

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
 Data File : BP002267.D
 Acq On : 13 May 2020 01:24
 Operator : CG/JU
 Sample : L2570-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFSP3

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-BP051220MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP002269.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	20	rVB	733797	784264	13.50%	1.127%
2	3.158	42	46	58	rVB	86914	115989	2.00%	0.167%
3	5.105	372	377	385	rBV	144434	201715	3.47%	0.290%
4	6.640	631	638	645	rBV2	28235	42314	0.73%	0.061%
5	6.793	658	664	673	rBV	403357	647168	11.14%	0.930%
6	6.958	685	692	702	rBV	1145686	1816884	31.29%	2.610%
7	7.152	718	725	734	rVV	1268205	1968015	33.89%	2.827%
8	7.228	734	738	743	rVV2	24596	44032	0.76%	0.063%
9	7.287	743	748	753	rVV2	57083	95646	1.65%	0.137%
10	7.617	796	804	813	rBV	981195	1573886	27.10%	2.261%
11	7.699	814	818	822	rVV3	20422	38605	0.66%	0.055%
12	7.964	855	863	871	rBV3	23845	46918	0.81%	0.067%
13	8.316	916	923	935	rVV	995348	1621276	27.92%	2.329%
14	8.622	969	975	983	rVB	36023	55942	0.96%	0.080%
15	8.752	990	997	1008	rVB	1291510	2071576	35.67%	2.976%
16	8.864	1008	1016	1021	rBV	43060	76668	1.32%	0.110%
17	9.016	1034	1042	1051	rVB6	19282	47071	0.81%	0.068%
18	9.258	1078	1083	1088	rVB	23094	35867	0.62%	0.052%
19	9.475	1112	1120	1133	rBV	1017270	1724386	29.69%	2.477%
20	9.646	1144	1149	1155	rVB2	83893	129638	2.23%	0.186%
21	9.846	1178	1183	1187	rVB2	24079	37765	0.65%	0.054%
22	10.010	1204	1211	1221	rBV	1581488	2628552	45.26%	3.776%
23	10.134	1222	1232	1236	rVB7	14815	41293	0.71%	0.059%
24	10.387	1268	1275	1287	rVB	1296864	2147137	36.97%	3.084%
25	10.516	1292	1297	1307	rVB	67480	129403	2.23%	0.186%
26	10.675	1316	1324	1332	rVB2	20232	51065	0.88%	0.073%
27	10.805	1332	1346	1355	rBV5	25816	77528	1.34%	0.111%
28	10.993	1373	1378	1390	rVB6	24629	54407	0.94%	0.078%
29	11.399	1443	1447	1456	rVB3	19706	45767	0.79%	0.066%
30	11.475	1456	1460	1472	rVB4	21214	54457	0.94%	0.078%
31	11.575	1472	1477	1488	rBV	127247	250908	4.32%	0.360%
32	11.663	1488	1492	1497	rVB7	25416	40796	0.70%	0.059%
33	11.728	1497	1503	1514	rVB3	89281	164191	2.83%	0.236%
34	11.993	1541	1548	1552	rVV4	59204	125390	2.16%	0.180%

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 Title : SVOA CALIBRATION

35	12.210	1580	1585	1588	rBV5	24414	44522	0.77%	0.064%
36	12.240	1588	1590	1599	rVB	35156	57592	0.99%	0.083%
37	12.346	1601	1608	1611	rBV8	17059	43422	0.75%	0.062%
38	12.381	1611	1614	1623	rVB5	31766	63745	1.10%	0.092%
39	12.504	1630	1635	1640	rBV	73604	113022	1.95%	0.162%
40	12.734	1668	1674	1681	rVV3	235607	406552	7.00%	0.584%
41	12.810	1682	1687	1690	rVV2	45729	79461	1.37%	0.114%
42	12.840	1690	1692	1696	rVB2	37228	44325	0.76%	0.064%
43	13.063	1725	1730	1735	rBV2	27927	51762	0.89%	0.074%
44	13.404	1784	1788	1791	rBV3	22333	36009	0.62%	0.052%
45	13.440	1791	1794	1802	rVB	38573	52639	0.91%	0.076%
46	13.622	1820	1825	1828	rBV4	22921	39715	0.68%	0.057%
47	13.675	1828	1834	1838	rVV	2410153	3478025	59.89%	4.996%
48	13.716	1838	1841	1847	rVB	373937	497229	8.56%	0.714%
49	13.787	1847	1853	1856	rBV6	27487	48273	0.83%	0.069%
50	13.940	1872	1879	1887	rBV	3119896	4889774	84.20%	7.024%
51	14.257	1926	1933	1939	rVV2	1669532	2612319	44.98%	3.753%
52	14.310	1939	1942	1949	rVB3	53620	77655	1.34%	0.112%
53	14.451	1960	1966	1973	rBV	222916	384485	6.62%	0.552%
54	14.510	1973	1976	1981	rVB4	32485	47020	0.81%	0.068%
55	14.822	2024	2029	2037	rVB3	47565	85471	1.47%	0.123%
56	14.928	2043	2047	2056	rVB2	59479	96996	1.67%	0.139%
57	15.022	2059	2063	2070	rVB5	21697	40720	0.70%	0.058%
58	15.104	2071	2077	2081	rBV6	30508	63785	1.10%	0.092%
59	15.251	2096	2102	2111	rBV	3490514	5241400	90.26%	7.529%
60	15.369	2116	2122	2131	rVV	964380	1320998	22.75%	1.898%
61	15.487	2137	2142	2145	rBV2	60214	122560	2.11%	0.176%
62	15.522	2145	2148	2156	rVB2	132105	244775	4.21%	0.352%
63	15.769	2186	2190	2192	rBV2	22709	35461	0.61%	0.051%
64	15.840	2197	2202	2205	rBV2	33847	50788	0.87%	0.073%
65	15.934	2215	2218	2225	rVB6	34278	63681	1.10%	0.091%
66	16.028	2230	2234	2240	rVB3	27511	49429	0.85%	0.071%
67	16.316	2279	2283	2289	rBV4	28161	48575	0.84%	0.070%
68	16.375	2289	2293	2297	rBV3	45187	66929	1.15%	0.096%
69	16.434	2300	2303	2311	rVB4	43977	82268	1.42%	0.118%
70	16.581	2324	2328	2332	rVB2	48750	63668	1.10%	0.091%
71	16.816	2365	2368	2374	rVB	54239	75366	1.30%	0.108%

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72	16.892	2377	2381	2389	rVB5	28043	43107	0.74%	0.062%
73	17.010	2394	2401	2410	rVB	1779310	2584148	44.50%	3.712%
74	17.104	2411	2417	2425	rBV2	3627607	5187378	89.33%	7.452%
75	17.204	2432	2434	2441	rVB7	21589	38674	0.67%	0.056%
76	17.934	2554	2558	2567	rBV2	176709	311159	5.36%	0.447%
77	18.069	2577	2581	2588	rVB	85053	117938	2.03%	0.169%
78	18.716	2687	2691	2695	rBV	99027	125215	2.16%	0.180%
79	18.875	2715	2718	2725	rVB	62884	80818	1.39%	0.116%
80	18.963	2730	2733	2736	rVV2	49472	69788	1.20%	0.100%
81	18.998	2736	2739	2743	rVB	49545	61974	1.07%	0.089%
82	19.086	2751	2754	2758	rVB2	49420	59309	1.02%	0.085%
83	19.139	2760	2763	2766	rVV	30258	38029	0.65%	0.055%
84	19.175	2766	2769	2775	rVV3	41152	62484	1.08%	0.090%
85	19.239	2777	2780	2788	rVB	39326	62033	1.07%	0.089%
86	19.351	2793	2799	2806	rBV	4280039	5802695	99.92%	8.336%
87	20.328	2961	2965	2970	rBV	337412	444869	7.66%	0.639%
88	20.369	2970	2972	2977	rVB2	61323	76255	1.31%	0.110%
89	20.851	3050	3054	3060	rBV	424861	546614	9.41%	0.785%
90	21.104	3092	3097	3109	rBV	2005667	2617927	45.08%	3.761%
91	21.222	3113	3117	3125	rBV	452905	595492	10.25%	0.855%
92	21.292	3126	3129	3134	rVB	92086	110302	1.90%	0.158%
93	21.728	3199	3203	3209	rBV	162119	215937	3.72%	0.310%
94	22.186	3277	3281	3286	rBV	230810	302284	5.21%	0.434%
95	22.692	3363	3367	3373	rBV	271515	383945	6.61%	0.552%
96	23.163	3440	3447	3457	rBV	3285306	5807249	100.00%	8.342%
97	23.263	3459	3464	3465	rVV	295500	388201	6.68%	0.558%
98	23.298	3465	3470	3487	rVB	1549851	3022295	52.04%	4.342%
99	23.916	3569	3575	3581	rBV2	230968	430196	7.41%	0.618%
100	24.669	3698	3703	3709	rVB	118449	219981	3.79%	0.316%

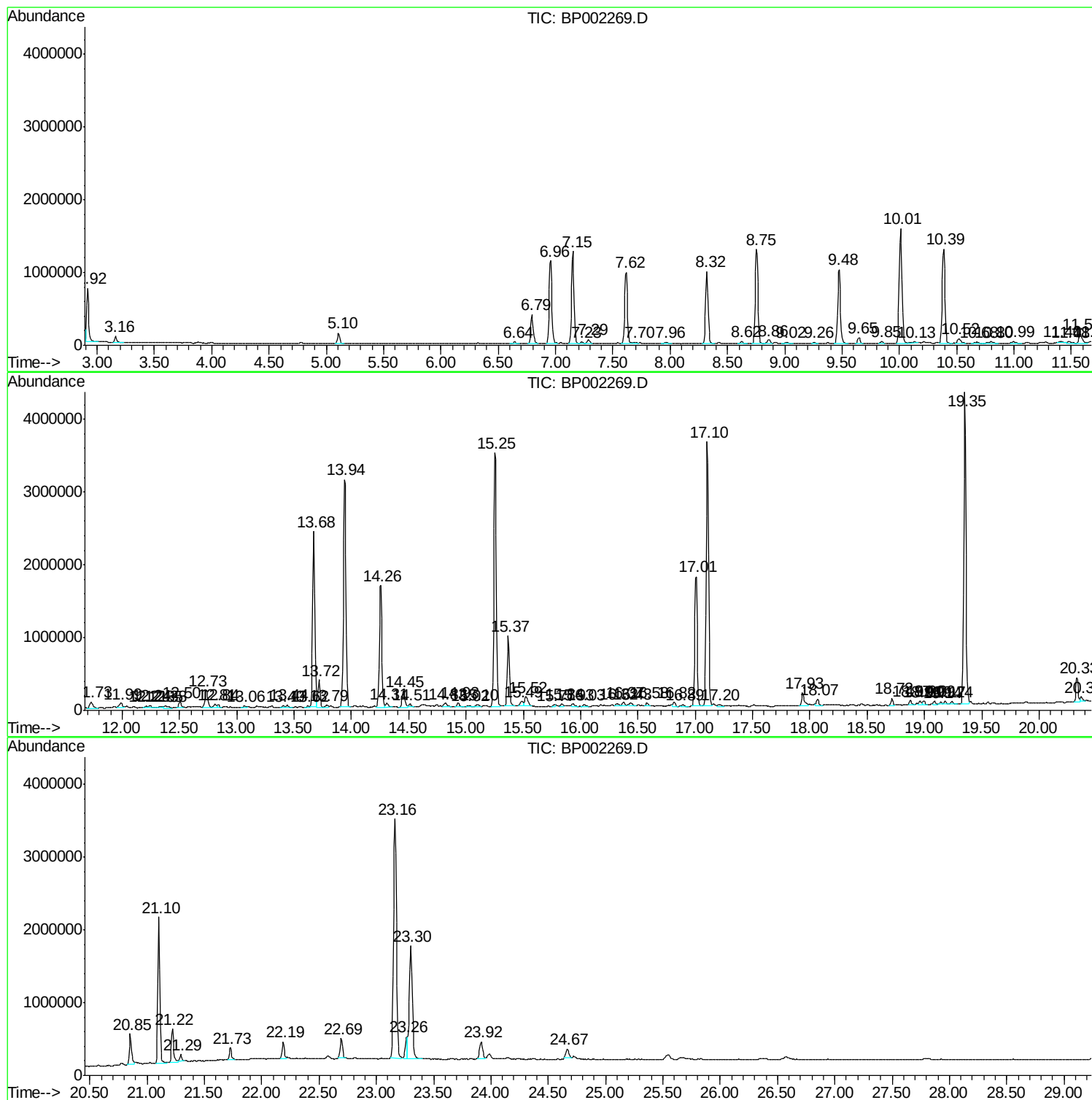
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP051220MA.M
Quant Title : SVOA CALIBRATION

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



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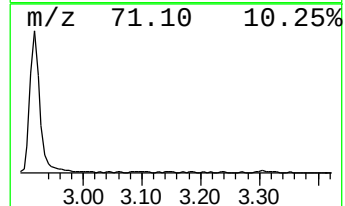
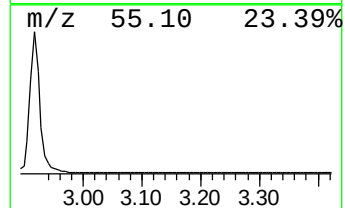
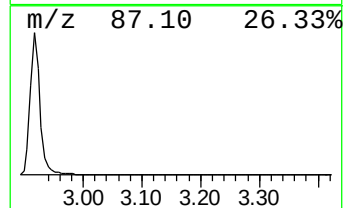
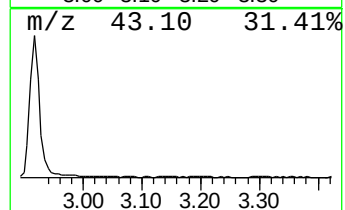
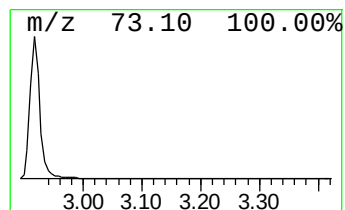
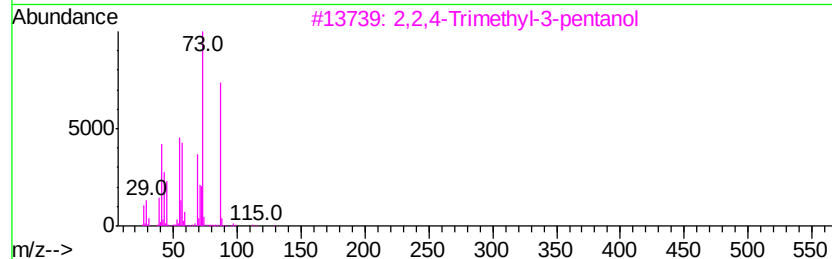
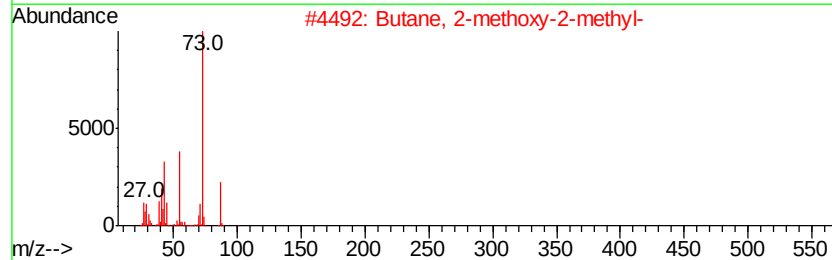
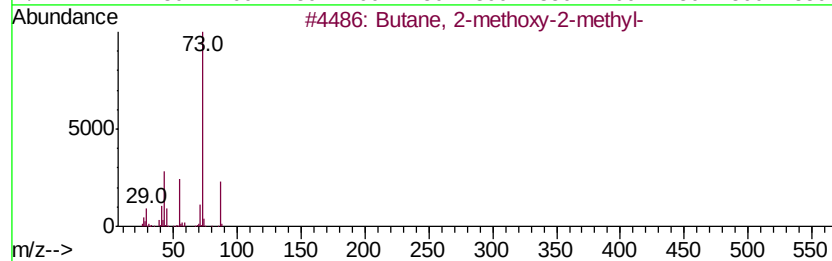
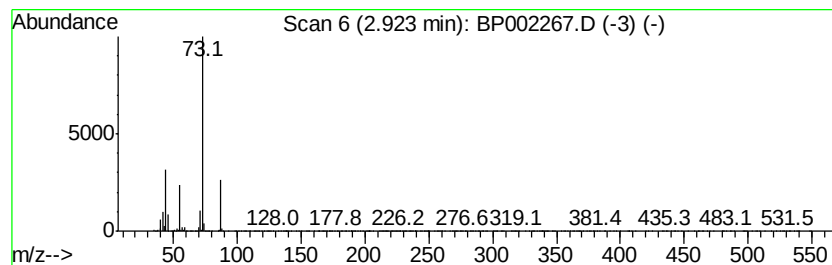
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP051220MA.M
Quant Title : SVOA CALIBRATION

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	9.97 ng/ul	784264	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23



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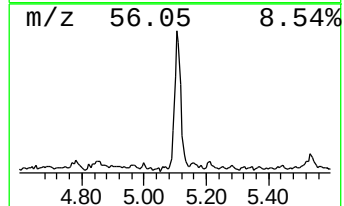
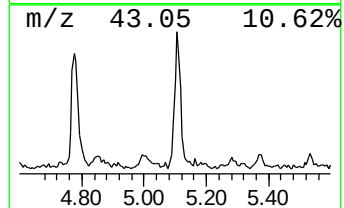
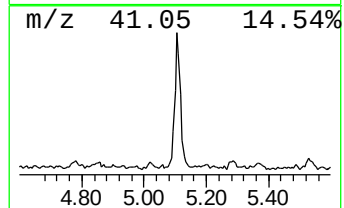
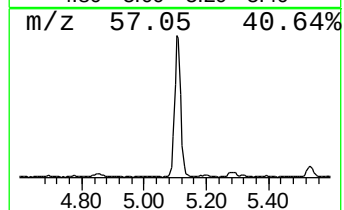
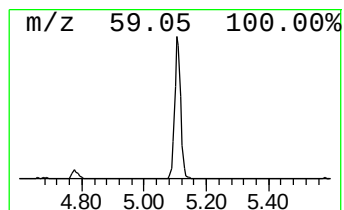
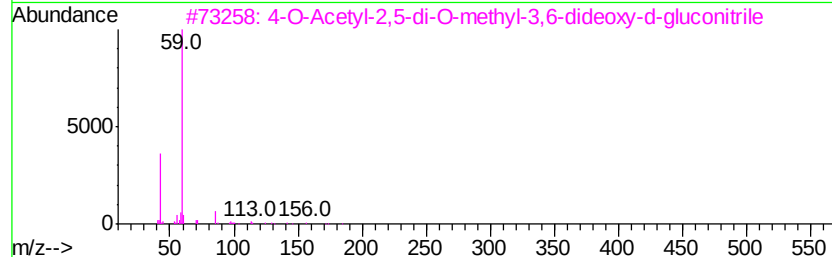
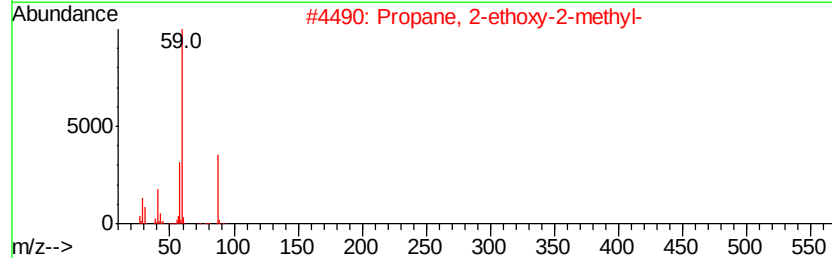
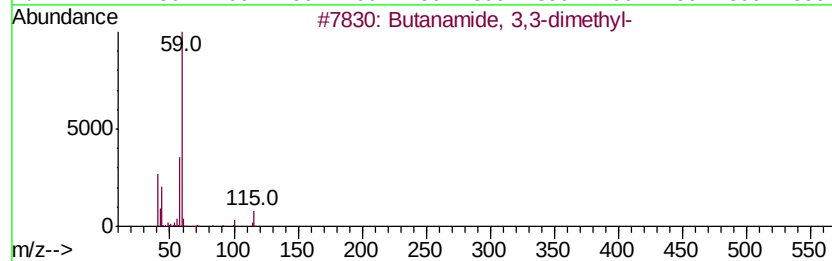
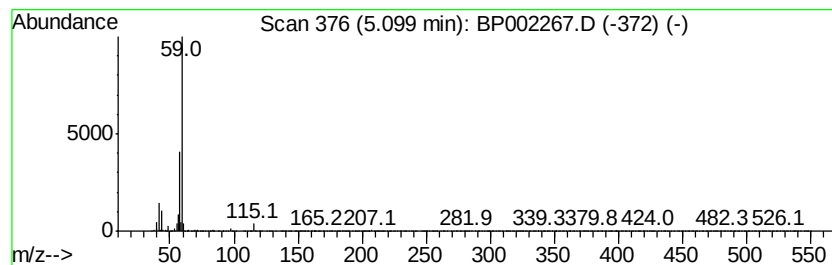
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP051220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butanamide, 3,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	2.56 ng/ul	201715	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	72
2		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	40
3		4-O-Acetyl-2,5-di-O-methyl-3,6-d...	215	C10H17NO4	1000101-80-3	39
4		2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9
5		2-Heptanol, 2-methyl-	130	C8H18O	000625-25-2	9



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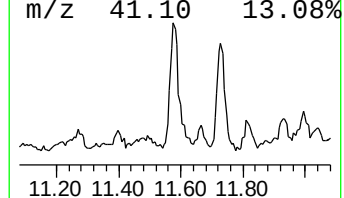
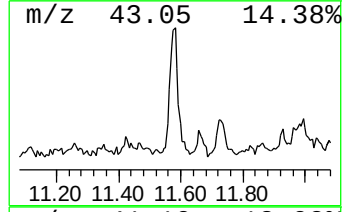
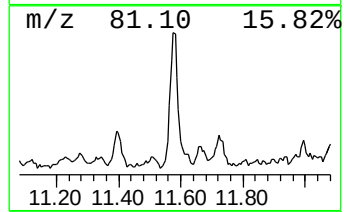
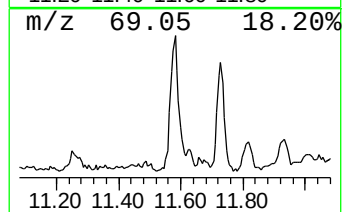
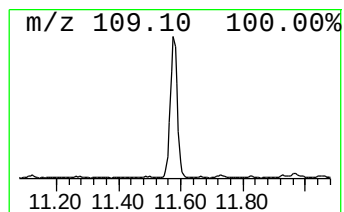
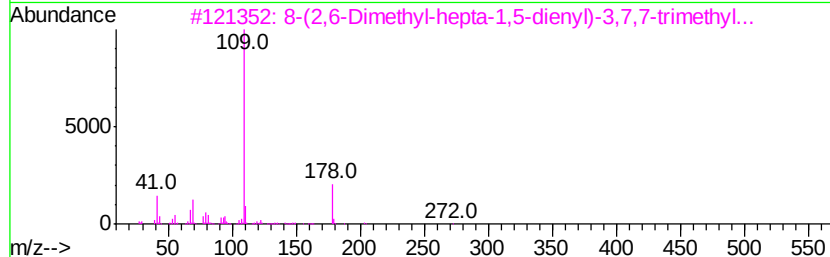
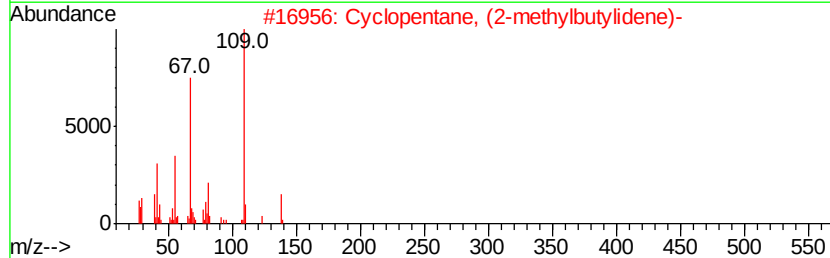
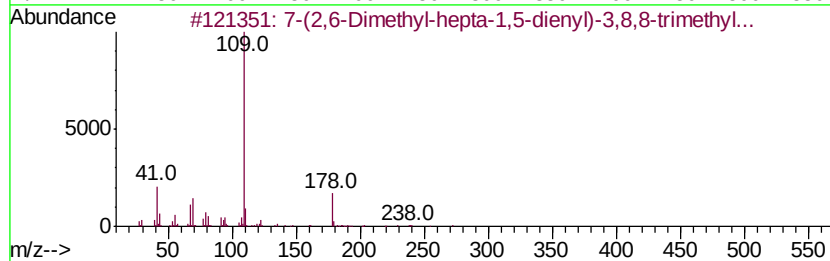
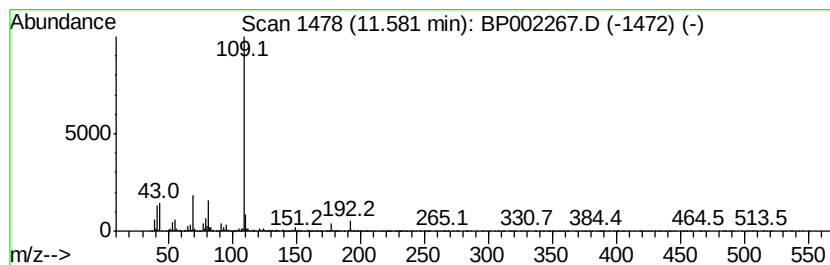
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 7-(2,6-Dimethyl-hepta-1,5-d... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.58	2.34 ng/ul	250908	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7-(2,6-Dimethyl-hepta-1,5-dienvl...	272	C20H32	1000190-62-8	64
2		Cyclopentane, (2-methylbutylidene)-	138	C10H18	053366-54-4	59
3		8-(2,6-Dimethyl-hepta-1,5-dienvl...	272	C20H32	113725-56-7	50
4		Phenol, 3-amino-	109	C6H7NO	000591-27-5	45
5		2,3-Pyridinediamine	109	C5H7N3	000452-58-4	45



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Data File : BP002267.D
Acq On : 13 May 2020 01:24
Operator : CG/JU
Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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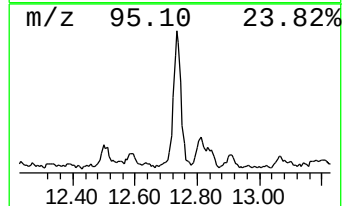
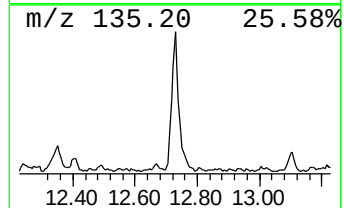
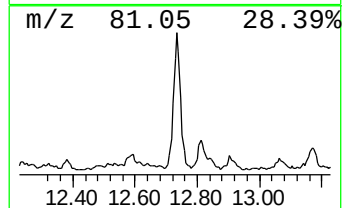
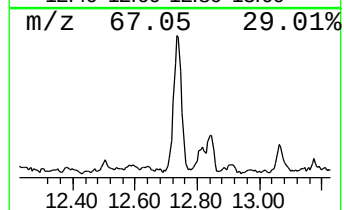
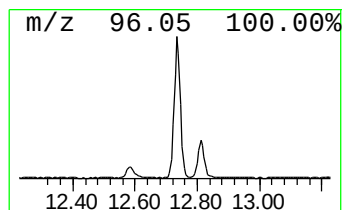
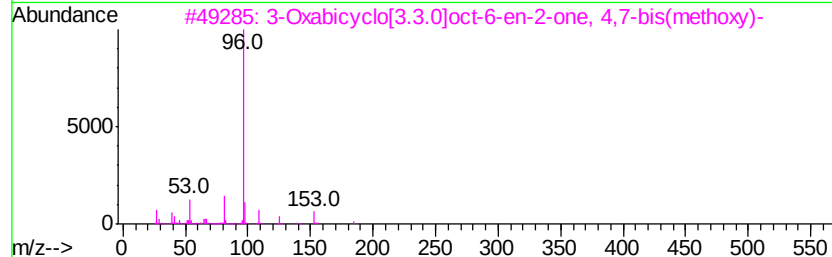
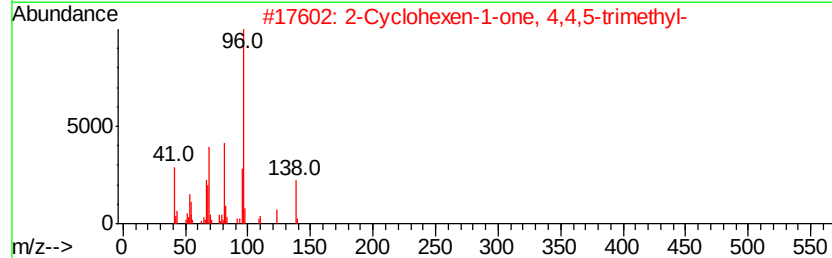
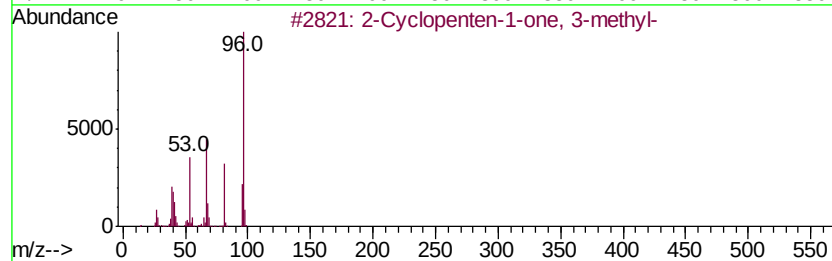
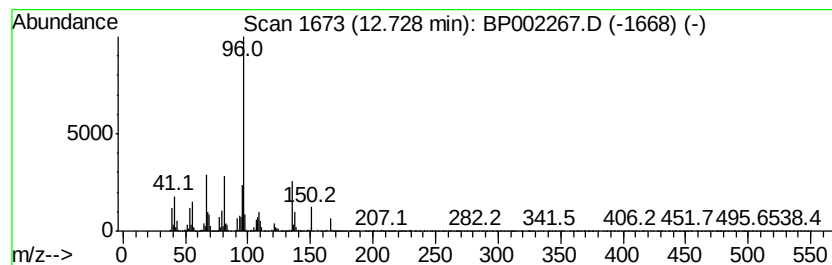
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Cyclopenten-1-one, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	3.11 ng/ul	406552	Acenaphthene-d10	14.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Cyclopenten-1-one, 3-methyl-	96	C6H8O	002758-18-1	53
2		2-Cyclohexen-1-one, 4,4,5-trimet...	138	C9H14O	017429-29-7	50
3		3-Oxabicyclo[3.3.0]oct-6-en-2-on...	184	C9H12O4	1000155-61-6	43
4		2-Cyclohexen-1-one, 4,4,6-trimet...	138	C9H14O	013395-73-8	42
5		12-Heptadecyn-1-ol	252	C17H32O	056554-76-8	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
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Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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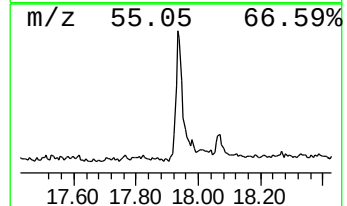
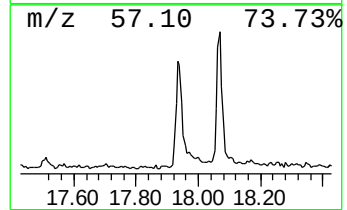
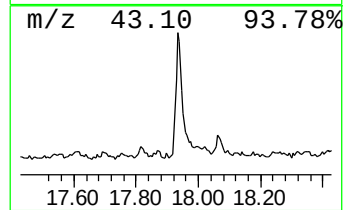
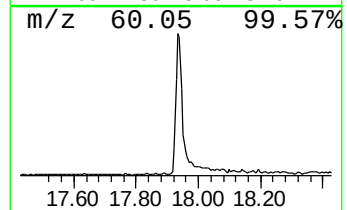
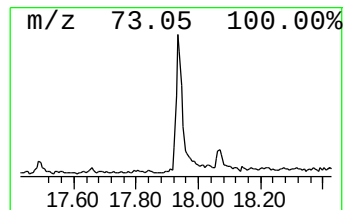
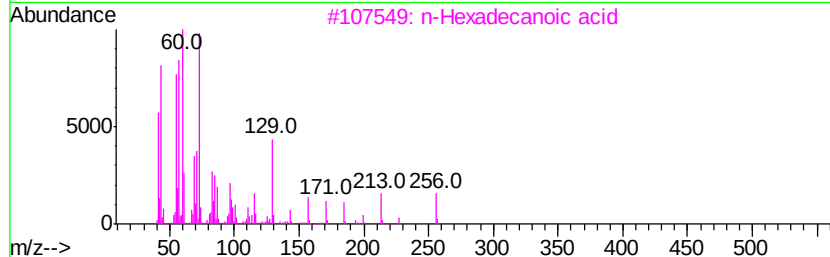
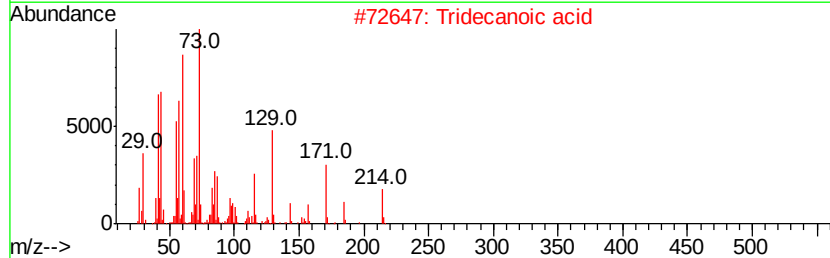
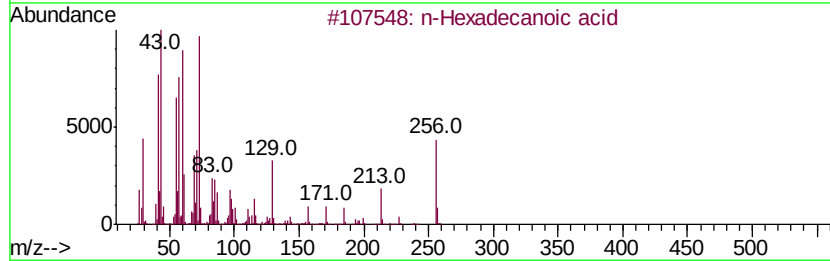
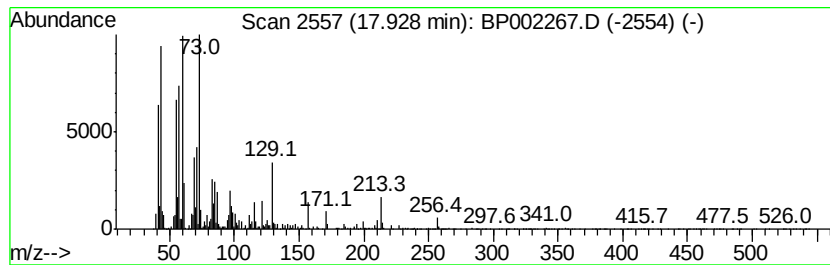
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.93	2.41 ng/ul	311159	Phenanthrene-d10	17.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5		Tridecanoic acid	214	C13H26O2	000638-53-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
Operator : CG/JU
Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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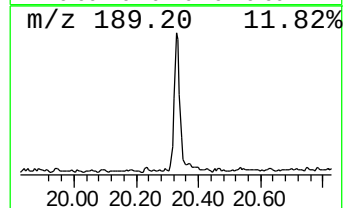
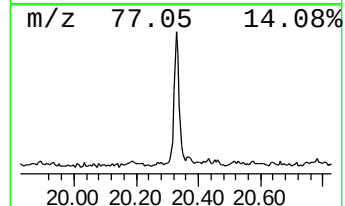
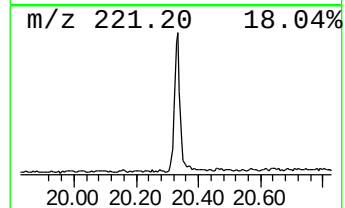
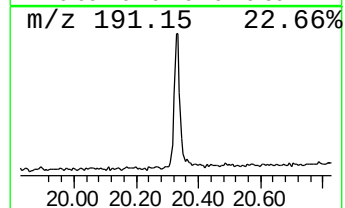
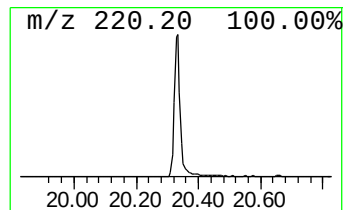
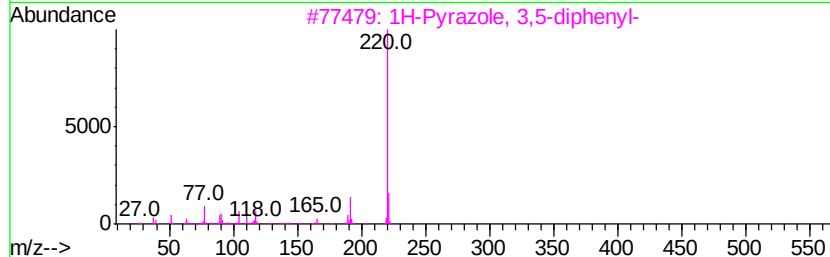
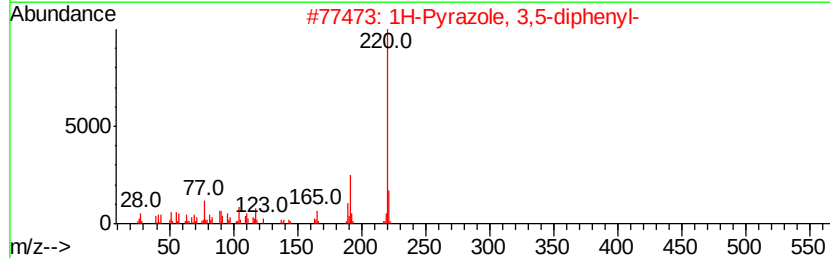
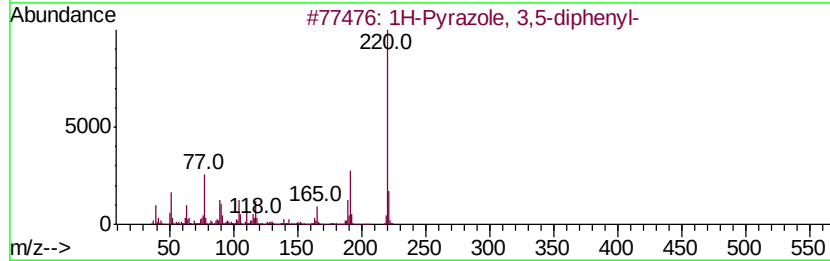
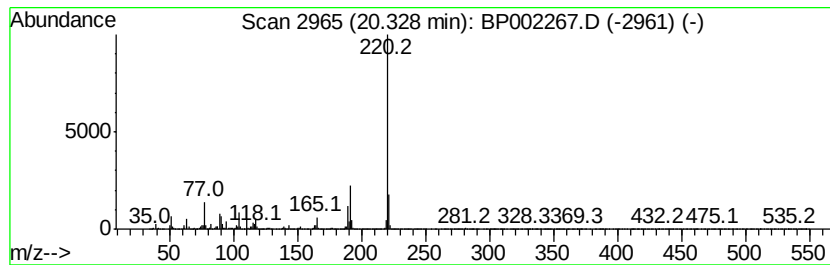
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Pyrazole, 3,5-diphenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.33	3.40 ng/ul	444869	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	97
2		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	96
3		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	94
4		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	91
5		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
Operator : CG/JU
Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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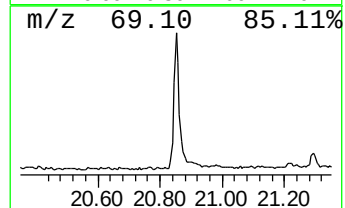
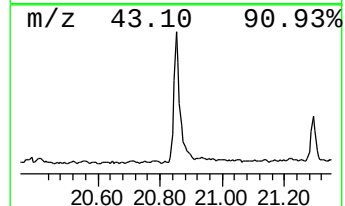
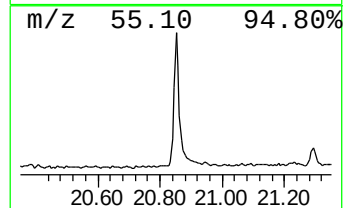
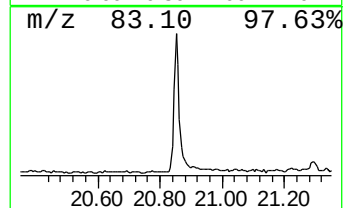
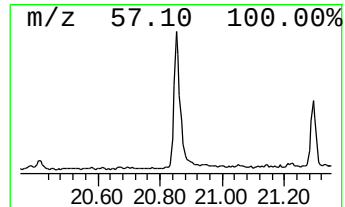
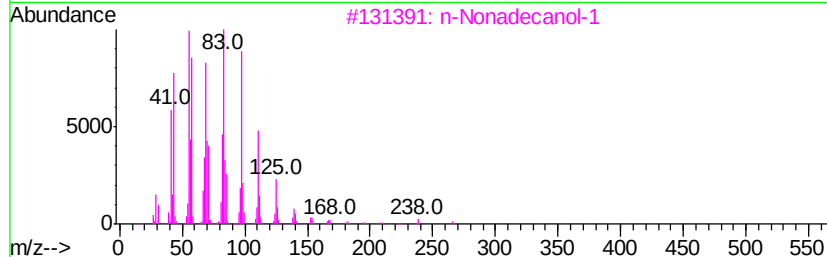
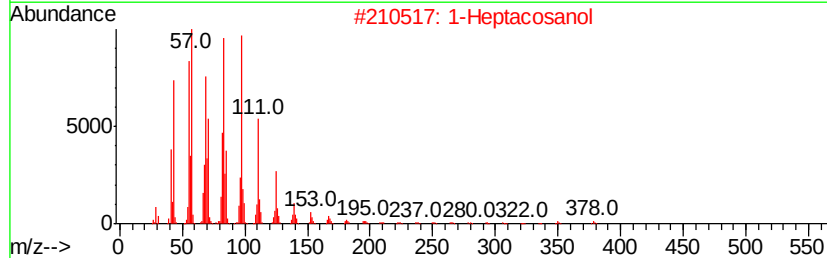
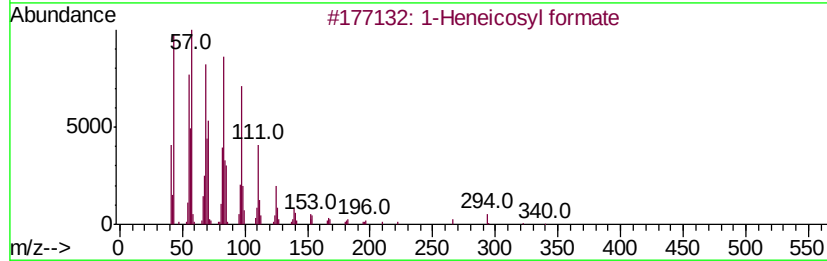
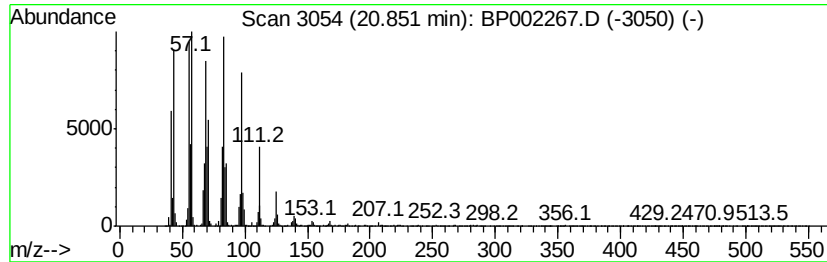
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	4.18 ng/ul	546614	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	96
2		1-Heptacosanol	396	C27H56O	002004-39-9	94
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
5		Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
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Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

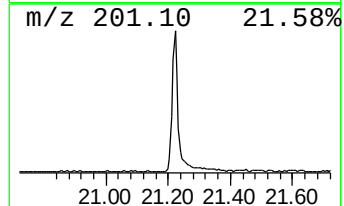
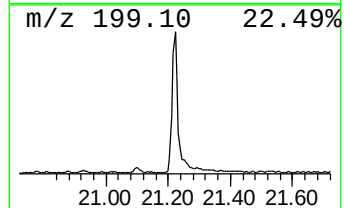
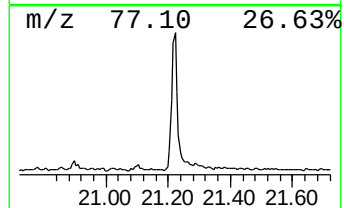
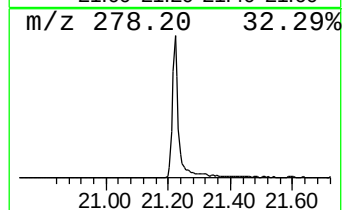
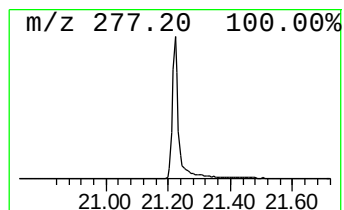
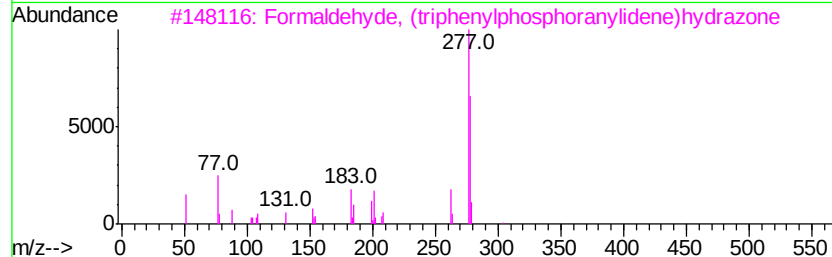
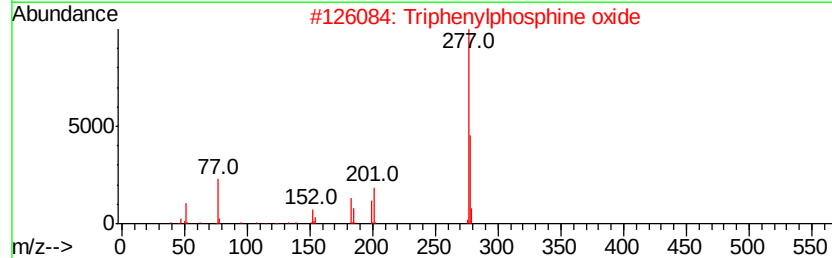
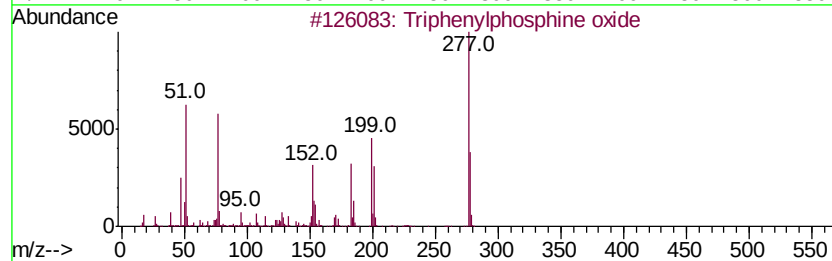
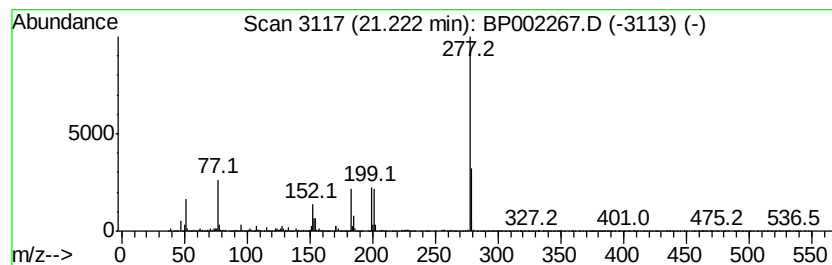
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Triphenylphosphine oxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.22	4.55 ng/ul	595492	Chrysene-d12		21.10	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triphenylphosphine oxide	278	C18H15OP	000791-28-6	93
2		Triphenylphosphine oxide	278	C18H15OP	000791-28-6	52
3		Formaldehydde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	50
4		Triphenylphosphine oxide	278	C18H15OP	000791-28-6	43
5		Triphenylphosphine oxide	278	C18H15OP	000791-28-6	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
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Sample : L2570-03
Misc :
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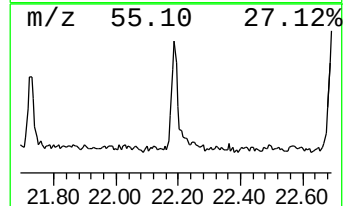
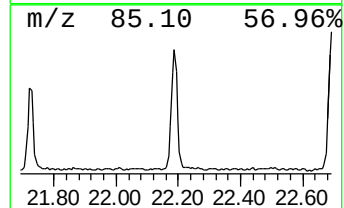
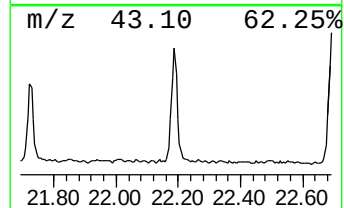
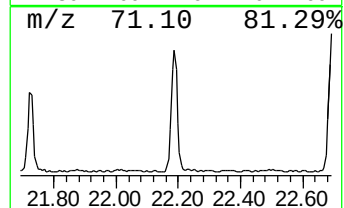
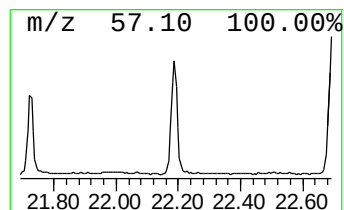
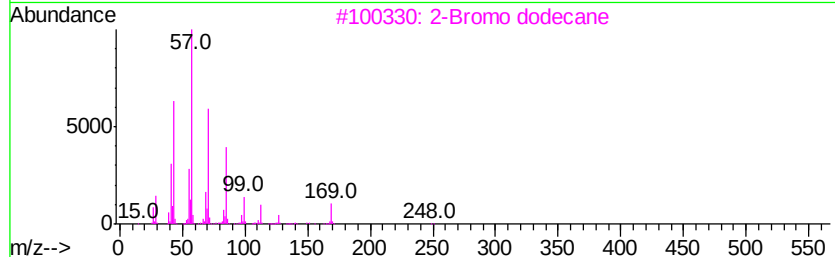
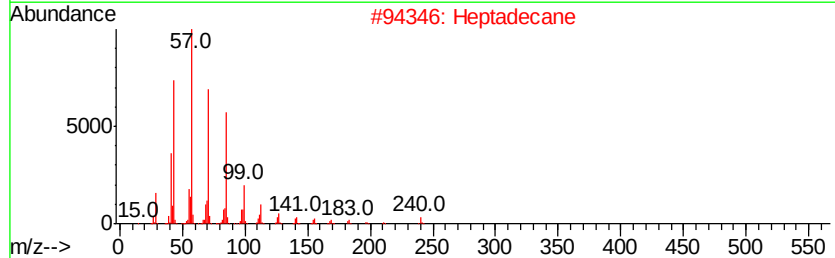
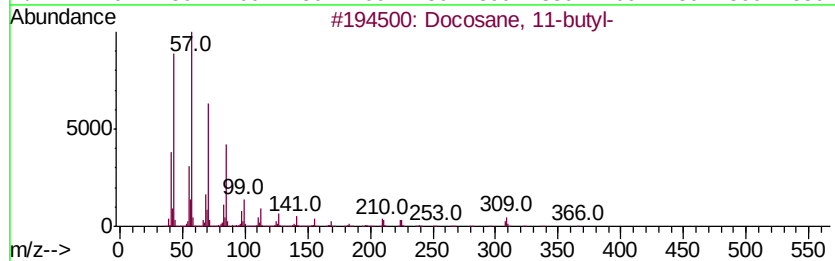
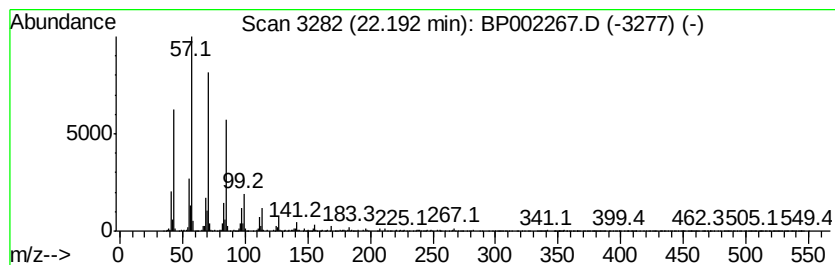
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.19	2.31 ng/ul	302284	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
2		Heptadecane	240	C17H36	000629-78-7	94
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	92
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
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Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

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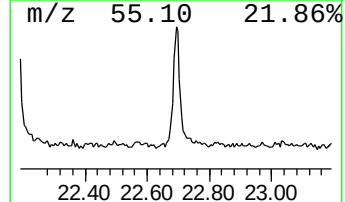
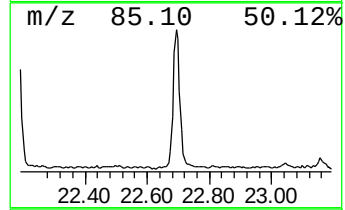
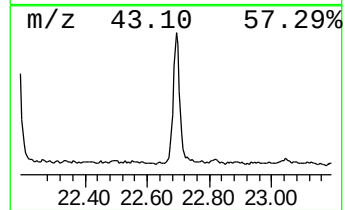
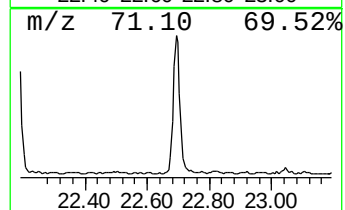
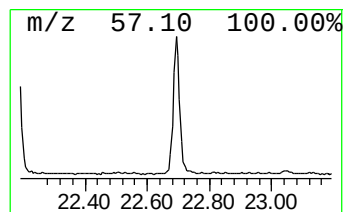
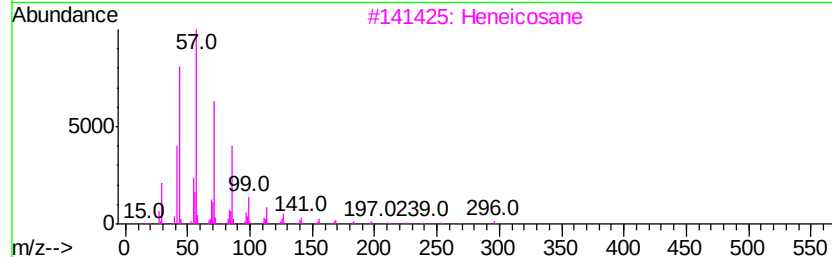
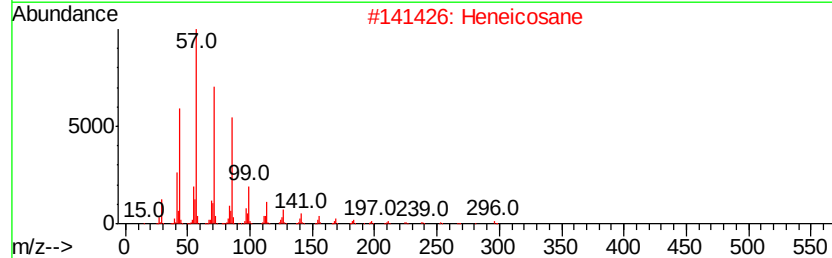
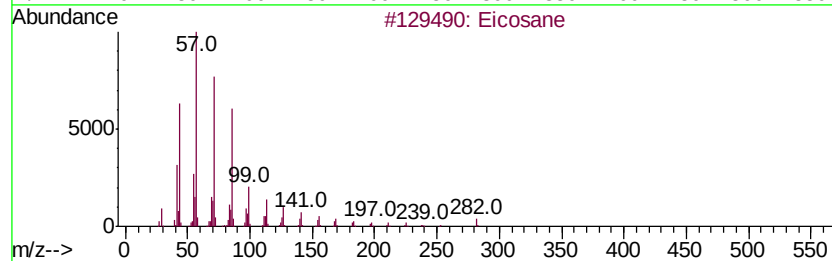
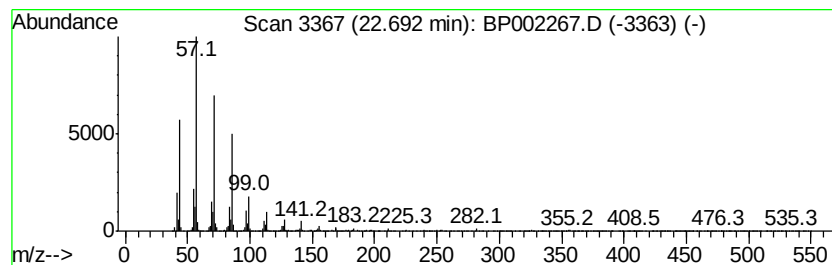
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	2.54 ng/ul	383945	Perylene-d12	23.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Heneicosane	296	C21H44	000629-94-7	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Hexatriacontane	507	C36H74	000630-06-8	90
5			Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
Operator : CG/JU
Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFSP3

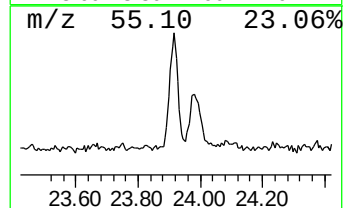
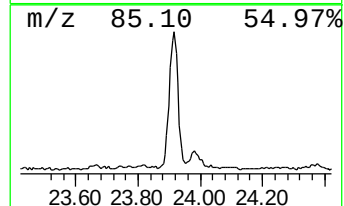
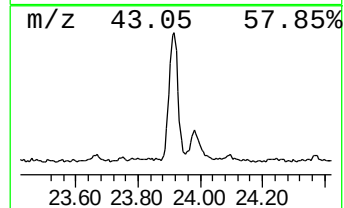
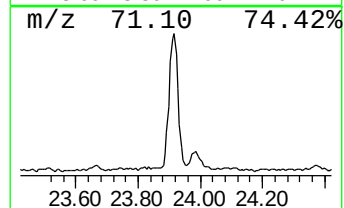
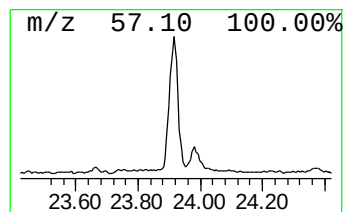
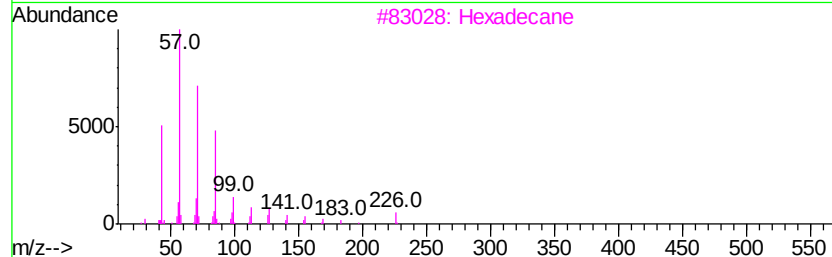
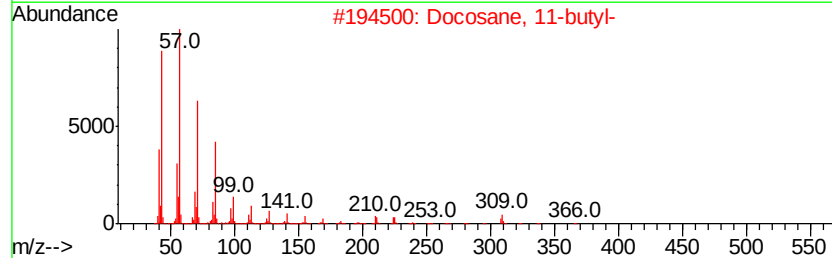
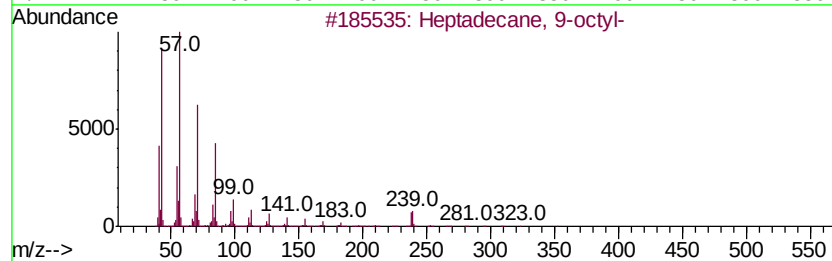
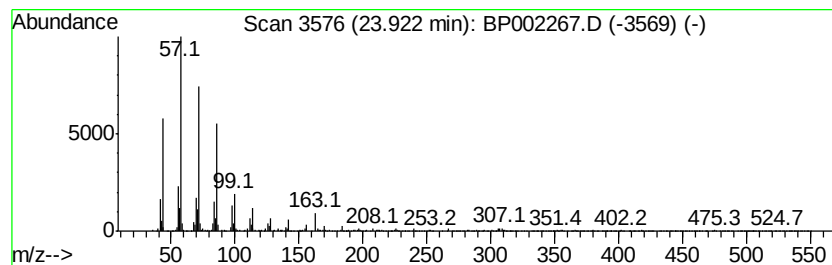
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.92	2.85 ng/ul	430196	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3		Hexadecane	226	C16H34	000544-76-3	90
4		Docosane, 5-butyl-	366	C26H54	055282-16-1	90
5		Docosane, 9-butyl-	366	C26H54	055282-14-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP051220\
Data File : BP002267.D
Acq On : 13 May 2020 01:24
Operator : CG/JU
Sample : L2570-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFSP3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP051220MA.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	10.0	ng/ul	784264	1	7.62	1573890	20.0
Butanamide, 3,3-d...	5.10	2.6	ng/ul	201715	1	7.62	1573890	20.0
7-(2,6-Dimethyl-h...	11.58	2.3	ng/ul	250908	2	10.39	2147140	20.0
2-Cyclopenten-1-o...	12.73	3.1	ng/ul	406552	3	14.26	2612320	20.0
n-Hexadecanoic acid	17.93	2.4	ng/ul	311159	4	17.00	2584150	20.0
1H-Pyrazole, 3,5-...	20.33	3.4	ng/ul	444869	5	21.10	2617930	20.0
1-Heneicosyl formate	20.85	4.2	ng/ul	546614	5	21.10	2617930	20.0
Triphenylphosphin...	21.22	4.5	ng/ul	595492	5	21.10	2617930	20.0
(DEL) Alkane: Str...	22.19	2.3	ng/ul	302284	5	21.10	2617930	20.0
(DEL) Alkane: Str...	22.69	2.5	ng/ul	383945	6	23.30	3022300	20.0
(DEL) Alkane: Str...	23.92	2.9	ng/ul	430196	6	23.30	3022300	20.0