

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
 Data File : BP001915.D
 Acq On : 26 Feb 2020 01:35
 Operator : CG/JU
 Sample : L1543-11
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBD93

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_P\METHODS\SOM-EPA-BP021820MA.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	2.922	3	6	30	rVB	2694141	4029664	36.63%	3.076%
2	3.169	44	48	62	rVB	70093	106147	0.96%	0.081%
3	3.810	153	157	165	rVV	23174	35402	0.32%	0.027%
4	3.899	169	172	179	rVB	14761	22013	0.20%	0.017%
5	4.440	260	264	271	rVB4	8097	13655	0.12%	0.010%
6	4.799	320	325	333	rBV2	88879	141243	1.28%	0.108%
7	6.828	661	670	686	rVB	1341188	2184315	19.85%	1.667%
8	6.987	691	697	710	rVB	1052419	1674156	15.22%	1.278%
9	7.187	721	731	743	rBV	1428865	2297230	20.88%	1.754%
10	7.304	745	751	759	rVB	6353	16663	0.15%	0.013%
11	7.651	802	810	819	rBV	1274144	2022218	18.38%	1.544%
12	8.187	897	901	908	rVB3	9008	17282	0.16%	0.013%
13	8.357	922	930	942	rBV	1721575	2812708	25.57%	2.147%
14	8.792	996	1004	1013	rBV	1269517	2092932	19.02%	1.598%
15	9.281	1081	1087	1094	rVB	46623	76260	0.69%	0.058%
16	9.510	1118	1126	1134	rBV	1061565	1739564	15.81%	1.328%
17	10.045	1209	1217	1233	rBV	1839039	3067913	27.89%	2.342%
18	10.422	1272	1281	1293	rBV	1868268	3065088	27.86%	2.340%
19	10.551	1296	1303	1320	rBV	1519450	2645231	24.04%	2.019%
20	10.951	1365	1371	1376	rBV7	9812	19108	0.17%	0.015%
21	11.957	1532	1542	1548	rBV3	25896	56651	0.51%	0.043%
22	12.016	1548	1552	1561	rVB2	32714	59382	0.54%	0.045%
23	12.651	1654	1660	1665	rBV7	7245	14967	0.14%	0.011%
24	13.033	1721	1725	1729	rBV4	17926	26205	0.24%	0.020%
25	13.086	1729	1734	1739	rVB4	11108	18214	0.17%	0.014%
26	13.204	1746	1754	1760	rBV2	40643	70767	0.64%	0.054%
27	13.610	1820	1823	1828	rVB6	7505	11258	0.10%	0.009%
28	13.698	1831	1838	1843	rBV	3369051	4646884	42.24%	3.547%
29	13.745	1843	1846	1855	rVB	590455	779964	7.09%	0.595%
30	13.975	1877	1885	1899	rBV	3900071	5813945	52.85%	4.438%
31	14.286	1931	1938	1946	rVV	2944386	4319832	39.27%	3.297%
32	14.374	1951	1953	1958	rVB6	10354	12321	0.11%	0.009%
33	14.480	1962	1971	1986	rBV	1378322	2087594	18.98%	1.594%
34	14.651	1995	2000	2005	rVV8	6821	12766	0.12%	0.010%

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
 Data File : BP001915.D
 Acq On : 26 Feb 2020 01:35
 Operator : CG/JU
 Sample : L1543-11
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBD93

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_P\METHODS\SOM-EPA-BP021820MA.M
 Title : SVOA CALIBRATION

35	14.710	2005	2010	2021	rVB2	72476	116360	1.06%	0.089%
36	15.121	2076	2080	2084	rVB3	12532	17389	0.16%	0.013%
37	15.180	2084	2090	2095	rBV7	10939	17124	0.16%	0.013%
38	15.286	2100	2108	2116	rBV	4947480	7397178	67.24%	5.646%
39	15.398	2121	2127	2144	rVV	1179400	1645159	14.95%	1.256%
40	15.527	2144	2149	2154	rVB2	55662	86740	0.79%	0.066%
41	15.739	2181	2185	2188	rBV5	8072	11041	0.10%	0.008%
42	15.851	2200	2204	2214	rBV5	6939	18147	0.16%	0.014%
43	16.069	2238	2241	2252	rVB6	25899	51333	0.47%	0.039%
44	16.304	2277	2281	2287	rBV5	7531	12438	0.11%	0.009%
45	16.721	2349	2352	2356	rVB3	13872	18548	0.17%	0.014%
46	16.833	2365	2371	2374	rBV2	17232	37211	0.34%	0.028%
47	17.039	2399	2406	2412	rBV	3822532	5445132	49.49%	4.156%
48	17.133	2416	2422	2435	rVB2	5401884	8229371	74.80%	6.282%
49	17.439	2466	2474	2482	rBV2	26670	76091	0.69%	0.058%
50	17.580	2494	2498	2501	rVB4	12122	14297	0.13%	0.011%
51	17.910	2551	2554	2558	rVB4	13272	15388	0.14%	0.012%
52	17.963	2558	2563	2579	rBV	531333	948730	8.62%	0.724%
53	18.251	2608	2612	2617	rBV2	56218	80703	0.73%	0.062%
54	18.433	2640	2643	2648	rVB	22926	25653	0.23%	0.020%
55	18.798	2703	2705	2711	rVB6	28099	39458	0.36%	0.030%
56	18.862	2713	2716	2720	rBV2	34126	41609	0.38%	0.032%
57	18.933	2724	2728	2737	rVB	34388	59742	0.54%	0.046%
58	19.027	2739	2744	2745	rBV	80258	103183	0.94%	0.079%
59	19.057	2745	2749	2755	rVB	323157	417834	3.80%	0.319%
60	19.198	2769	2773	2780	rBV2	105927	172963	1.57%	0.132%
61	19.268	2781	2785	2792	rVB2	89281	141885	1.29%	0.108%
62	19.380	2798	2804	2815	rBV	7528536	11001664	100.00%	8.398%
63	19.457	2815	2817	2827	rVB3	61852	88490	0.80%	0.068%
64	19.639	2845	2848	2854	rVB	56147	78528	0.71%	0.060%
65	19.733	2860	2864	2869	rBV2	33666	62726	0.57%	0.048%
66	19.857	2882	2885	2888	rBV4	41671	68337	0.62%	0.052%
67	19.945	2897	2900	2904	rVB2	76056	81907	0.74%	0.063%
68	19.986	2905	2907	2911	rVB2	35773	40776	0.37%	0.031%
69	20.180	2937	2940	2944	rBV2	49728	57586	0.52%	0.044%
70	20.427	2979	2982	2985	rVB	138358	131098	1.19%	0.100%
71	20.621	3012	3015	3020	rBV3	47605	89792	0.82%	0.069%

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

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_P\METHODS\SOM-EPA-BP021820MA.M
Title : SVOA CALIBRATION

72	20.868	3051	3057	3077	rVB3	1442154	2311469	21.01%	1.764%
73	21.009	3078	3081	3087	rBV2	98471	134678	1.22%	0.103%
74	21.127	3095	3101	3105	rBV	5458082	7229867	65.72%	5.519%
75	21.162	3105	3107	3115	rVB	442987	536260	4.87%	0.409%
76	21.245	3117	3121	3125	rBV	193983	279022	2.54%	0.213%
77	21.304	3128	3131	3137	rVB	372635	436153	3.96%	0.333%
78	21.739	3200	3205	3221	rVB2	585507	1072647	9.75%	0.819%
79	22.198	3279	3283	3289	rVB	548154	741336	6.74%	0.566%
80	22.327	3302	3305	3310	rVB2	98925	131051	1.19%	0.100%
81	22.533	3336	3340	3346	rVB	144118	218987	1.99%	0.167%
82	22.715	3359	3371	3388	rBV2	5034510	10227502	92.96%	7.807%
83	23.045	3422	3427	3434	rBV	294431	631521	5.74%	0.482%
84	23.151	3441	3445	3447	rBV	176471	282995	2.57%	0.216%
85	23.198	3447	3453	3462	rVV	5485486	10308238	93.70%	7.868%
86	23.274	3463	3466	3471	rVV	551250	923323	8.39%	0.705%
87	23.339	3471	3477	3495	rVB2	3487755	6765132	61.49%	5.164%
88	23.592	3514	3520	3530	rBV	922002	1796813	16.33%	1.372%
89	23.927	3571	3577	3584	rVB	526272	1006981	9.15%	0.769%
90	24.003	3586	3590	3601	rVB	180643	382539	3.48%	0.292%
91	24.680	3699	3705	3711	rVB	317022	672505	6.11%	0.513%
92	25.562	3848	3855	3869	rVB3	335026	1026133	9.33%	0.783%
93	26.409	3993	3999	4015	rVB5	294060	940264	8.55%	0.718%

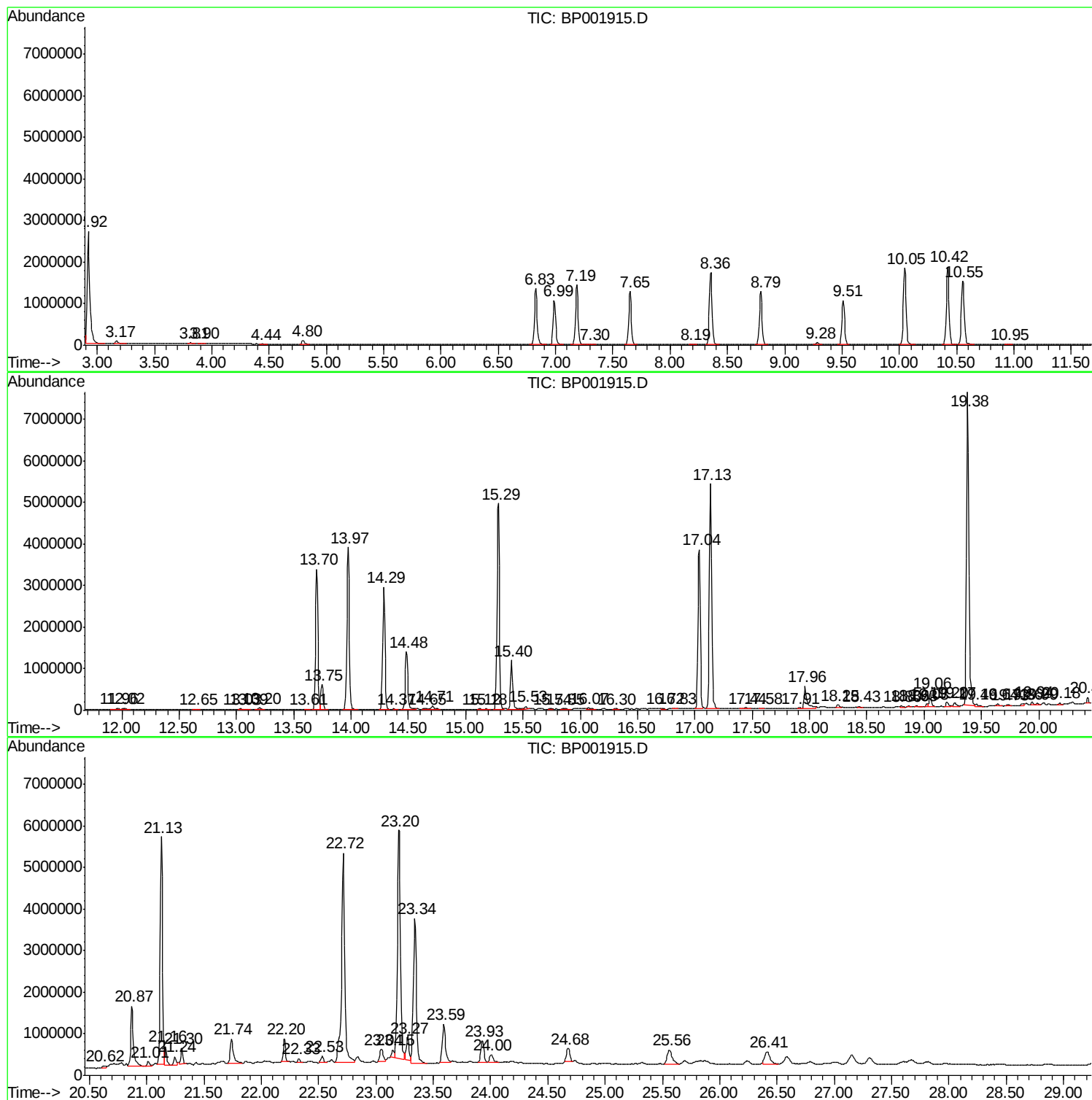
Sum of corrected areas: 131006569

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

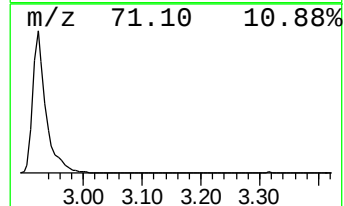
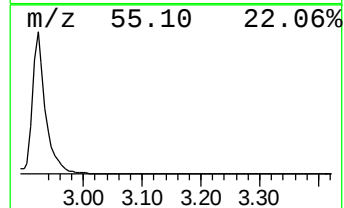
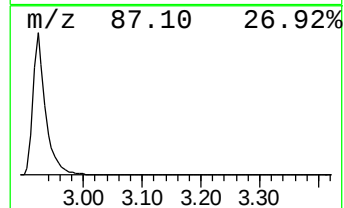
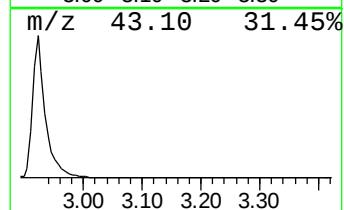
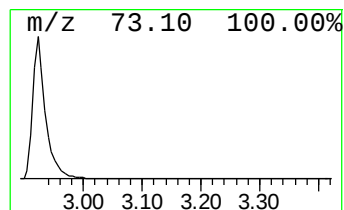
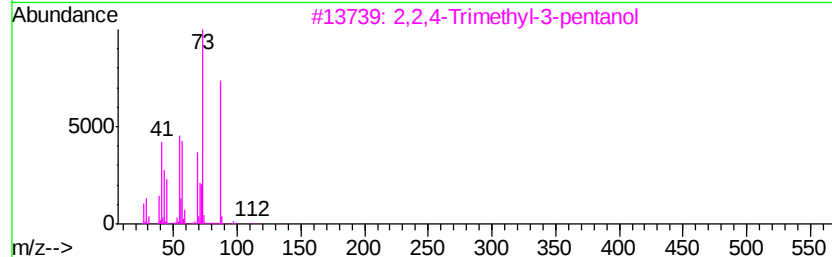
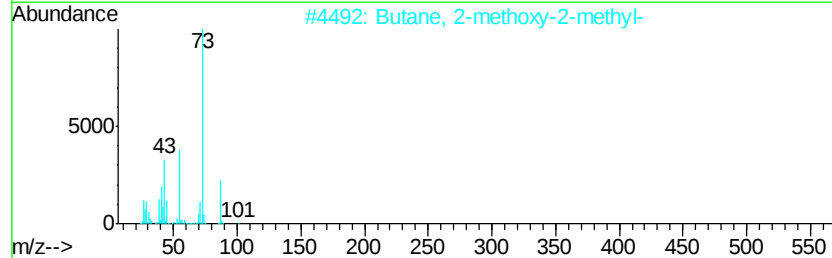
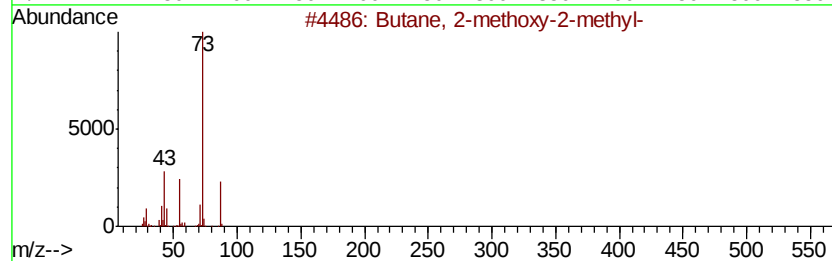
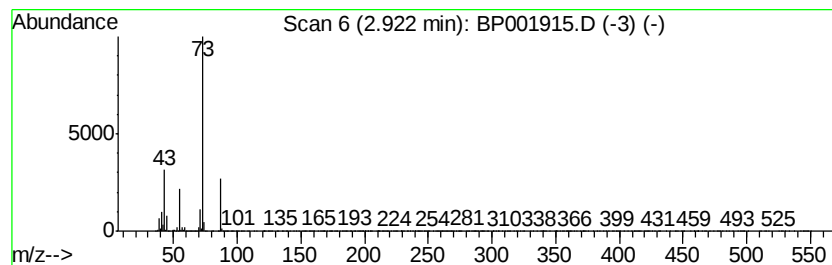
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
2.92	39.85 ng/ul	4029660	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

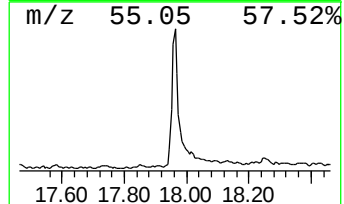
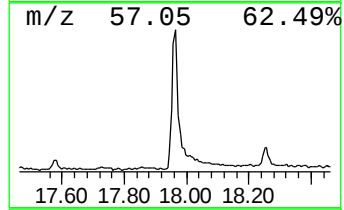
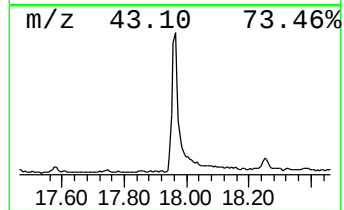
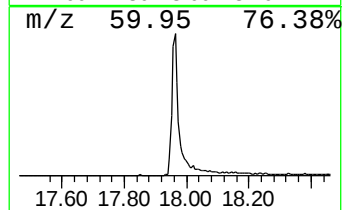
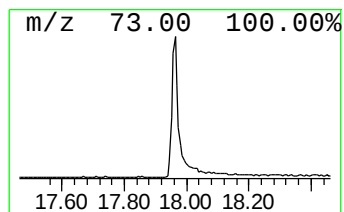
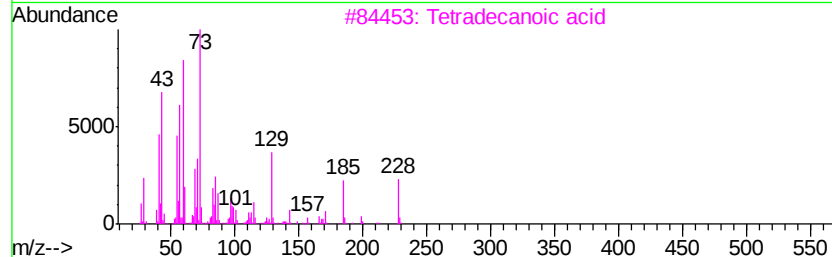
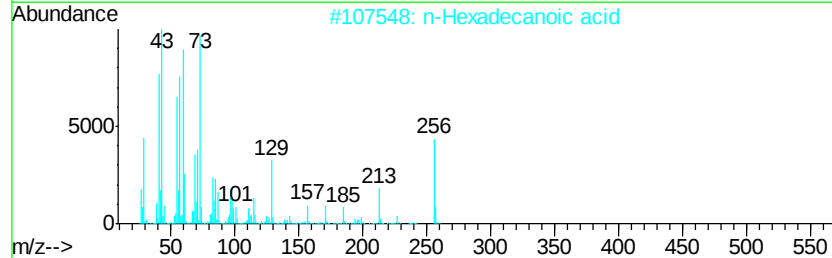
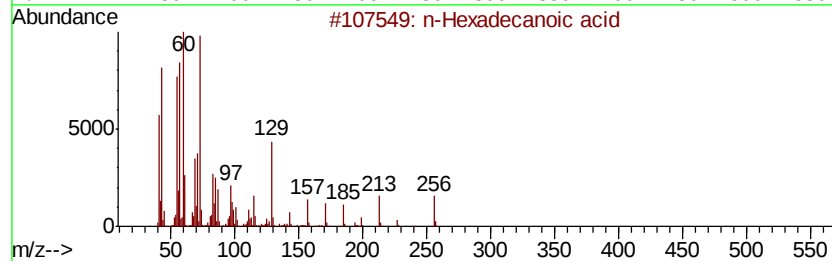
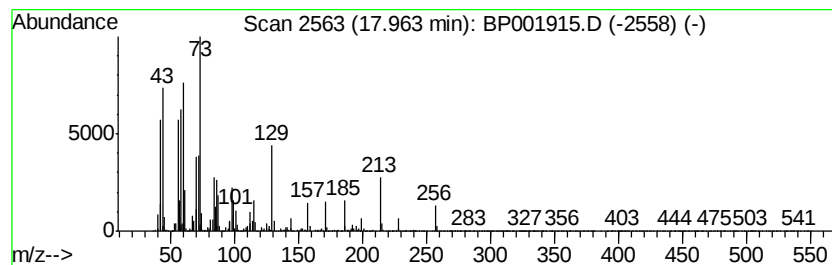
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	3.48 ng/ul	948730	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

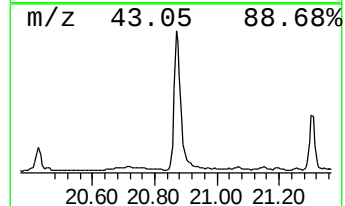
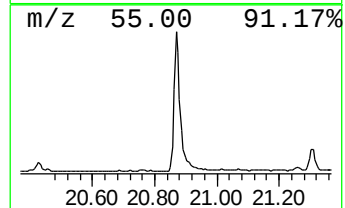
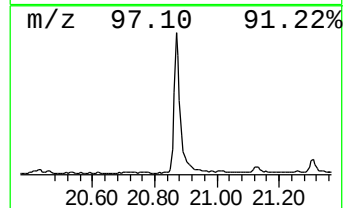
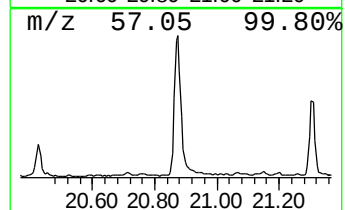
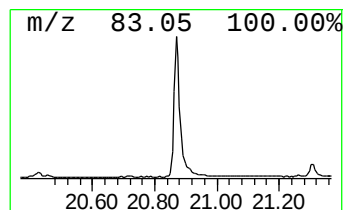
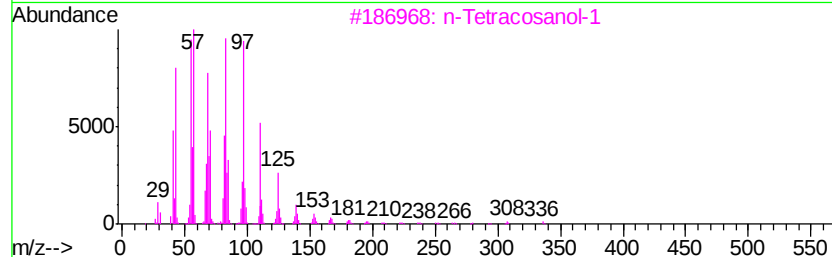
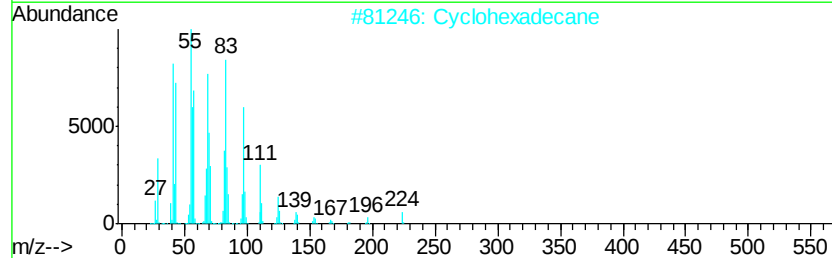
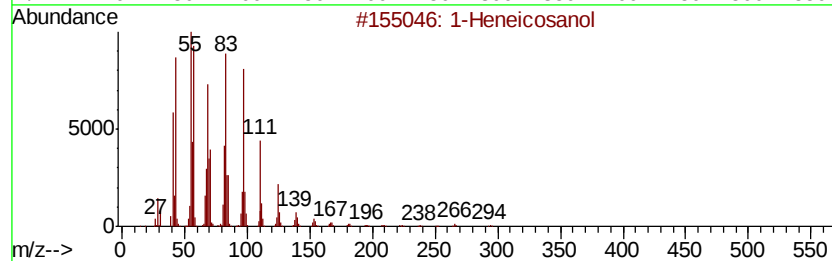
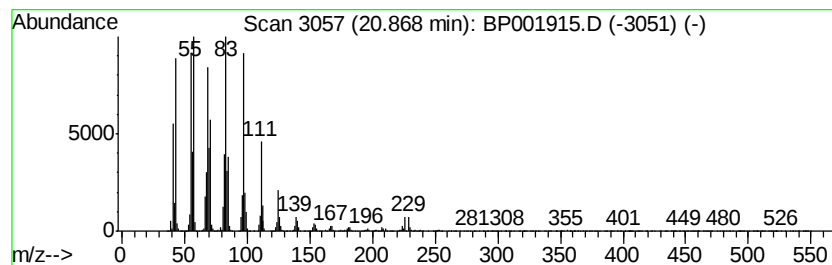
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	6.39 ng/ul	2311470	Chrysene-d12	21.13

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		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

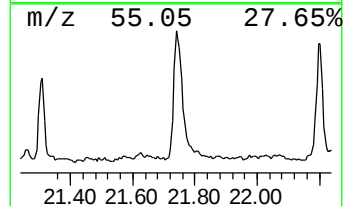
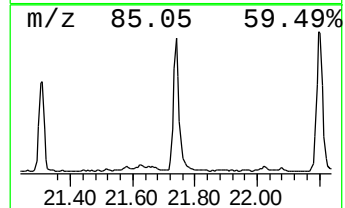
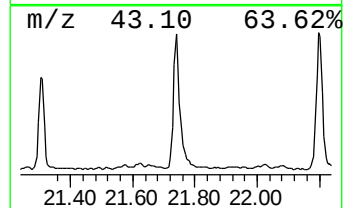
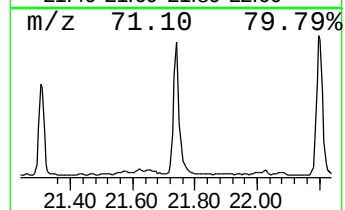
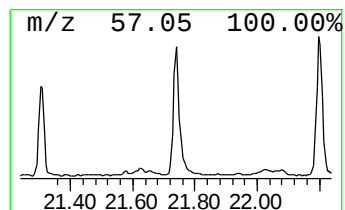
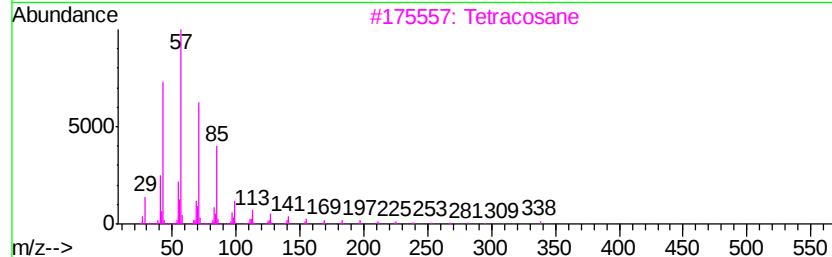
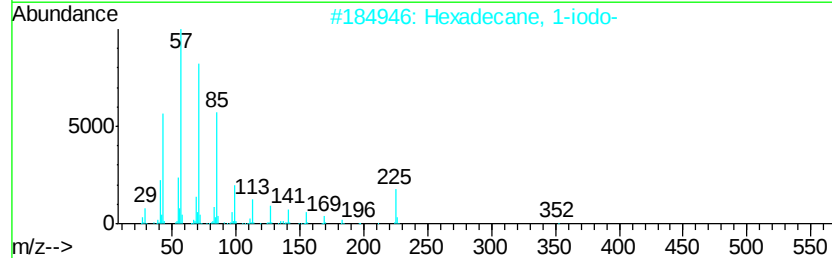
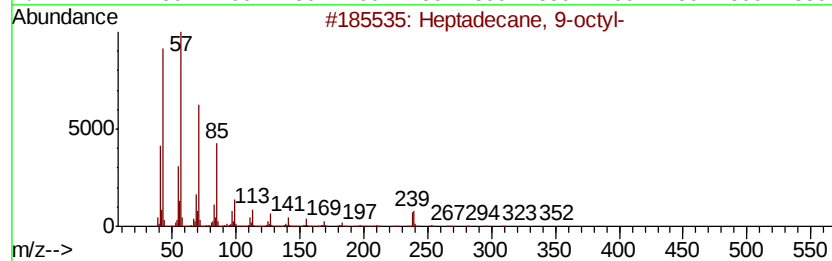
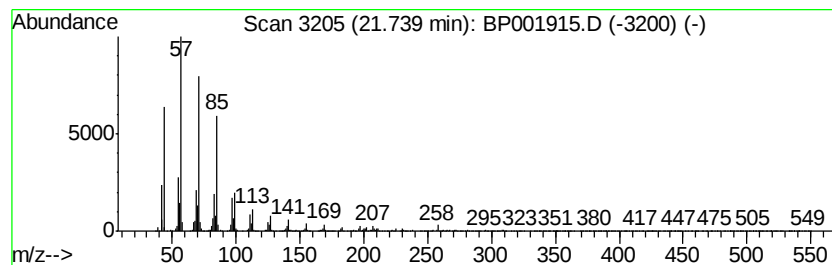
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.74	2.97 ng/ul	1072650	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
3		Tetracosane	338	C24H50	000646-31-1	93
4		Octacosane	394	C28H58	000630-02-4	93
5		Octacosane	394	C28H58	000630-02-4	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

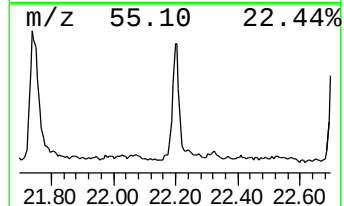
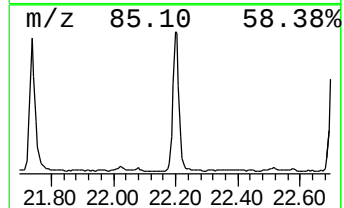
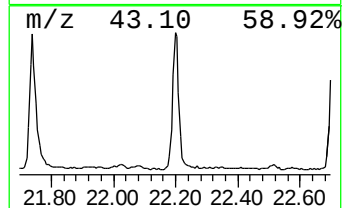
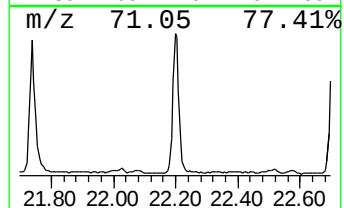
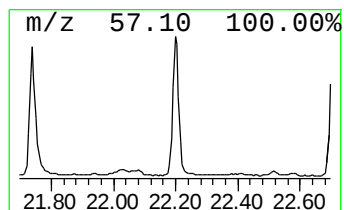
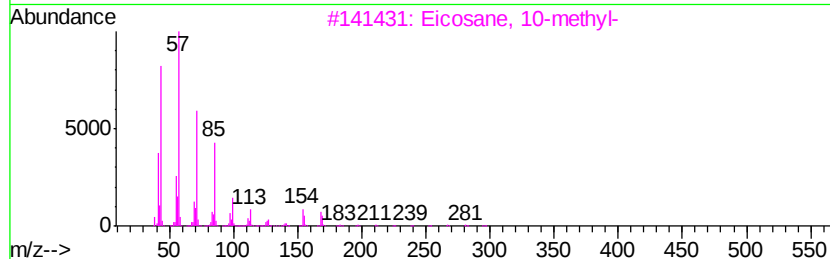
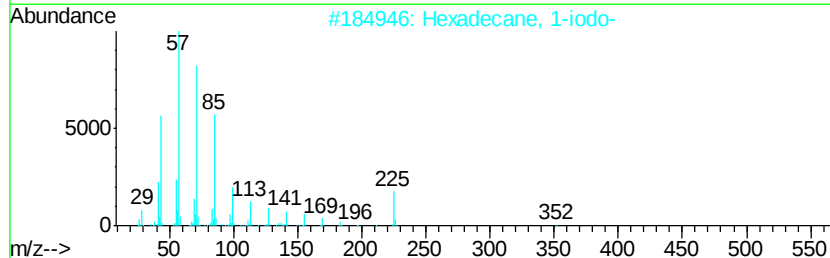
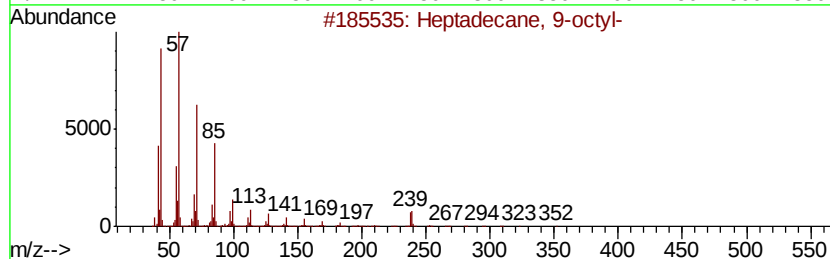
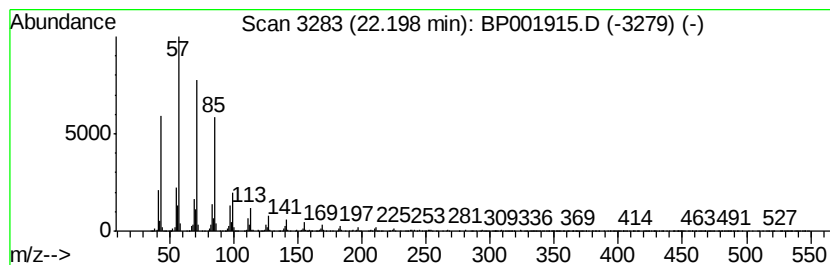
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.20	2.05 ng/ul	741336	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	97
2		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

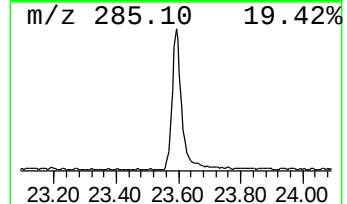
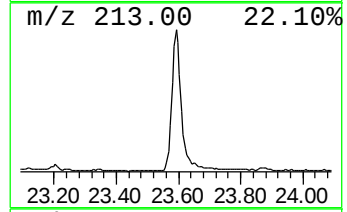
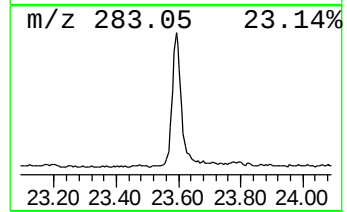
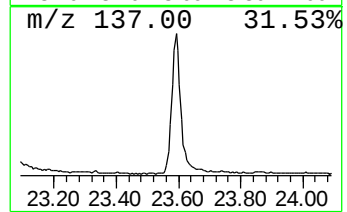
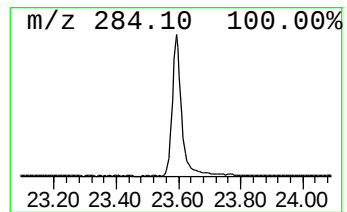
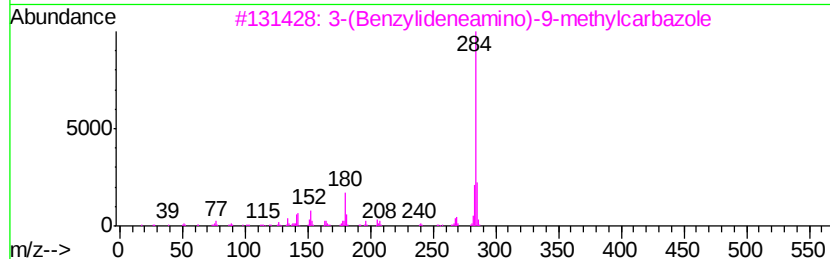
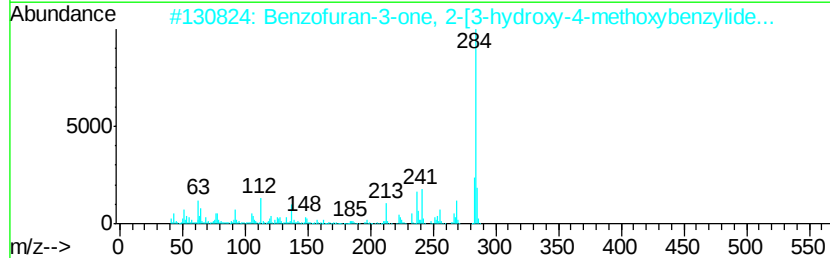
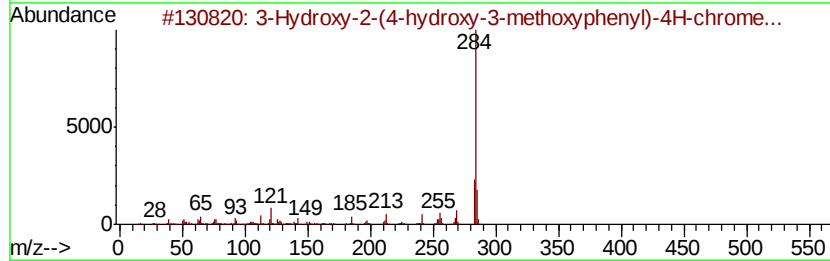
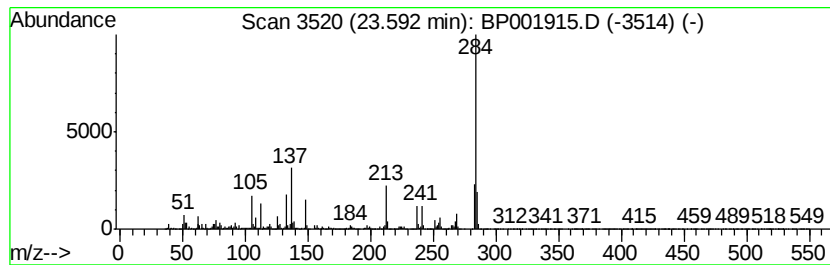
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 3-Hydroxy-2-(4-hydroxy-3-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.59	5.31 ng/ul	1796810	Perylene-d12	23.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Hydroxy-2-(4-hydroxy-3-methoxy...	284	C16H12O5	159184-95-9	92
2		Benzofuran-3-one, 2-[3-hydroxy-4...	284	C16H12O5	1000128-62-8	68
3		3-(Benzylideneamino)-9-methylcar...	284	C20H16N2	035473-72-4	58
4		7-Hydroxy-2-(3-hydroxy-4-methoxy...	284	C16H12O5	020575-57-9	53
5		Iron, (.eta.-5-cyclohepta-1,4-di...	284	C17H24Fe	1000162-47-6	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

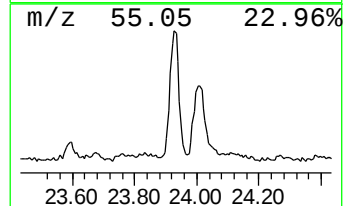
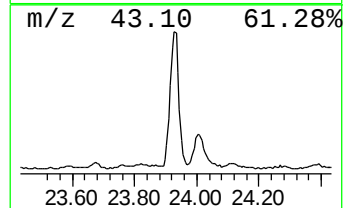
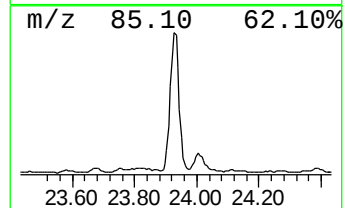
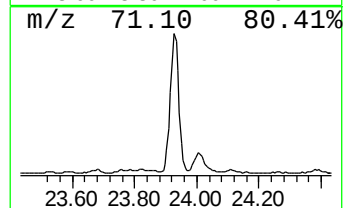
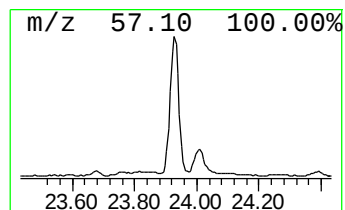
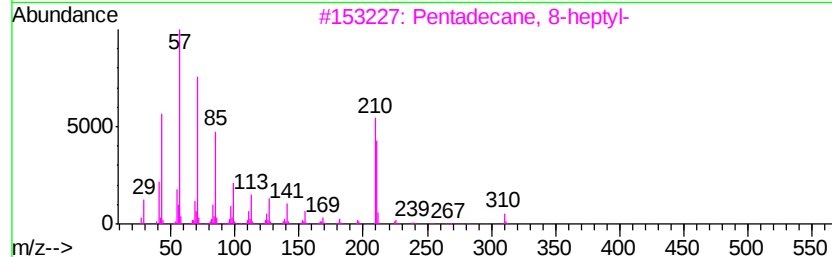
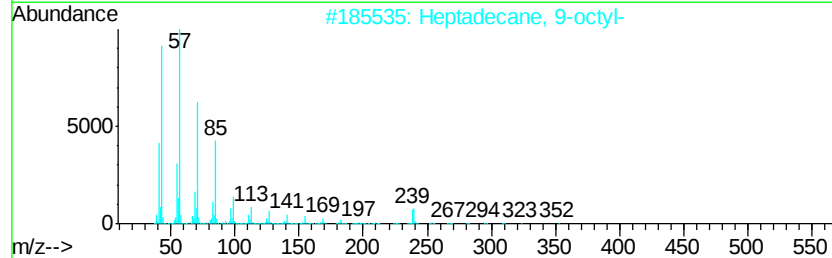
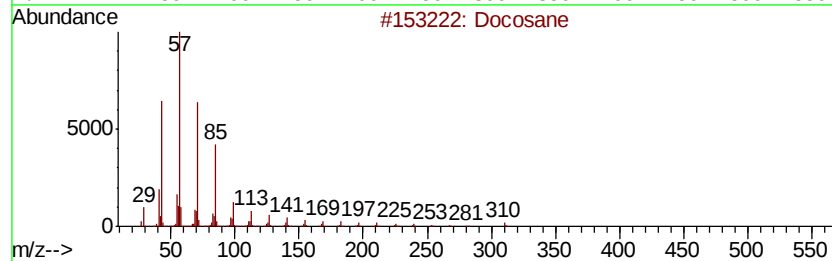
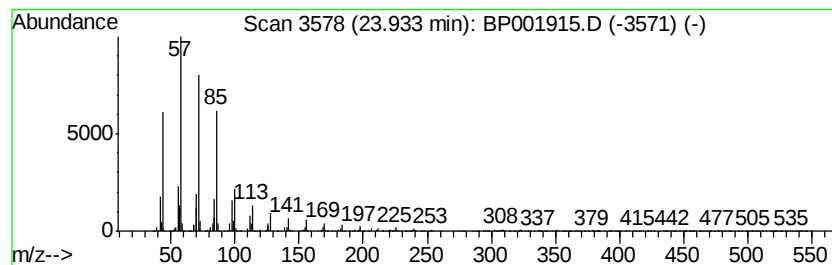
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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.
23.93	2.98 ng/ul	1006980	Perylene-d12	23.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	95
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	93
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

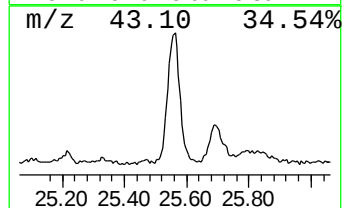
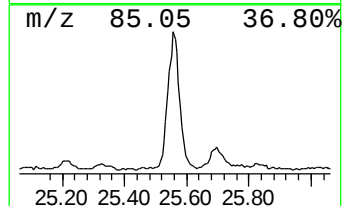
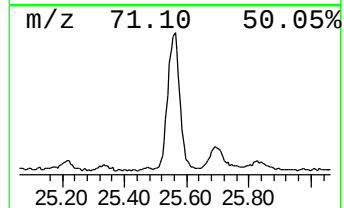
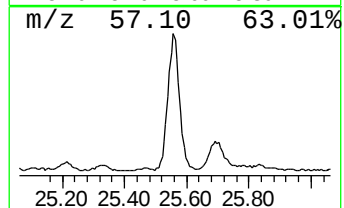
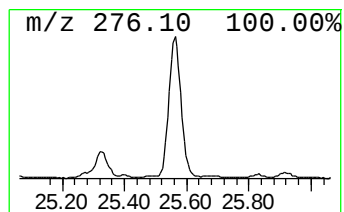
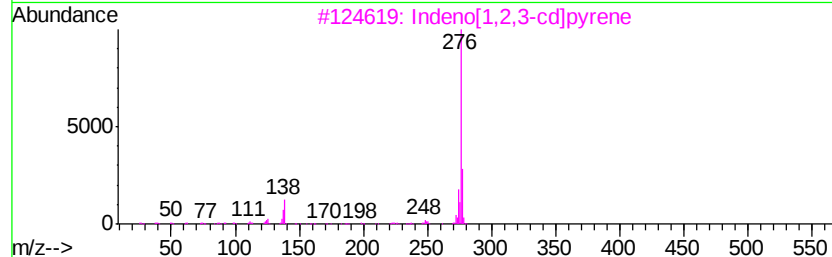
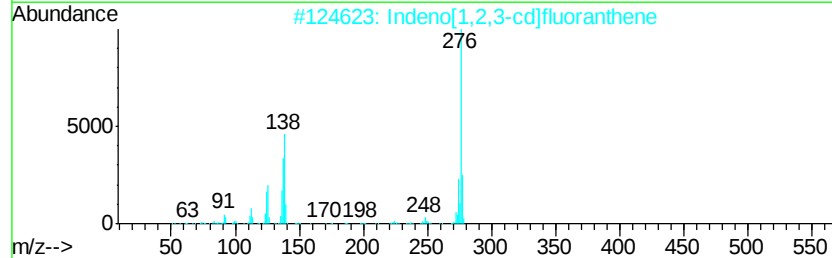
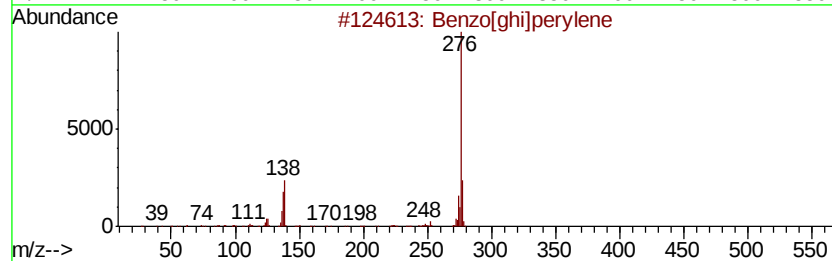
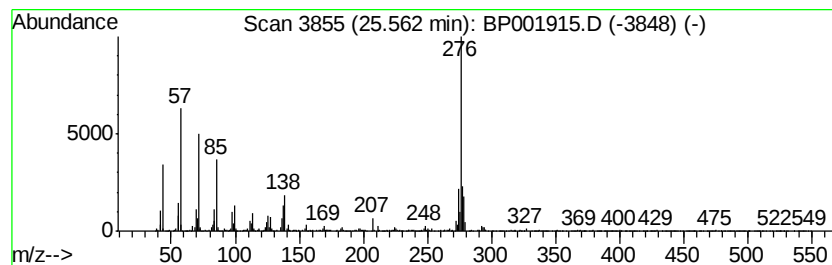
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Indeno[1,2,3-cd]fluoranthene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.56	3.03 ng/ul	1026130	Perylene-d12	23.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[ghi]perylene	276	C22H12	000191-24-2	95
2		Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	91
3		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	91
4		Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	87
5		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

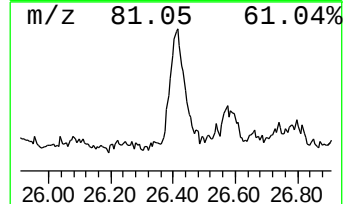
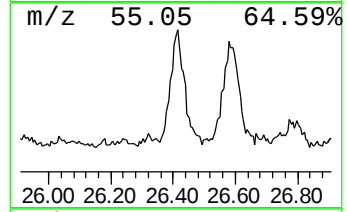
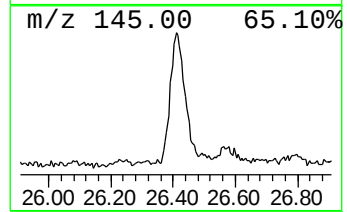
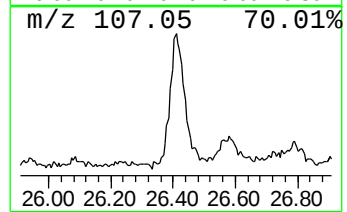
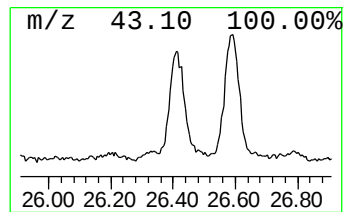
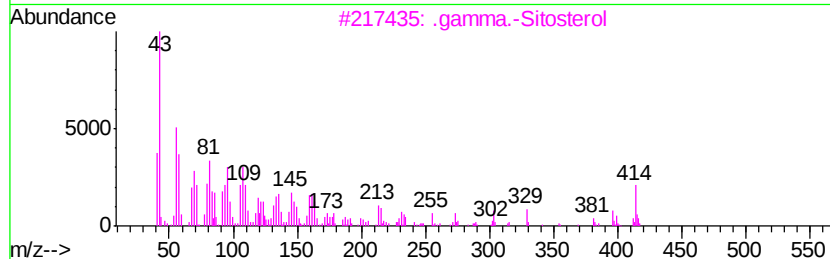
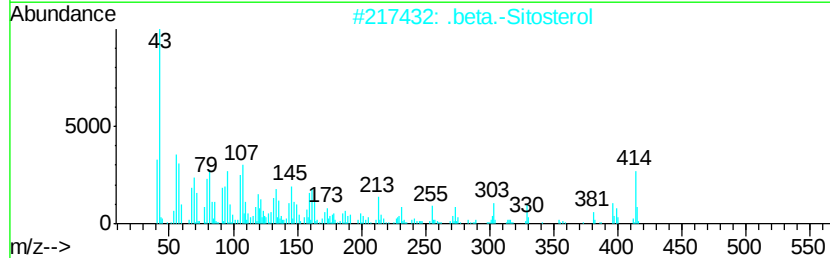
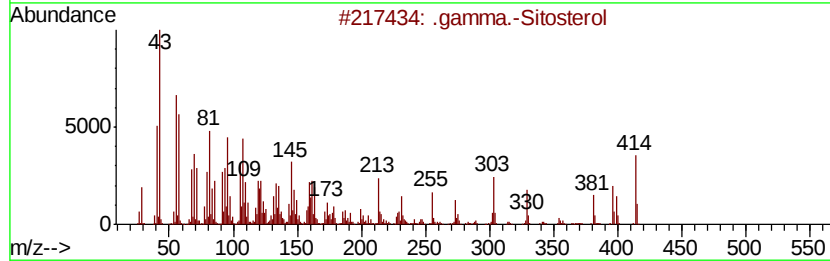
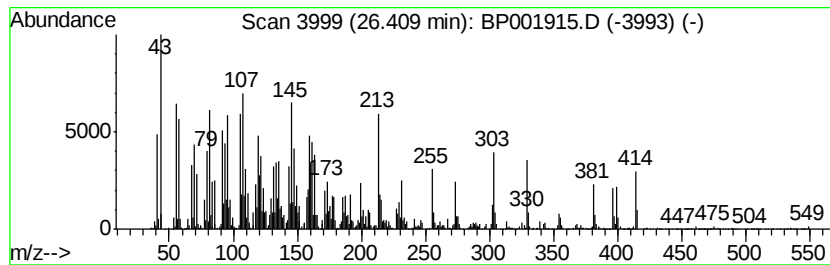
Instrument :
BNA_P
ClientSampleId :
DBD93

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 .gamma.-Sitosterol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.41	2.78 ng/ul	940264	Perylene-d12	23.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.gamma.-Sitosterol	414 C29H50O	000083-47-6 99
2		.beta.-Sitosterol	414 C29H50O	000083-46-5 95
3		.gamma.-Sitosterol	414 C29H50O	000083-47-6 86
4		.beta.-Sitosterol	414 C29H50O	000083-46-5 83
5		5-Cholestene-3-ol, 24-methyl-	400 C28H48O	1000214-17-4 14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
Data File : BP001915.D
Acq On : 26 Feb 2020 01:35
Operator : CG/JU
Sample : L1543-11
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBD93

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.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
Butane, 2-methoxy...	2.92	39.9	ng/ul	4029660	1	7.65	2022220	20.0
n-Hexadecanoic acid	17.96	3.5	ng/ul	948730	4	17.04	5445130	20.0
1-Heneicosanol	20.87	6.4	ng/ul	2311470	5	21.13	7229870	20.0
(DEL) Alkane: Str...	21.74	3.0	ng/ul	1072650	5	21.13	7229870	20.0
(DEL) Alkane: Str...	22.20	2.0	ng/ul	741336	5	21.13	7229870	20.0
3-Hydroxy-2-(4-hy...	23.59	5.3	ng/ul	1796810	6	23.34	6765130	20.0
(DEL) Alkane: Str...	23.93	3.0	ng/ul	1006980	6	23.34	6765130	20.0
Indeno[1,2,3-cd]f...	25.56	3.0	ng/ul	1026130	6	23.34	6765130	20.0
.gamma.-Sitosterol	26.41	2.8	ng/ul	940264	6	23.34	6765130	20.0