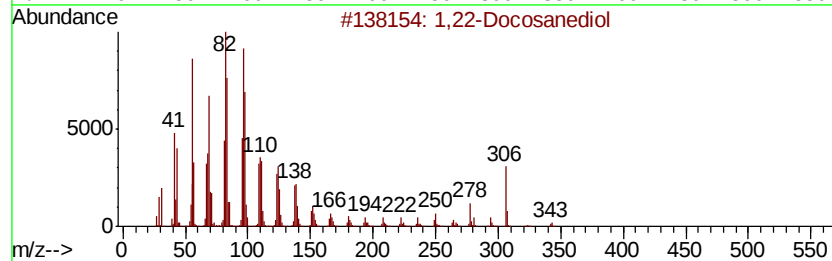
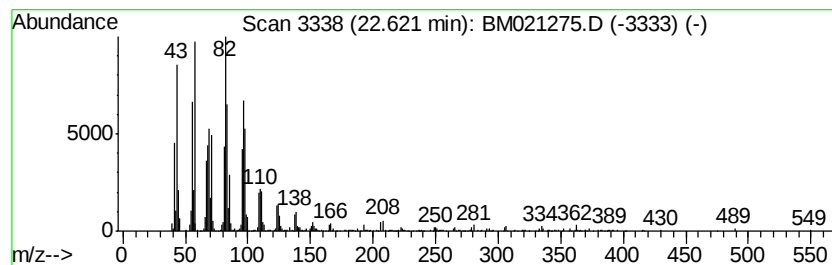
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,22-Docosanediol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.62	5.88 ng/ul	1001980	Perylene-d12	23.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,22-Docosanediol	342	C22H46O2	022513-81-1	93
2		1,19-Eicosadiene	278	C20H38	014811-95-1	93
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4		Octadecanal	268	C18H36O	000638-66-4	91
5		17-Octadecenal	266	C18H34O	056554-86-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021275.D
Acq On : 29 Jun 2019 23:05
Operator : HP/JU
Sample : K3593-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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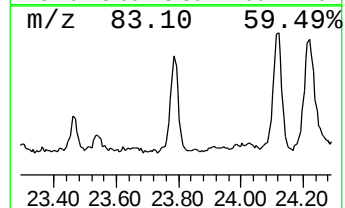
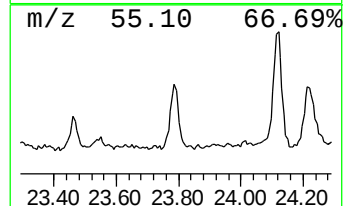
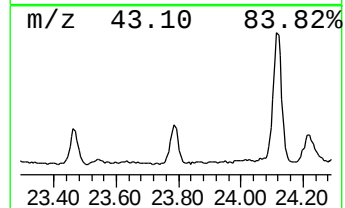
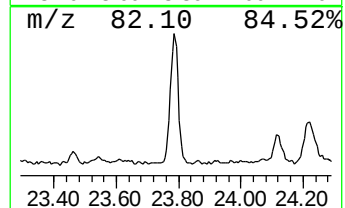
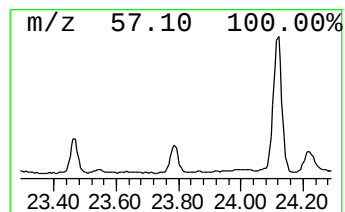
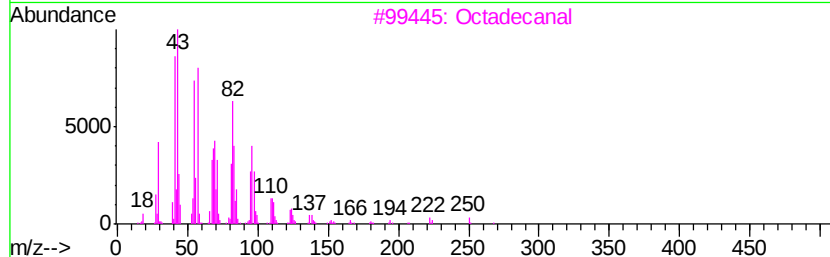
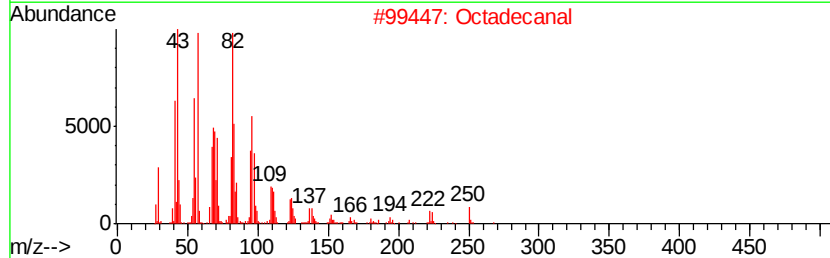
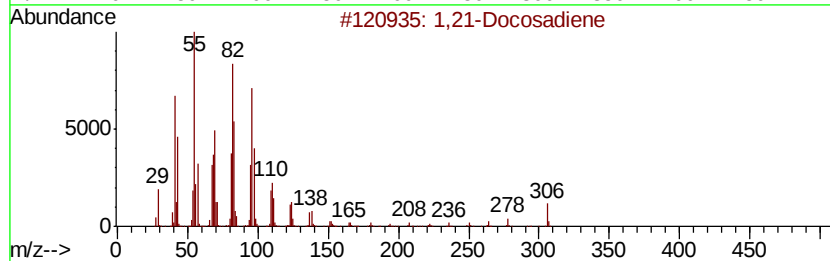
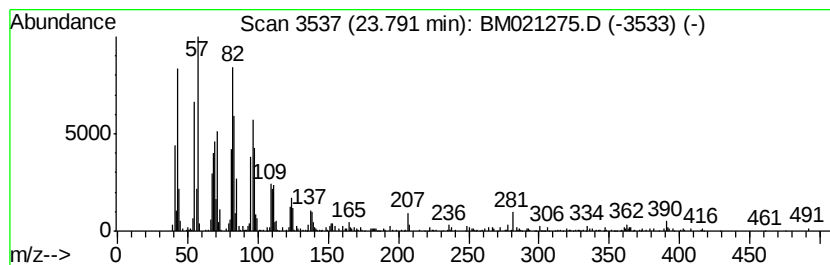
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,21-Docosadiene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.79	4.26 ng/ul	725822	Perylene-d12	23.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,21-Docosadiene	306	C22H42	053057-53-7	99
2		Octadecanal	268	C18H36O	000638-66-4	95
3		Octadecanal	268	C18H36O	000638-66-4	94
4		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
5		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\
Data File : BM021275.D
Acq On : 29 Jun 2019 23:05
Operator : HP/JU
Sample : K3593-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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ClientSampleId :
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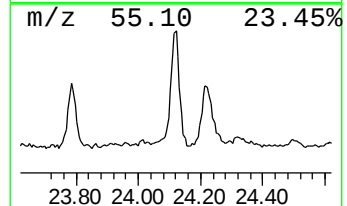
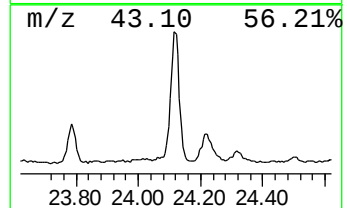
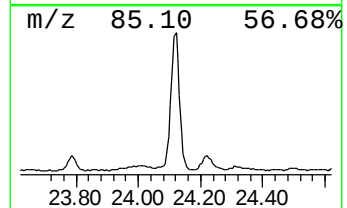
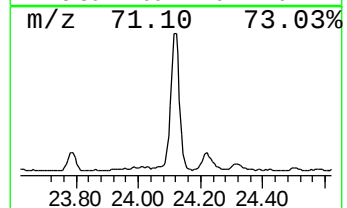
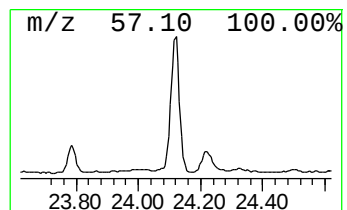
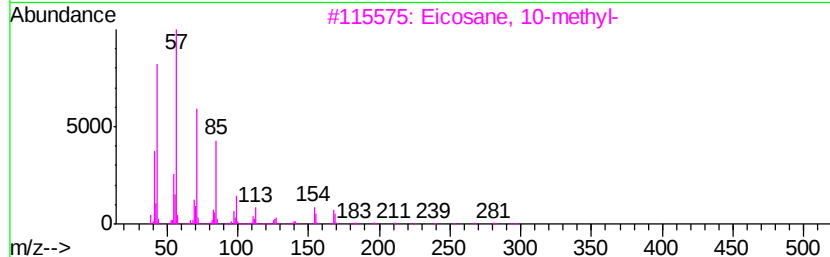
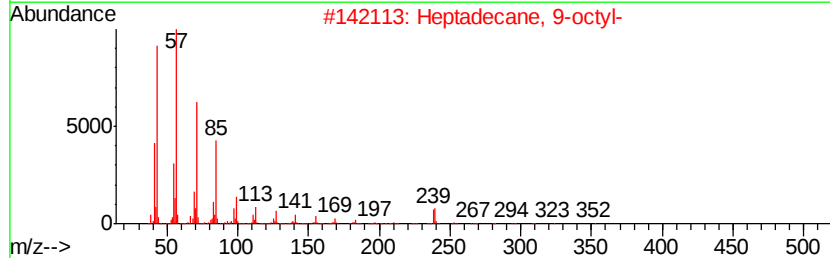
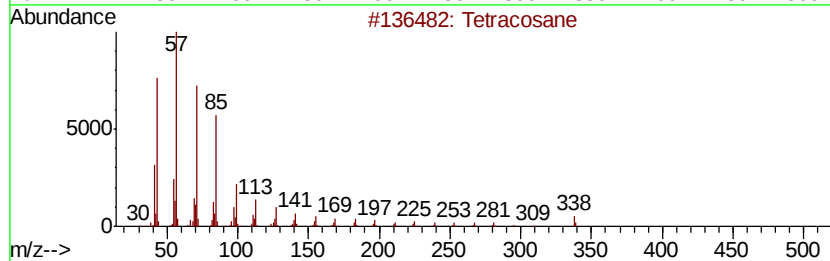
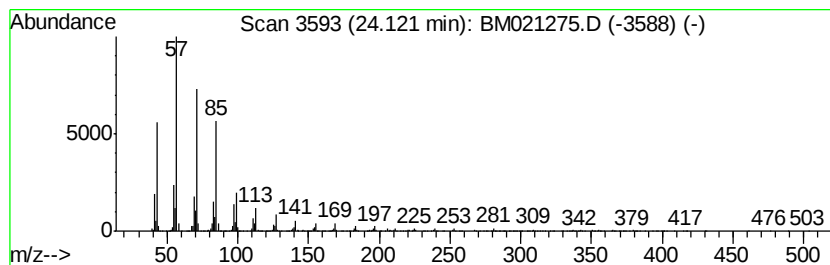
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.12	11.13 ng/ul	1898600	Perylene-d12	23.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062919\
Data File : BM021275.D
Acq On : 29 Jun 2019 23:05
Operator : HP/JU
Sample : K3593-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5BC5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.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
Butane, 2-methoxy...	3.02	28.4	ng/ul	2240420	1	7.76	1579870	20.0
Cyclotrisiloxane,...	4.40	2.5	ng/ul	194386	1	7.76	1579870	20.0
(DEL) Alkane: Cyc...	21.04	4.2	ng/ul	687027	5	21.33	3280320	20.0
(DEL) Alkane: Cyc...	21.94	4.7	ng/ul	777267	5	21.33	3280320	20.0
1,22-Docosanediol	22.62	5.9	ng/ul	1001980	6	23.60	3410950	20.0
1,21-Docosadiene	23.79	4.3	ng/ul	725822	6	23.60	3410950	20.0
(DEL) Alkane: Str...	24.12	11.1	ng/ul	1898600	6	23.60	3410950	20.0