

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
 Data File : BP020364.D
 Acq On : 13 May 2024 22:13
 Operator : MA/JU
 Sample : P2369-13MEDL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43N7MEDL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020364.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.646	91	95	105	rBV2	16884	43032	2.94%	0.216%
2	5.587	420	425	429	rBV2	17191	27088	1.85%	0.136%
3	5.646	429	435	443	rVB3	29135	52592	3.59%	0.264%
4	5.887	469	476	484	rBV4	37401	98470	6.72%	0.494%
5	6.005	489	496	503	rVB2	30604	60026	4.09%	0.301%
6	6.093	503	511	513	rBV2	30644	47776	3.26%	0.240%
7	6.164	519	523	528	rBV	67016	90262	6.16%	0.453%
8	6.211	528	531	533	rBV2	19668	26404	1.80%	0.133%
9	6.269	533	541	550	rVB3	76619	208131	14.20%	1.045%
10	6.358	551	556	561	rVB3	23180	36886	2.52%	0.185%
11	6.422	561	567	572	rBV4	28452	46798	3.19%	0.235%
12	6.481	572	577	578	rBV2	36600	55882	3.81%	0.281%
13	6.552	585	589	592	rBV3	29410	46213	3.15%	0.232%
14	6.593	592	596	601	rVB	75581	110529	7.54%	0.555%
15	6.646	601	605	609	rVB	63723	84896	5.79%	0.426%
16	6.693	609	613	616	rBV3	34552	50470	3.44%	0.253%
17	6.752	616	623	629	rVB3	114756	234882	16.02%	1.179%
18	6.834	632	637	642	rBV2	59553	97159	6.63%	0.488%
19	6.917	647	651	656	rVB4	40975	74128	5.06%	0.372%
20	6.975	656	661	669	rBV4	67348	192093	13.10%	0.964%
21	7.040	670	672	675	rVV3	38584	54520	3.72%	0.274%
22	7.075	675	678	683	rVV3	69616	124411	8.49%	0.625%
23	7.152	686	691	697	rVV3	71775	152629	10.41%	0.766%
24	7.211	697	701	706	rVV	308966	447155	30.50%	2.245%
25	7.258	706	709	713	rVV2	43855	67741	4.62%	0.340%
26	7.305	713	717	722	rVB3	46615	76190	5.20%	0.382%
27	7.411	727	735	740	rVB5	47718	125231	8.54%	0.629%
28	7.499	740	750	754	rBV4	92342	242334	16.53%	1.216%
29	7.558	755	760	764	rVV	291365	402590	27.46%	2.021%
30	7.611	764	769	780	rVB2	254647	603525	41.17%	3.030%
31	7.711	780	786	790	rBV2	89878	182548	12.45%	0.916%
32	7.764	791	795	799	rVB2	102831	150575	10.27%	0.756%
33	7.828	799	806	809	rBV2	66134	122568	8.36%	0.615%
34	7.864	809	812	819	rVB2	102061	161786	11.04%	0.812%
35	7.922	819	822	829	rVB4	46429	90729	6.19%	0.455%
36	8.016	832	838	844	rVB5	44413	101234	6.91%	0.508%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
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37	8.099	844	852	856	rBV3	159982	352734	24.06%	1.771%
38	8.164	860	863	868	rVB	137563	187832	12.81%	0.943%
39	8.234	868	875	880	rBV2	213033	383096	26.13%	1.923%
40	8.340	888	893	898	rBV2	162944	279294	19.05%	1.402%
41	8.405	898	904	909	rVB2	96264	216098	14.74%	1.085%
42	8.540	920	927	932	rBV2	62517	138550	9.45%	0.695%
43	8.587	932	935	940	rVB	68068	96706	6.60%	0.485%
44	8.646	940	945	946	rBV2	60246	80439	5.49%	0.404%
45	8.799	962	971	980	rVB	671580	1275242	86.99%	6.401%
46	8.881	980	985	987	rBV3	62984	106347	7.25%	0.534%
47	8.928	987	993	999	rVB8	110030	250863	17.11%	1.259%
48	9.022	1001	1009	1017	rBV6	77367	214723	14.65%	1.078%
49	9.210	1036	1041	1047	rBV3	76155	134437	9.17%	0.675%
50	9.281	1047	1053	1058	rVB5	49938	105577	7.20%	0.530%
51	9.446	1074	1081	1085	rBV4	40822	103795	7.08%	0.521%
52	9.493	1085	1089	1094	rVV4	67714	167560	11.43%	0.841%
53	9.540	1094	1097	1101	rVV4	61805	110425	7.53%	0.554%
54	9.581	1101	1104	1108	rVV3	45303	87560	5.97%	0.440%
55	9.640	1109	1114	1119	rVV4	48278	103215	7.04%	0.518%
56	9.710	1121	1126	1133	rVV3	47887	103309	7.05%	0.519%
57	9.787	1134	1139	1145	rVV3	77756	170723	11.65%	0.857%
58	9.846	1145	1149	1152	rVV2	22237	40002	2.73%	0.201%
59	9.893	1152	1157	1161	rVV	38491	57949	3.95%	0.291%
60	9.952	1161	1167	1173	rVB5	32661	83326	5.68%	0.418%
61	10.081	1182	1189	1195	rBV7	34231	85252	5.82%	0.428%
62	10.234	1211	1215	1219	rBV6	10299	22478	1.53%	0.113%
63	10.340	1227	1233	1236	rBV	129093	227877	15.55%	1.144%
64	10.381	1236	1240	1253	rVB	327057	653753	44.60%	3.282%
65	10.522	1260	1264	1269	rVB	26766	45224	3.09%	0.227%
66	10.587	1269	1275	1284	rBV3	25136	61711	4.21%	0.310%
67	10.746	1297	1302	1310	rVB7	10939	23561	1.61%	0.118%
68	11.663	1452	1458	1464	rBV5	10849	25173	1.72%	0.126%
69	11.804	1476	1482	1487	rBV3	13033	24788	1.69%	0.124%
70	12.246	1551	1557	1562	rBV5	16370	33968	2.32%	0.171%
71	12.769	1635	1646	1660	rVB5	8388	32962	2.25%	0.165%
72	13.163	1707	1713	1719	rBV8	8229	21931	1.50%	0.110%
73	13.322	1733	1740	1746	rBV6	9348	20965	1.43%	0.105%
74	13.704	1799	1805	1817	rBV2	79126	183493	12.52%	0.921%
75	13.969	1845	1850	1857	rBV	110524	183154	12.49%	0.919%
76	14.275	1896	1902	1910	rBV	553342	918478	62.66%	4.611%

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 Title : SVOA CALIBRATION

77	14.346	1910	1914	1918	rVB2	28281	39930	2.72%	0.200%
78	14.493	1934	1939	1940	rBV3	16332	22945	1.57%	0.115%
79	14.516	1940	1943	1947	rVB2	20495	28747	1.96%	0.144%
80	14.681	1966	1971	1975	rBV5	13648	22159	1.51%	0.111%
81	14.875	1999	2004	2009	rVV8	12088	19828	1.35%	0.100%
82	14.945	2011	2016	2027	rVB6	24649	57691	3.94%	0.290%
83	15.140	2044	2049	2053	rVV3	14167	19935	1.36%	0.100%
84	15.310	2070	2078	2088	rBV	137996	259304	17.69%	1.302%
85	15.493	2104	2109	2114	rBV6	19182	41938	2.86%	0.211%
86	15.922	2178	2182	2189	rVB2	15050	25116	1.71%	0.126%
87	16.769	2319	2326	2341	rBV	711773	1204098	82.14%	6.044%
88	17.122	2380	2386	2399	rBV	770778	1212212	82.69%	6.085%
89	17.228	2399	2404	2415	rVB	168681	285597	19.48%	1.434%
90	18.075	2543	2548	2558	rVB3	31805	54529	3.72%	0.274%
91	19.422	2774	2777	2784	rBV6	14829	23864	1.63%	0.120%
92	19.628	2806	2812	2823	rBV	191580	409216	27.92%	2.054%
93	20.228	2910	2914	2920	rBV	93613	122741	8.37%	0.616%
94	20.745	2998	3002	3010	rVB	157847	195834	13.36%	0.983%
95	21.292	3090	3095	3103	rBV	147041	212855	14.52%	1.068%
96	21.610	3143	3149	3162	rBV	787726	1392129	94.97%	6.988%
97	21.869	3189	3193	3200	rVB	90001	140468	9.58%	0.705%
98	22.516	3299	3303	3312	rVB	64824	104263	7.11%	0.523%
99	24.739	3675	3681	3695	rVB2	123173	356067	24.29%	1.787%
100	24.968	3712	3720	3733	rVB2	524084	1465888	100.00%	7.358%

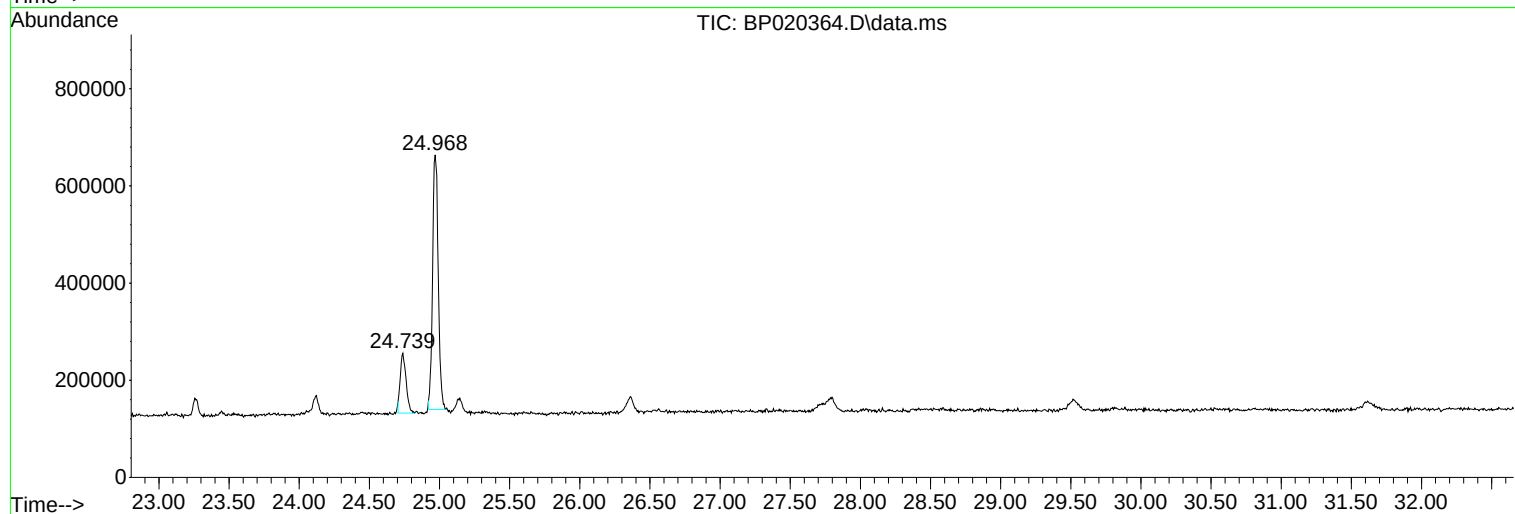
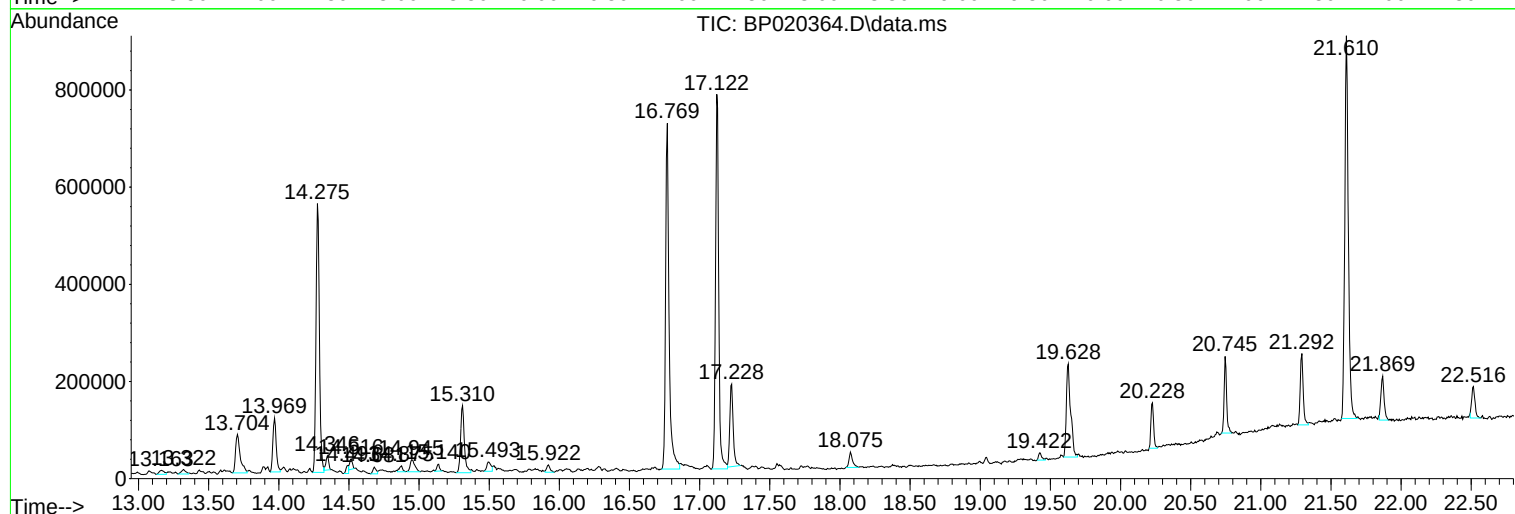
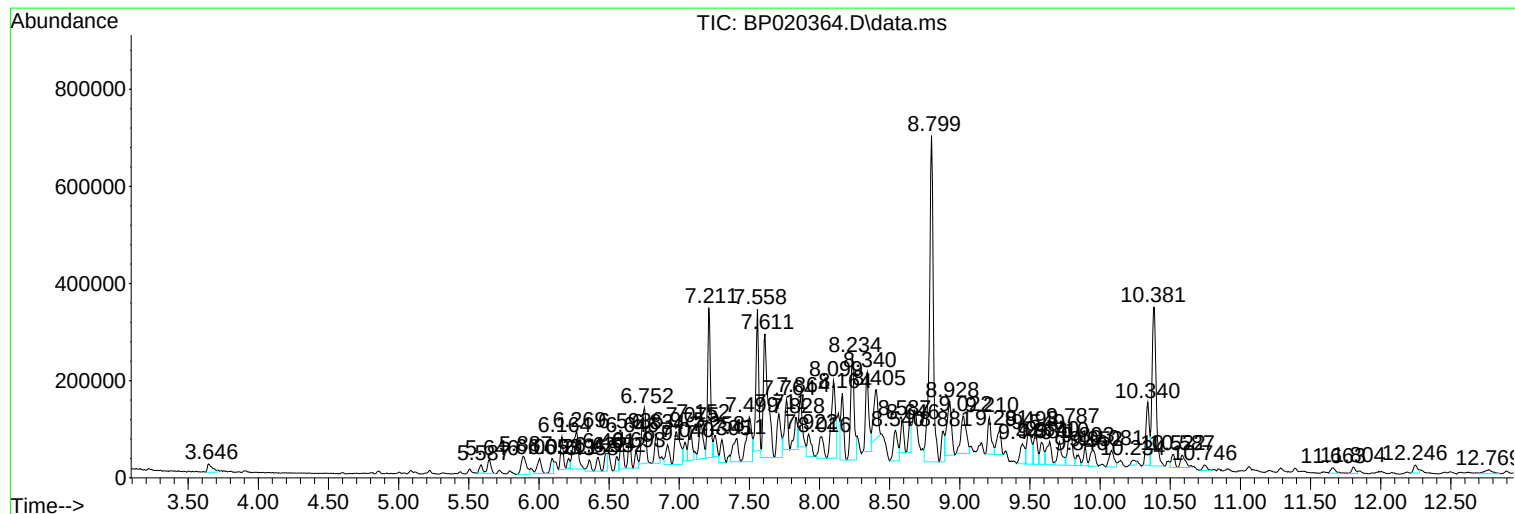
Sum of corrected areas: 19921407

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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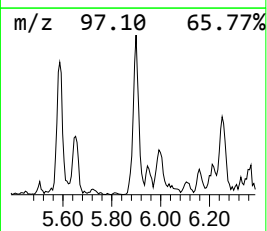
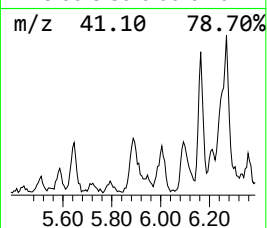
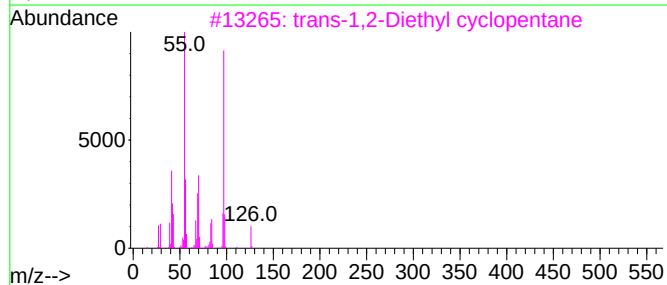
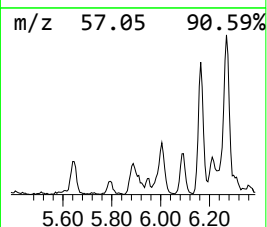
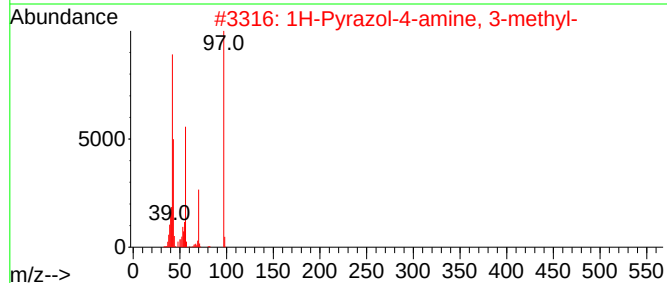
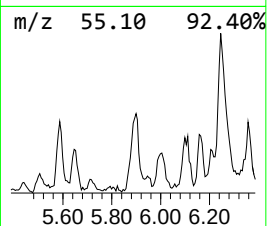
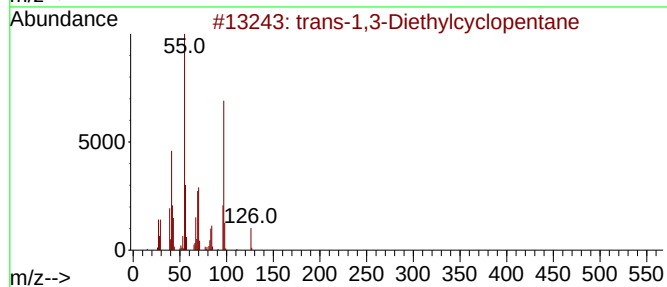
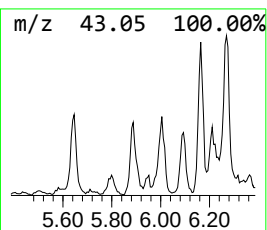
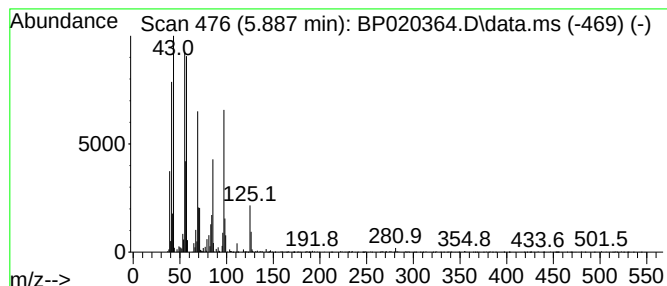
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Cyclic5.887 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.887	3.26 ng/ul	98470	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-1,3-Diethylcyclopentane	126	C9H18	1000113-87-1	45
2			1H-Pyrazol-4-amine, 3-methyl-	97	C4H7N3	1010338-28-2	38
3			trans-1,2-Diethyl cyclopentane	126	C9H18	000932-40-1	30
4			Cyclohexanone, 4-ethyl-	126	C8H14O	005441-51-0	18
5			3-Heptene, 2,6-dimethyl-	126	C9H18	002738-18-3	18



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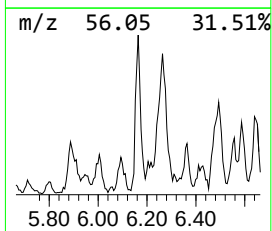
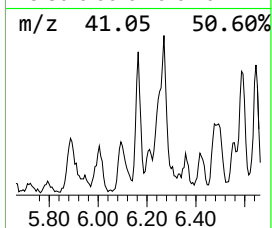
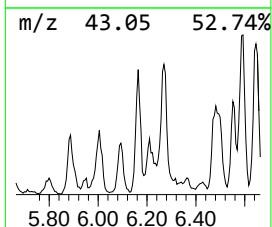
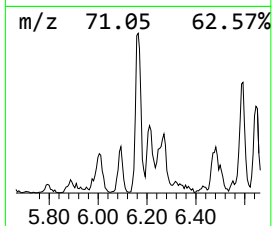
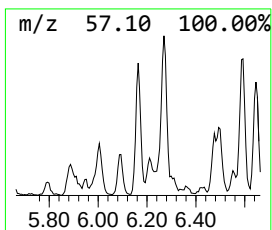
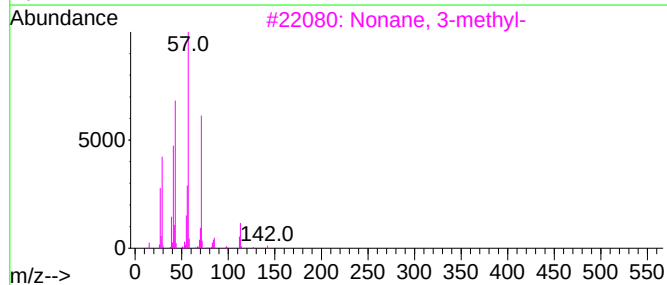
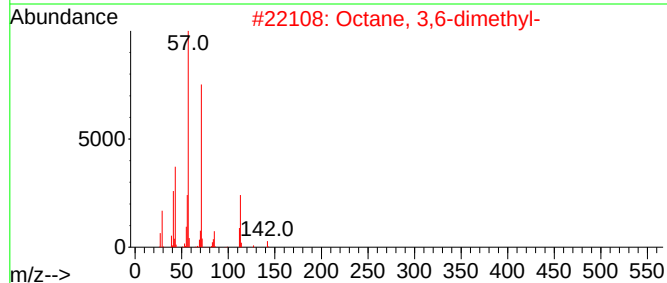
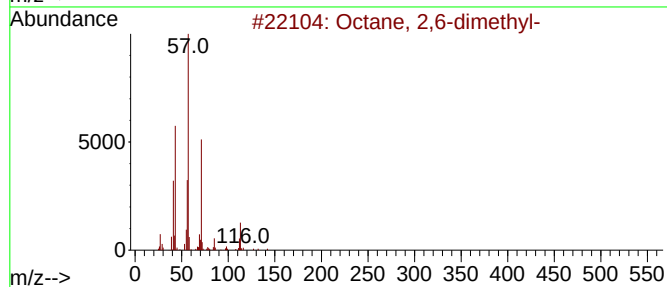
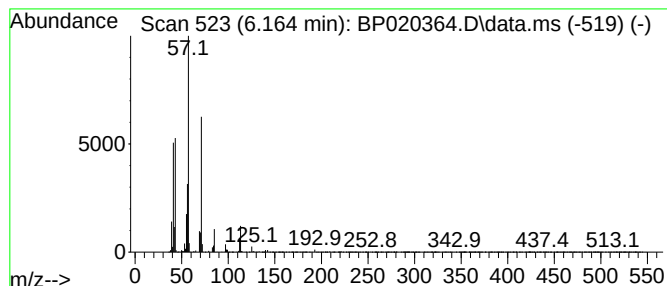
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.164	2.99 ng/ul	90262	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	91	
2	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	87	
3	Nonane, 3-methyl-		142	C10H22	005911-04-6	87	
4	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	87	
5	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	83	



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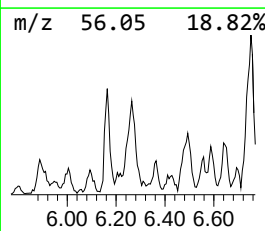
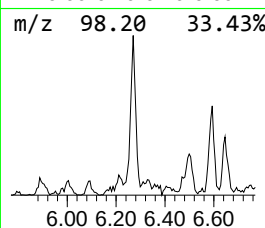
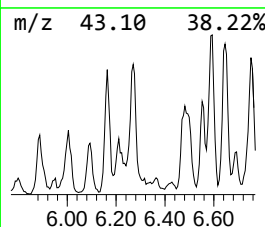
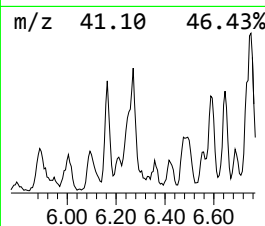
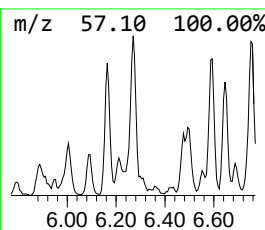
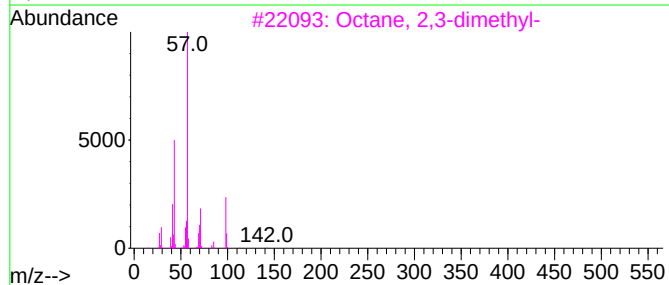
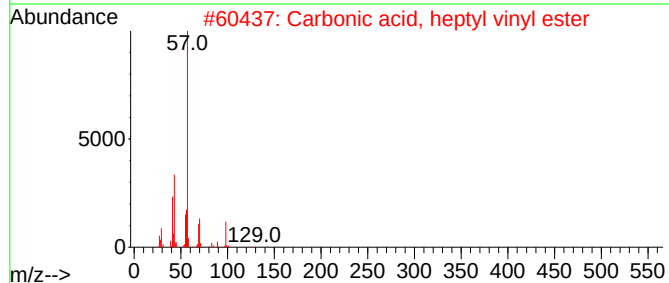
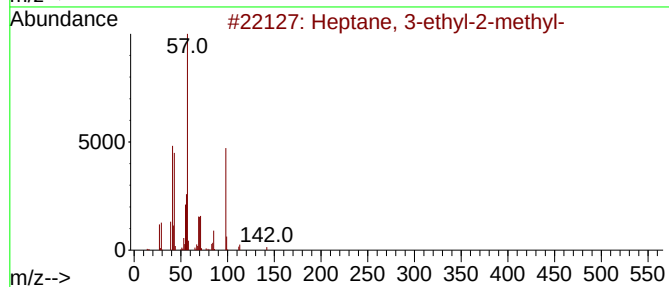
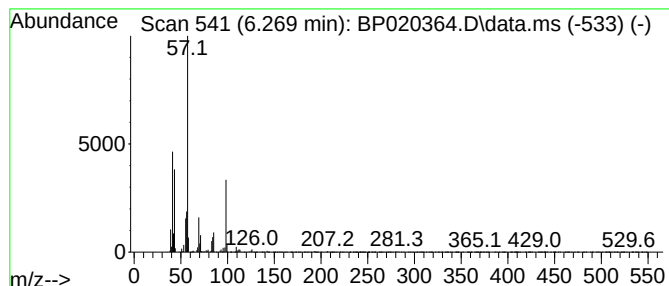
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.269	6.90 ng/ul	208131	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	64
2			Carbonic acid, heptyl vinyl ester	186	C10H18O3	1010383-25-4	50
3			Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	50
4			Ether, tert-butyl isopropylidene...	154	C10H18O	024524-56-9	50
5			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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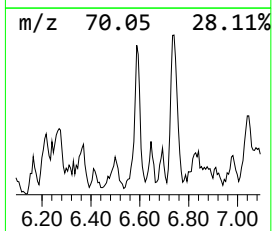
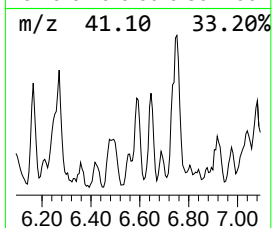
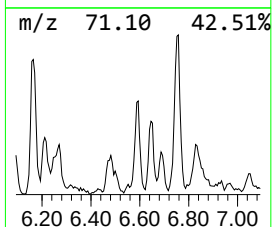
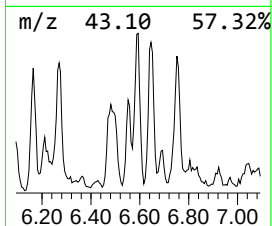
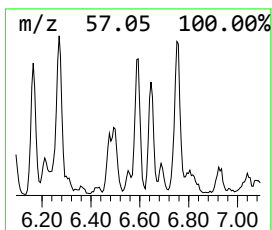
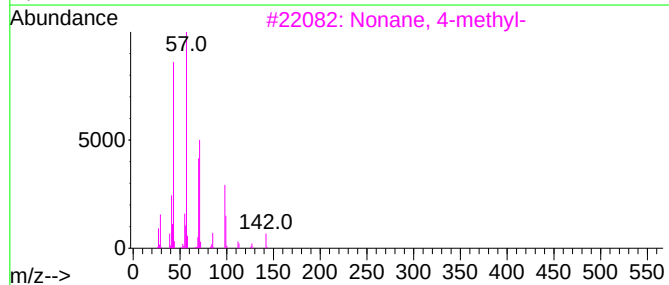
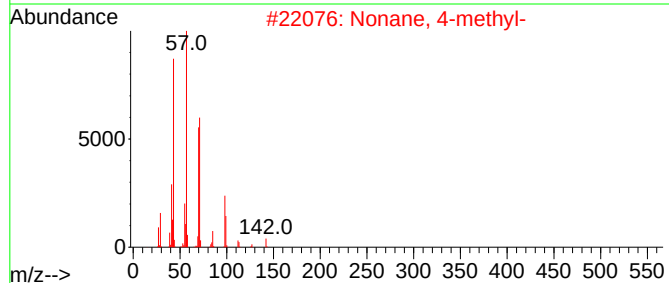
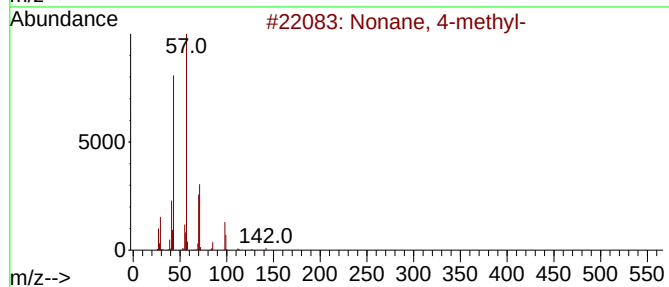
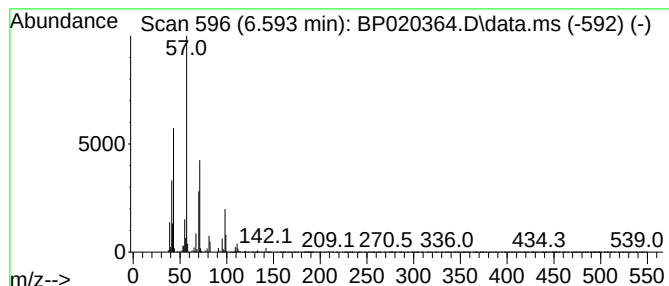
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.593	3.66 ng/ul	110529	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	72
2			Nonane, 4-methyl-	142	C10H22	017301-94-9	59
3			Nonane, 4-methyl-	142	C10H22	017301-94-9	59
4			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	50
5			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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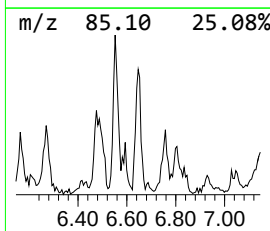
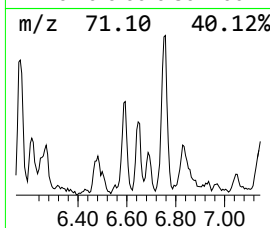
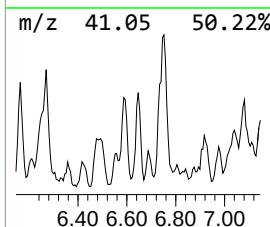
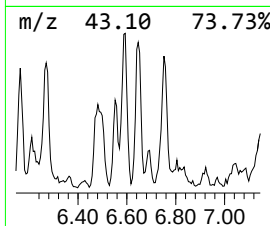
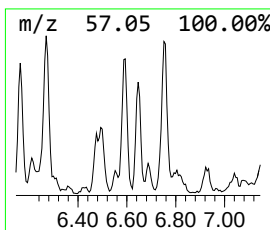
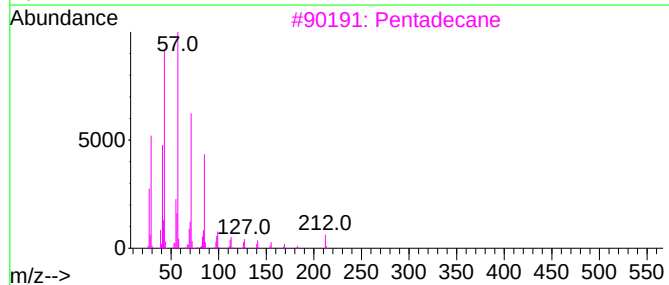
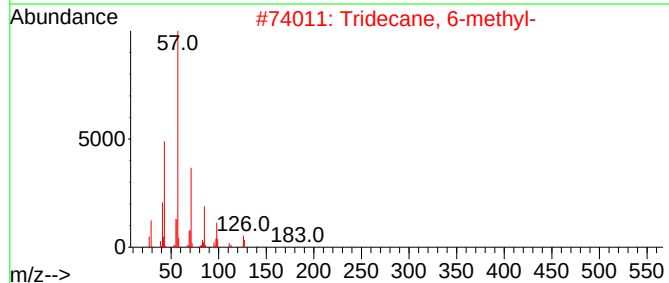
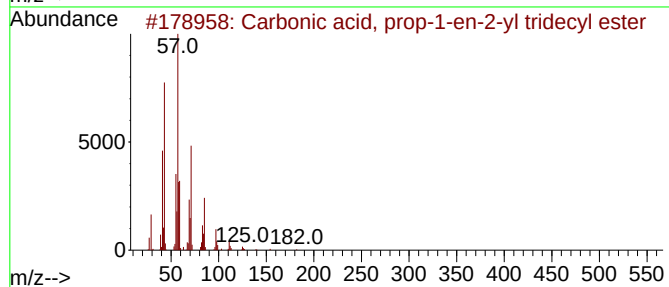
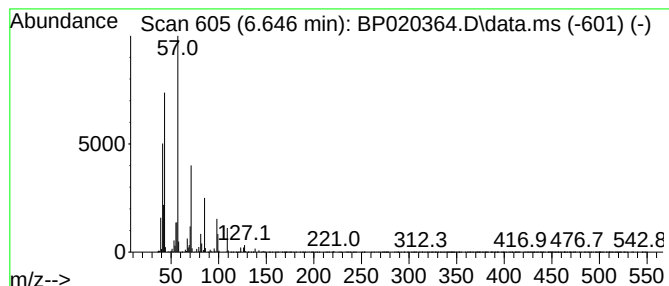
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Carbonic acid, prop-1-en-2-... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.646	2.81 ng/ul	84896	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	59
2			Tridecane, 6-methyl-	198	C14H30	013287-21-3	59
3			Pentadecane	212	C15H32	000629-62-9	59
4			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	58
5			Nonane, 2-methyl-	142	C10H22	000871-83-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

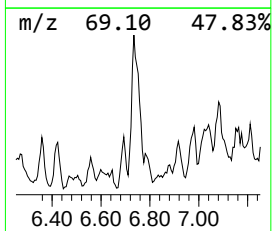
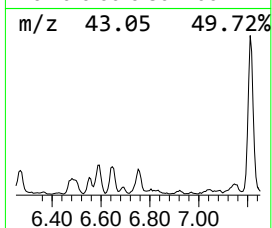
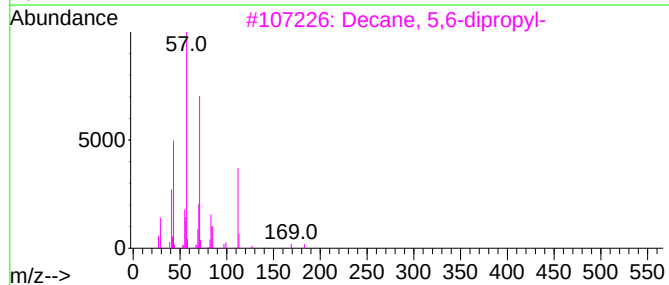
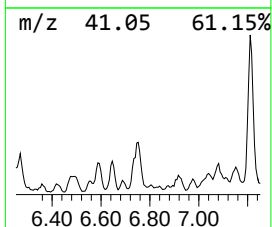
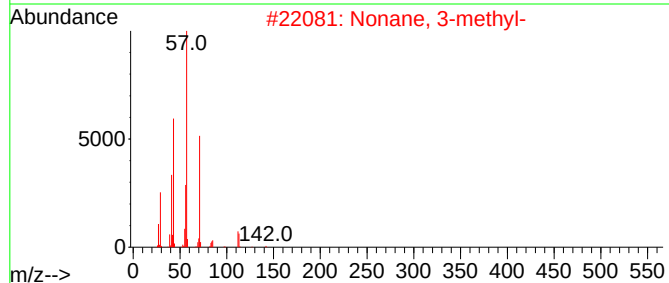
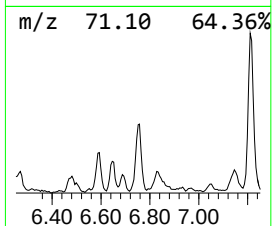
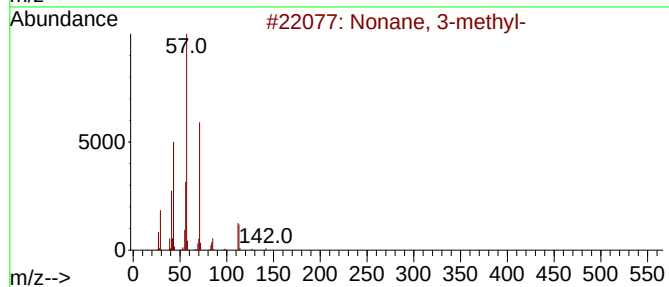
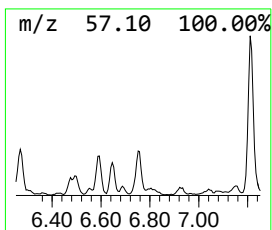
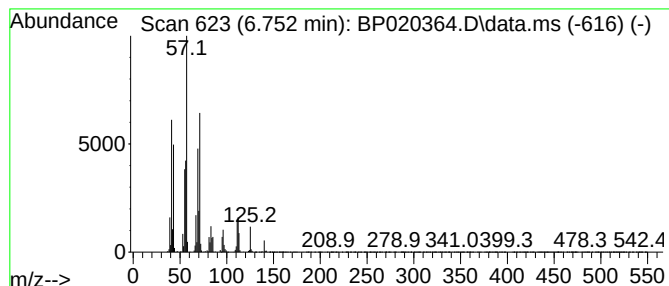
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.752	7.78 ng/ul	234882	1,4-Dichlorobenzene-d4			7.611
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-		142	C10H22	005911-04-6	58
2	Nonane, 3-methyl-		142	C10H22	005911-04-6	52
3	Decane, 5,6-dipropyl-		226	C16H34	119209-20-0	47
4	2-Ethylhexyl mercaptoacetate		204	C10H20O2S	007659-86-1	43
5	Octane, 2,3,7-trimethyl-		156	C11H24	062016-34-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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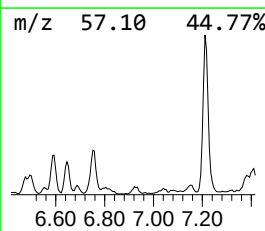
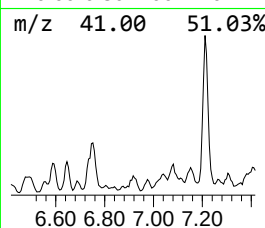
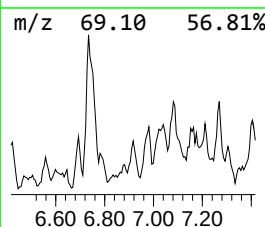
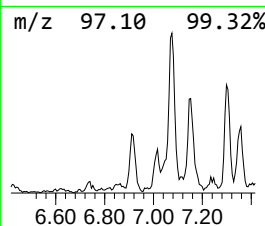
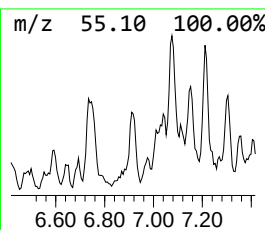
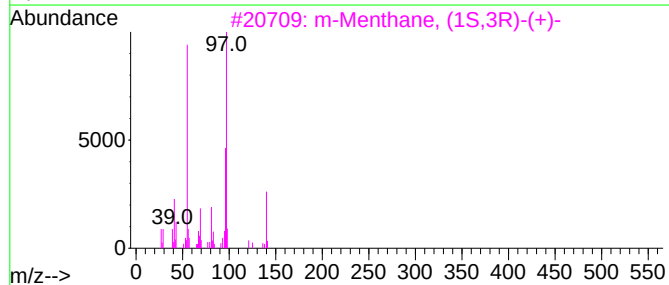
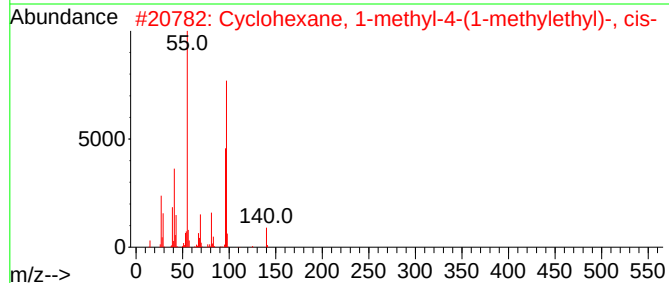
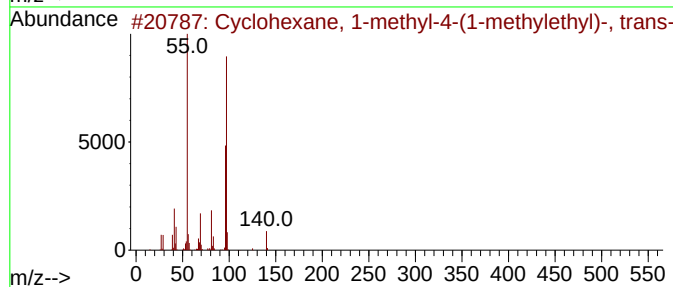
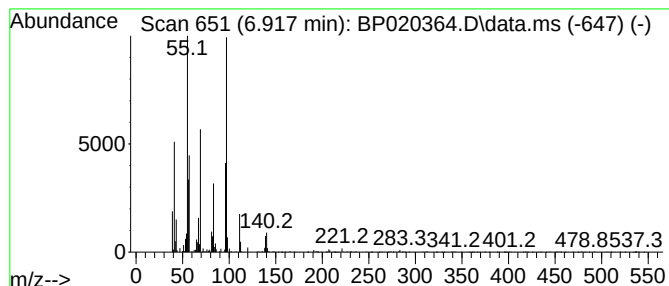
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Cyclic6.917 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.917	2.46 ng/ul	74128	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	64
2			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	64
3			m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	59
4			m-Menthane, (1S,3S)-(+)-	140	C10H20	013837-67-7	59
5			Oxalic acid, cyclohexylmethyl et...	214	C11H18O4	1000309-68-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Sample : P2369-13MEDL 5X
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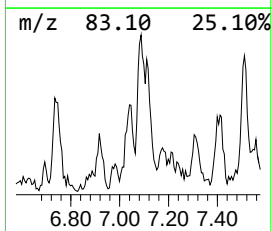
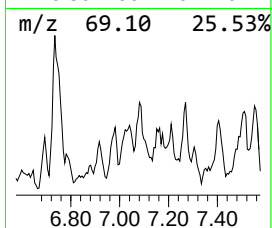
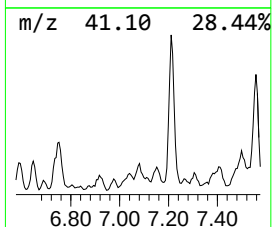
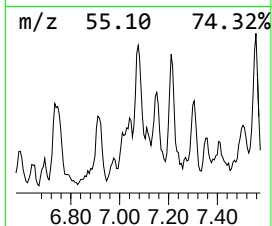
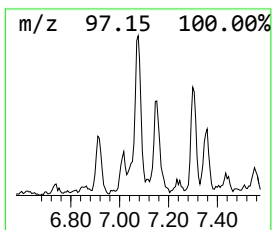
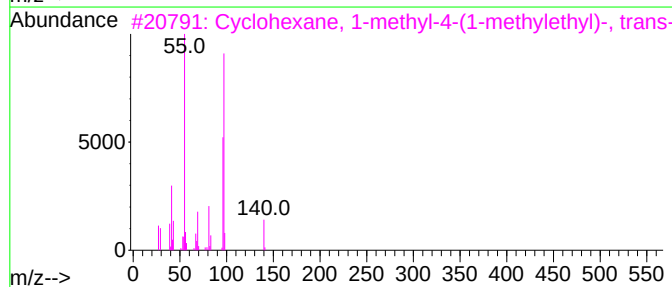
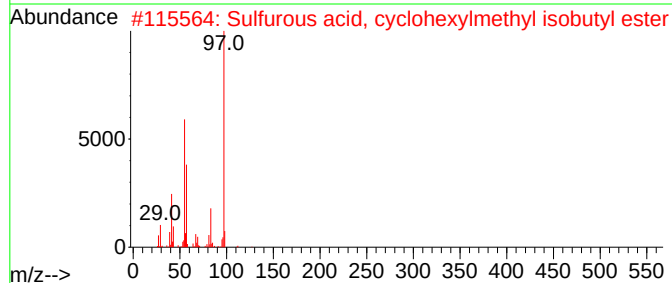
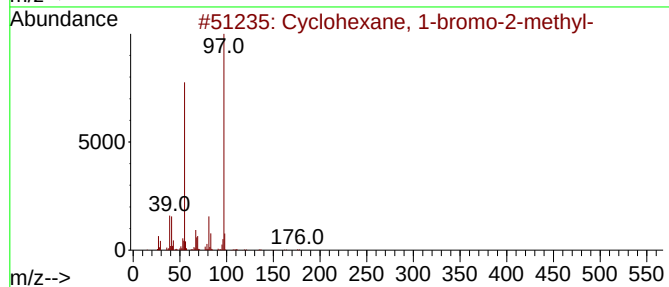
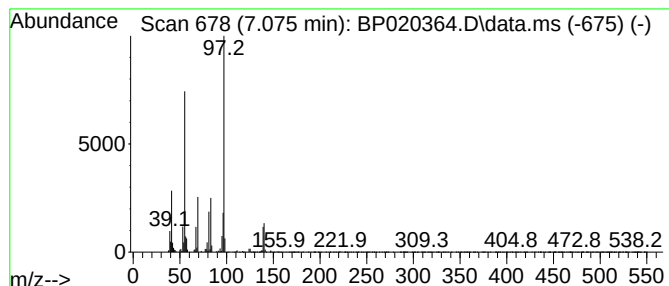
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Cyclohexane, 1-bromo-2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.075	4.12 ng/ul	124411	1,4-Dichlorobenzene-d4	7.611	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-bromo-2-methyl-	176	C7H13Br	006294-39-9	59
2	Sulfurous acid, cyclohexylmethyl...	234	C11H22O3S	1000309-21-3	59
3	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	59
4	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	59
5	Sulfurous acid, butyl cyclohexyl...	234	C11H22O3S	1000309-21-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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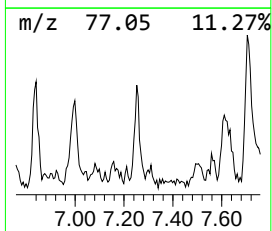
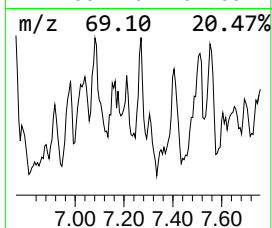
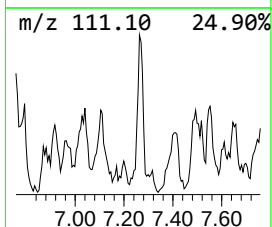
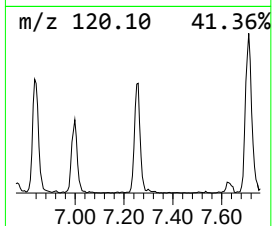
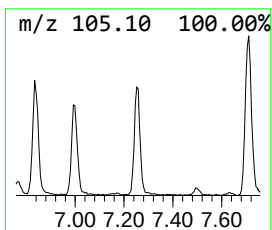
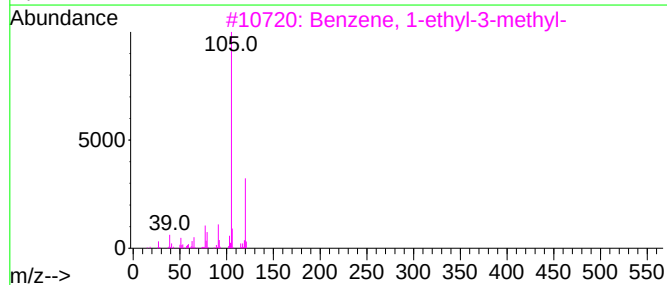
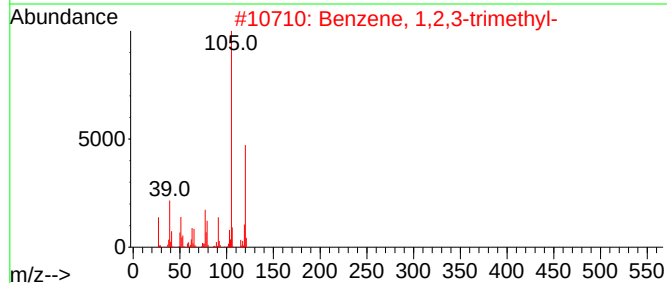
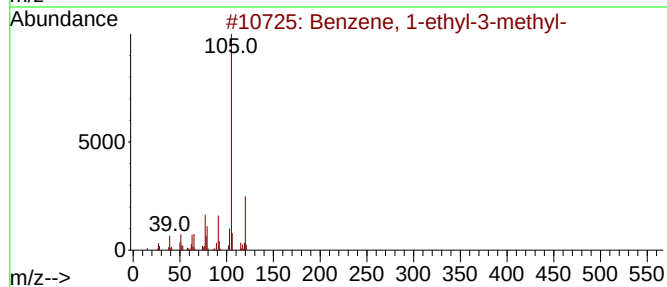
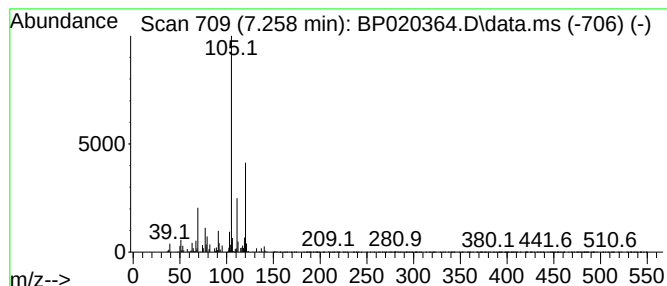
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.258	2.24 ng/ul	67741	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	81
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	76
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N7MEDL

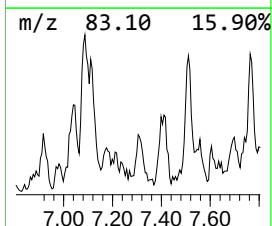
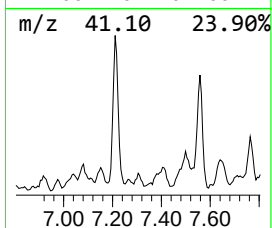
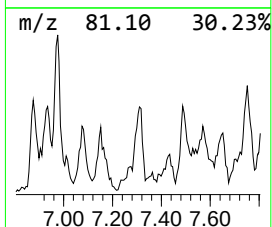
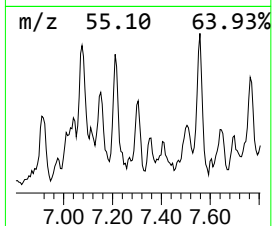
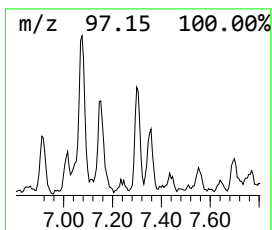
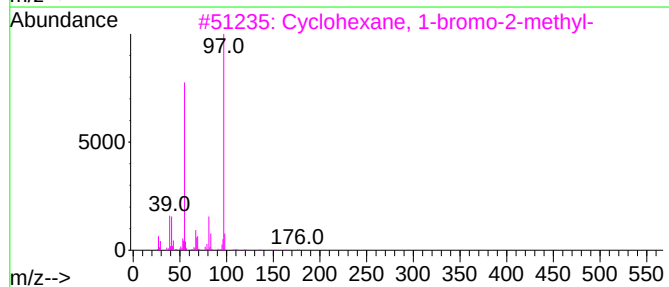
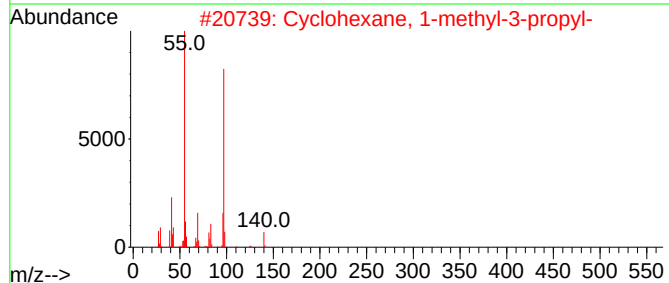
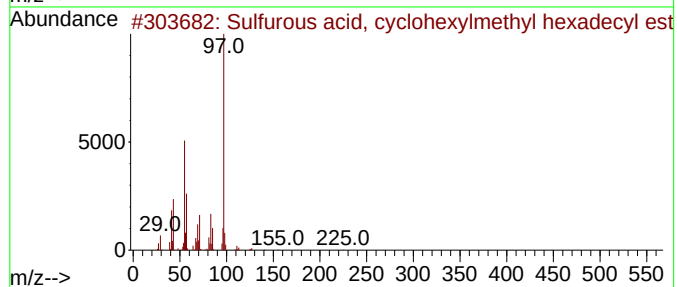
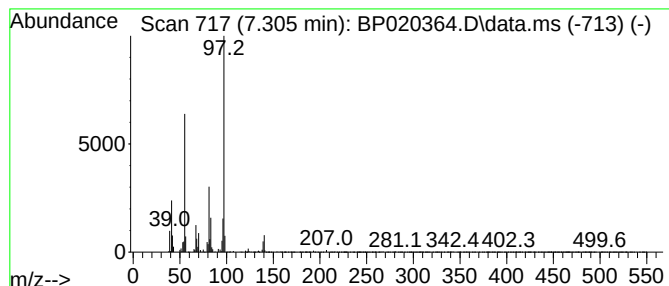
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Sulfurous acid, cyclohexylm... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.305	2.52 ng/ul	76190	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	72
2			Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	72
3			Cyclohexane, 1-bromo-2-methyl-	176	C7H13Br	006294-39-9	72
4			Succinic acid, cyclohexylmethyl ...	315	C18H21NO4	1000389-81-4	72
5			Sulfurous acid, cyclohexylmethyl...	332	C18H36O3S	1000309-21-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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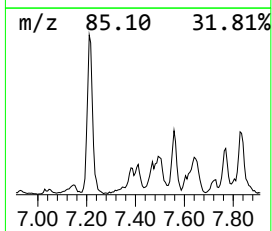
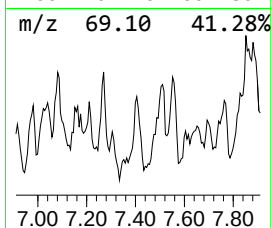
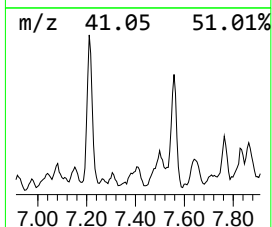
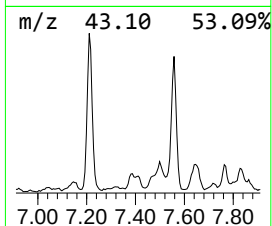
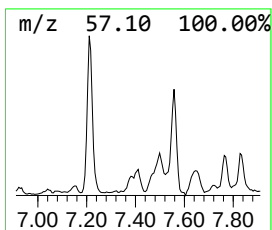
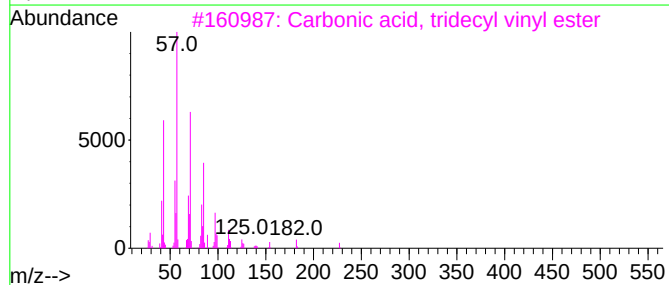
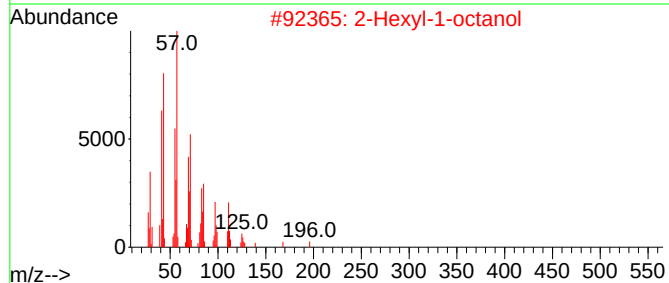
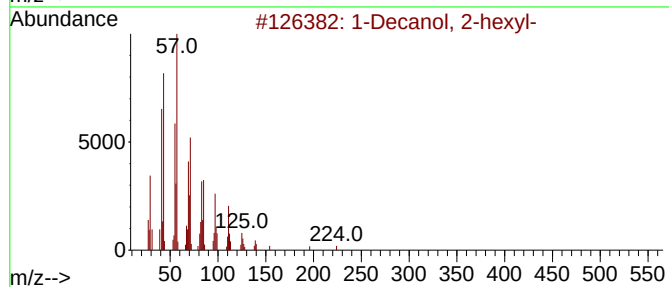
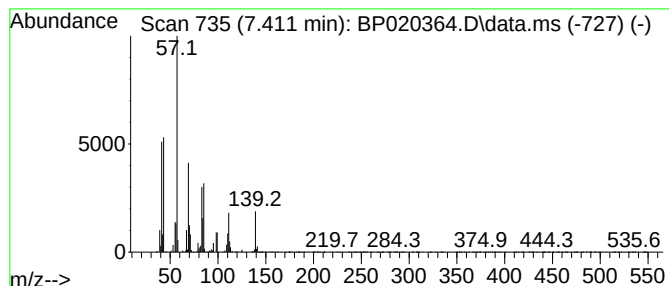
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Decanol, 2-hexyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.411	4.15 ng/ul	125231	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-			242	C16H34O	002425-77-6	83
2	2-Hexyl-1-octanol			214	C14H30O	019780-79-1	72
3	Carbonic acid, tridecyl vinyl ester			270	C16H30O3	1000382-54-7	59
4	Carbonic acid, eicosyl vinyl ester			368	C23H44O3	1000382-54-3	53
5	Hexadecane, 1-chloro-			260	C16H33Cl	004860-03-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Sample : P2369-13MEDL 5X
Misc :
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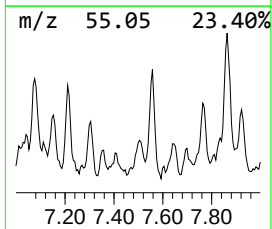
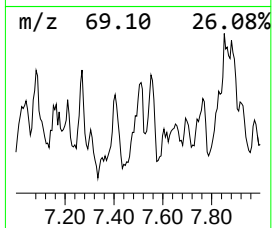
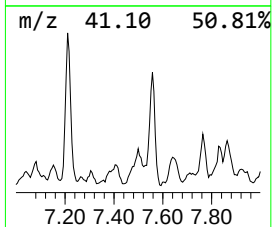
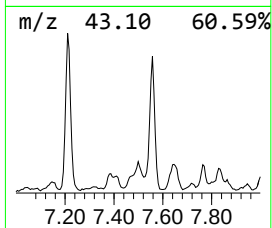
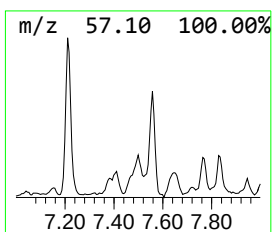
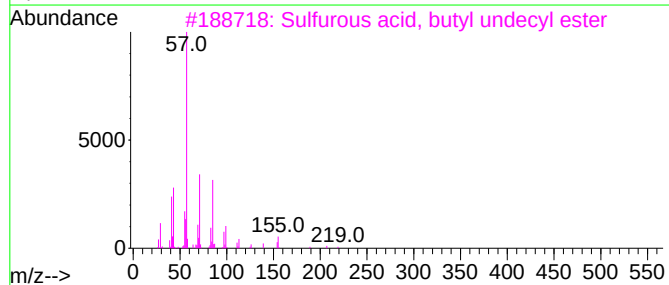
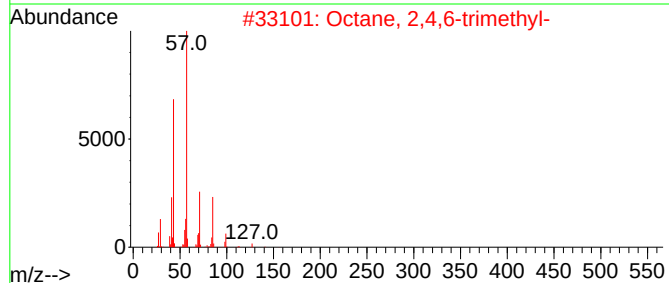
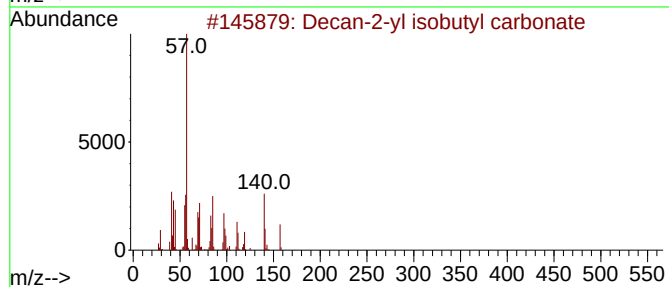
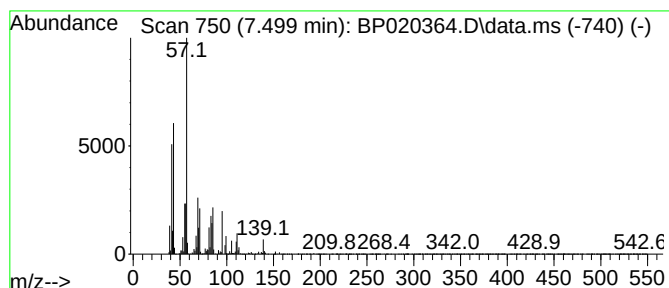
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Decan-2-yl isobutyl carbonate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.499	8.03 ng/ul	242334	1,4-Dichlorobenzene-d4	7.611	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decan-2-yl isobutyl carbonate	258	C15H30O3	1000372-79-5	59
2	Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	47
3	Sulfurous acid, butyl undecyl ester	292	C15H32O3S	1000309-17-8	47
4	Hexyl octyl ether	214	C14H30O	017071-54-4	43
5	Nonane	128	C9H20	000111-84-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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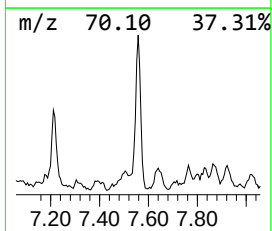
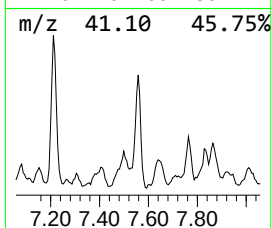
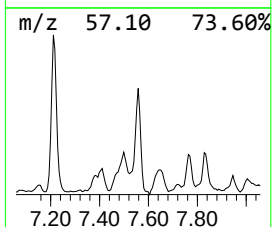
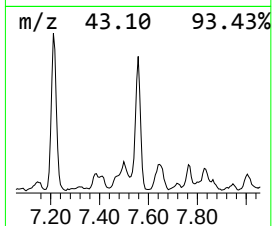
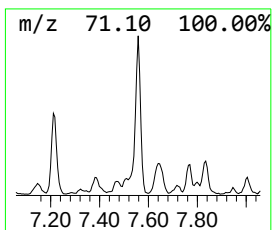
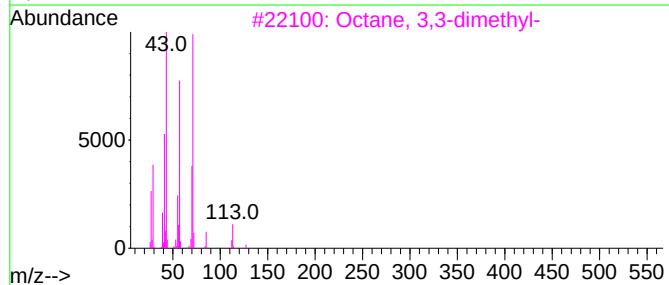
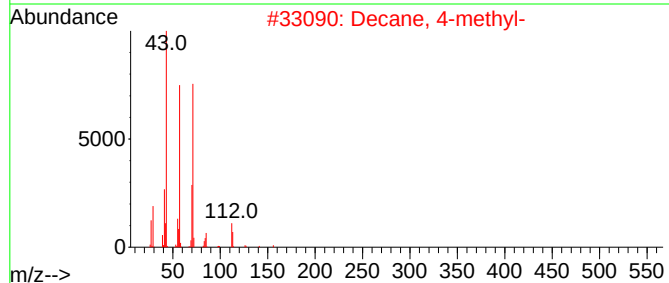
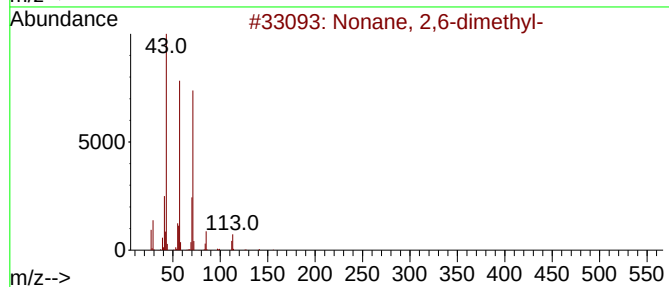
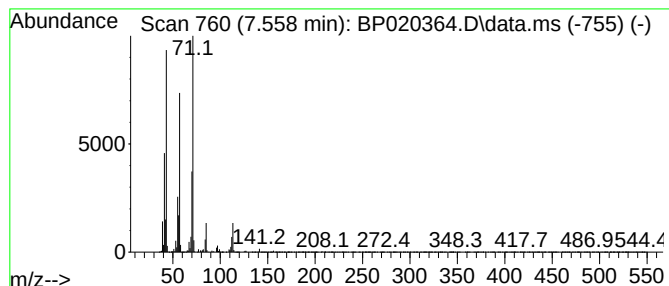
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.558	13.34 ng/ul	402590	1,4-Dichlorobenzene-d4	7.611

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	91
2		Decane, 4-methyl-	156	C11H24	002847-72-5	91
3		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	78
4		Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	78
5		Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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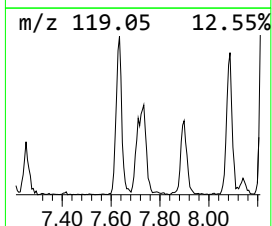
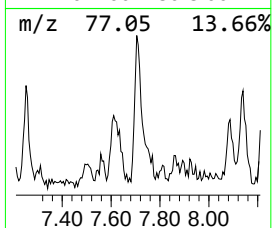
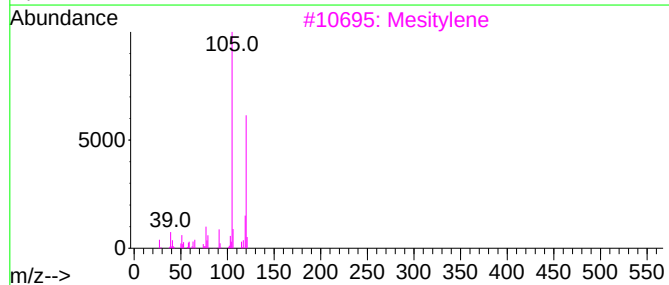
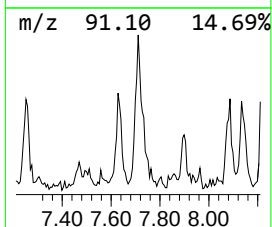
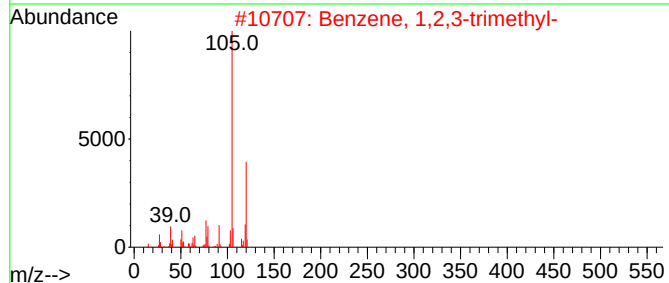
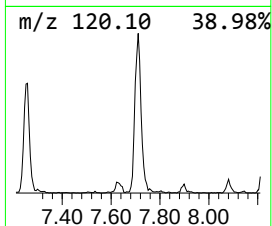
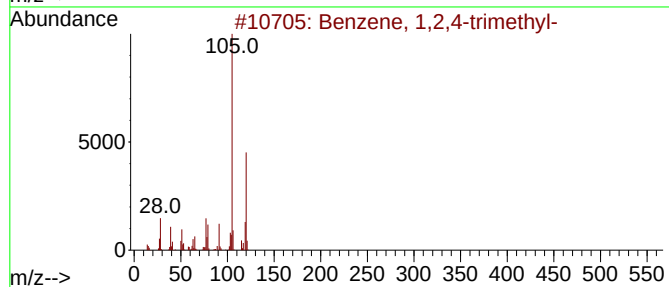
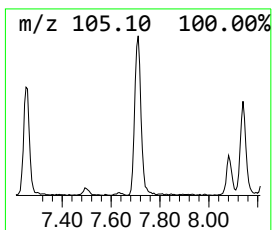
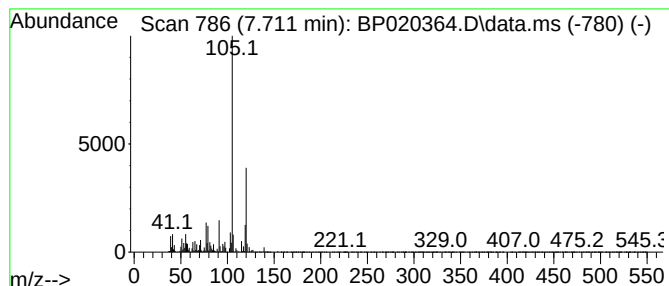
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 1,2,4-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.711	6.05 ng/ul	182548	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3			Mesitylene	120	C9H12	000108-67-8	93
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
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Misc :
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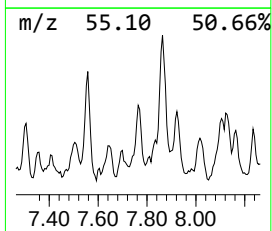
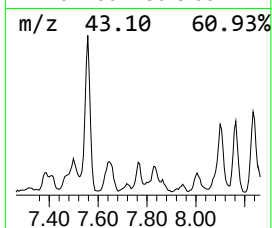
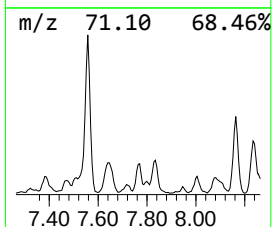
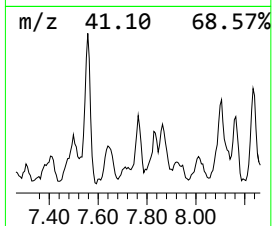
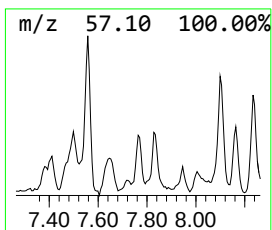
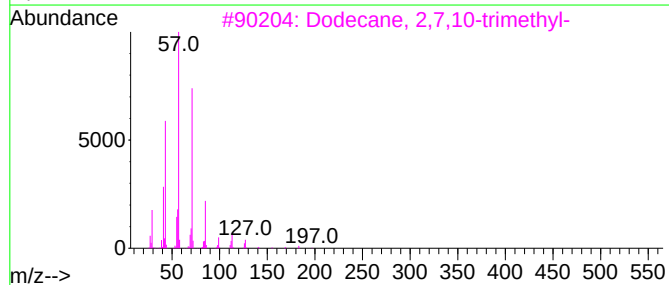
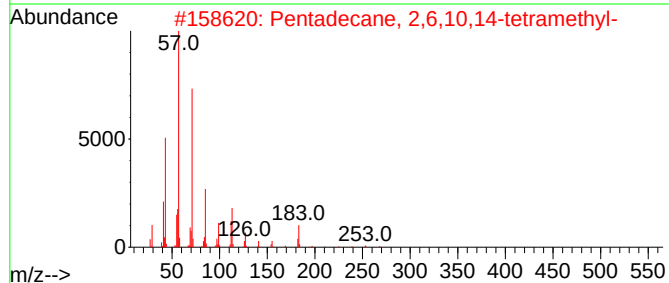
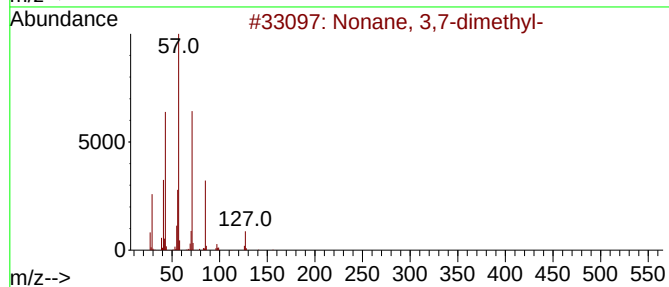
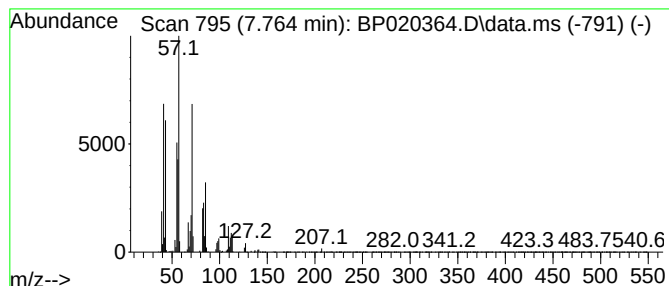
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.764	4.99 ng/ul	150575	1,4-Dichlorobenzene-d4	7.611

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	46
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	46
3		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	43
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	43
5		Undecane, 3-methyl-	170	C12H26	001002-43-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Operator : MA/JU
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Misc :
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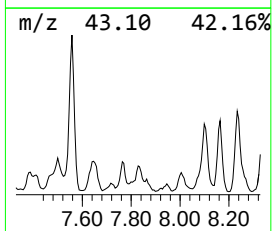
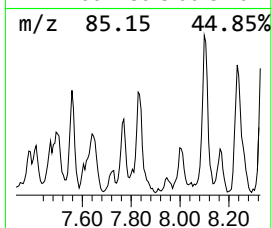
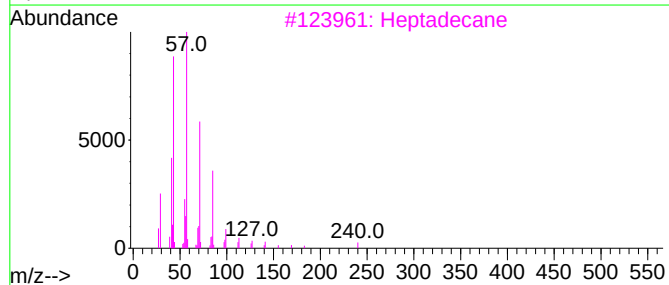
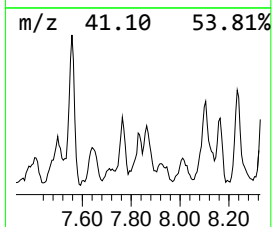
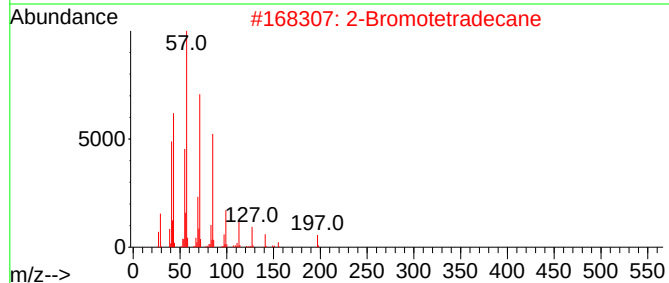
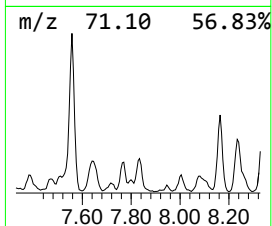
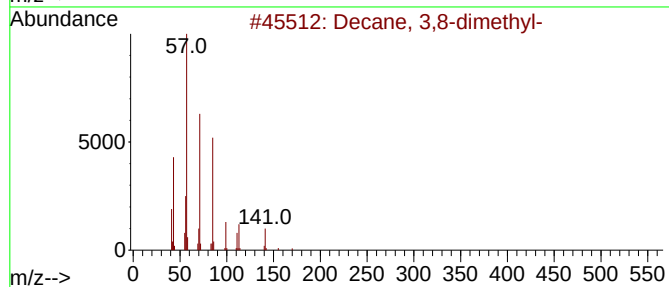
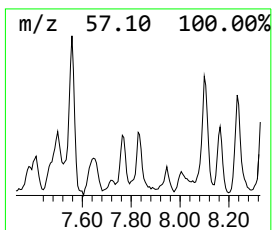
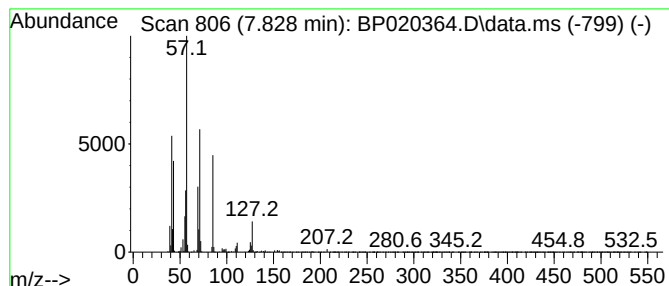
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	4.06 ng/ul	122568	1,4-Dichlorobenzene-d4	7.611

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
2		2-Bromotetradecane	276	C14H29Br	074036-95-6	64
3		Heptadecane	240	C17H36	000629-78-7	64
4		Dodecane, 3-methyl-	184	C13H28	017312-57-1	64
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

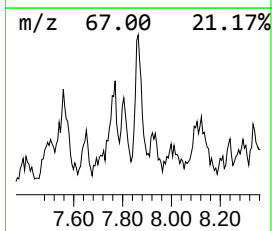
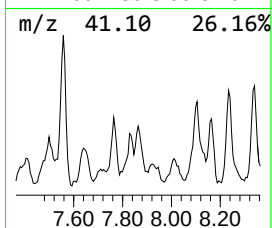
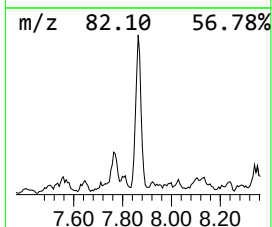
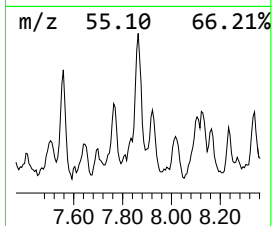
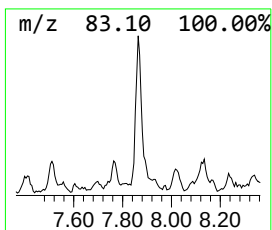
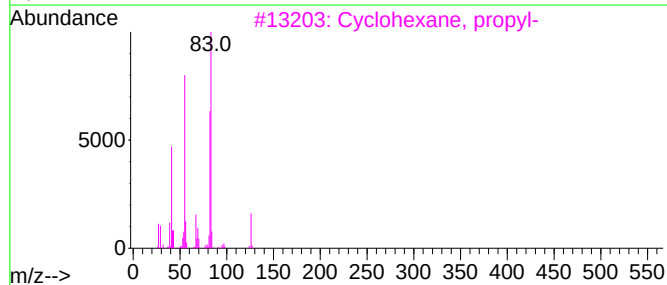
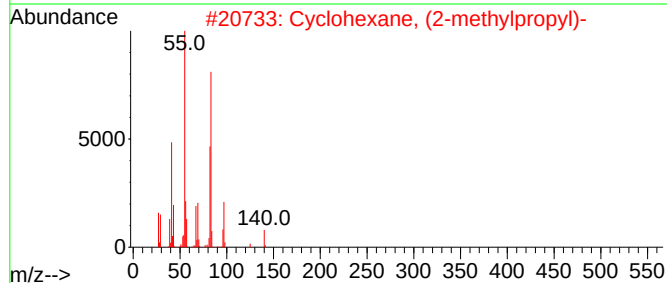
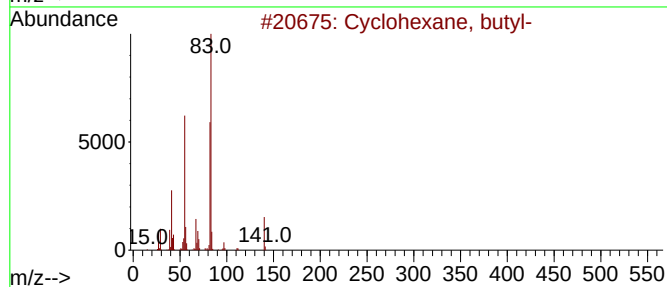
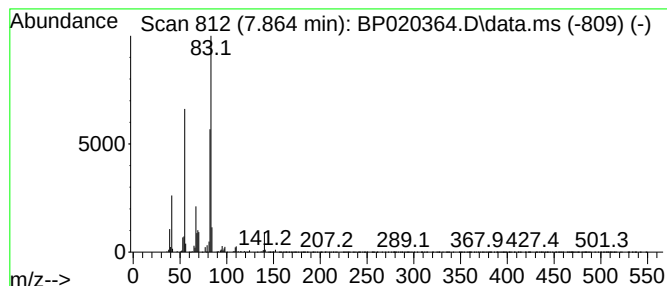
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Cyclic7.864 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.864	5.36 ng/ul	161786	1,4-Dichlorobenzene-d4			7.611
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane, butyl-		140	C10H20	001678-93-9	78
2	Cyclohexane, (2-methylpropyl)-		140	C10H20	001678-98-4	72
3	Cyclohexane, propyl-		126	C9H18	001678-92-8	72
4	Cyclohexane, (2-methylpropyl)-		140	C10H20	001678-98-4	64
5	Succinic acid, 2-methylpent-3-yl...		284	C16H28O4	1000391-11-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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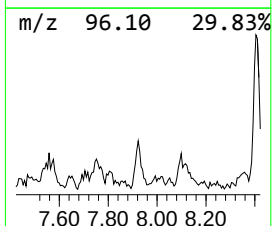
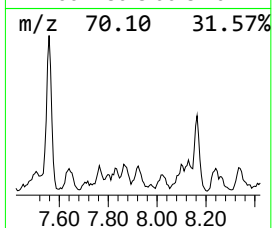
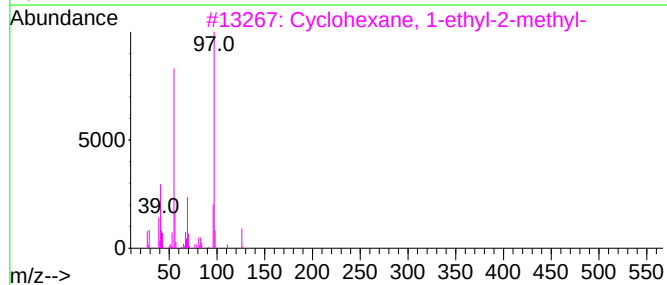
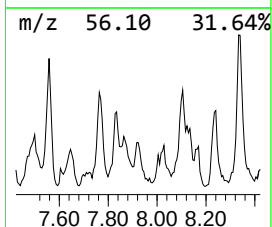
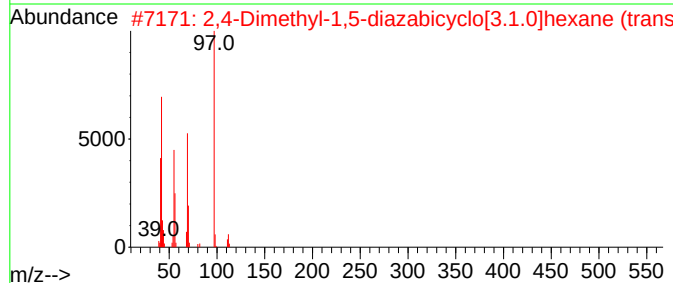
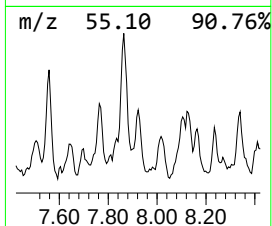
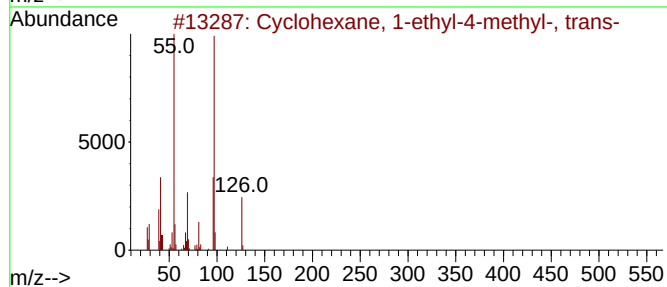
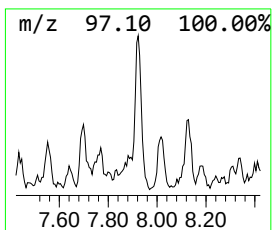
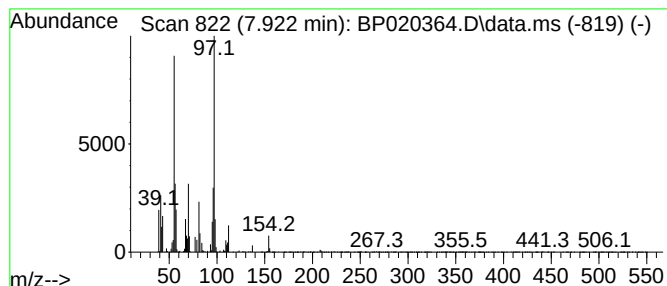
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Cyclic7.922 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.922	3.01 ng/ul	90729	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	59
2			2,4-Dimethyl-1,5-diazabicyclo[3....	112	C6H12N2	1000283-20-9	59
3			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	59
4			Succinic acid, 2-ethylhexyl tran...	326	C19H34O4	1000390-07-1	53
5			Oxalic acid, cyclohexylmethyl pr...	228	C12H20O4	1010309-68-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

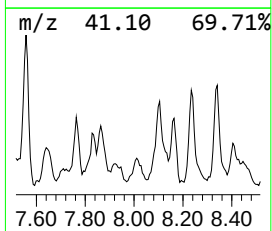
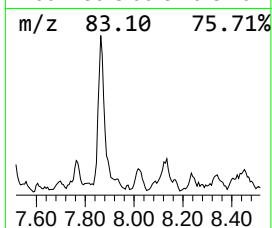
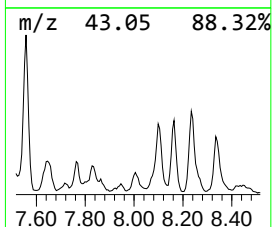
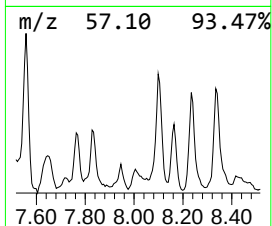
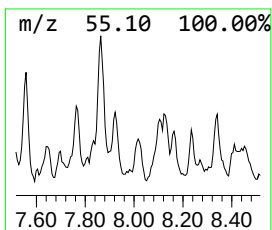
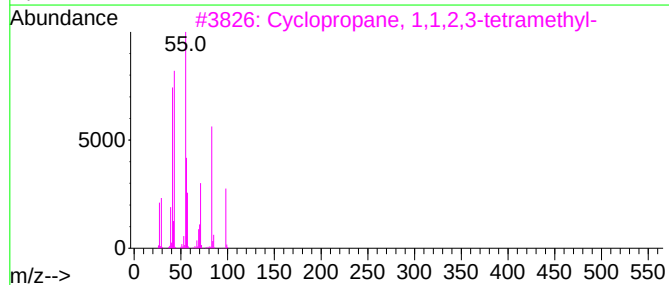
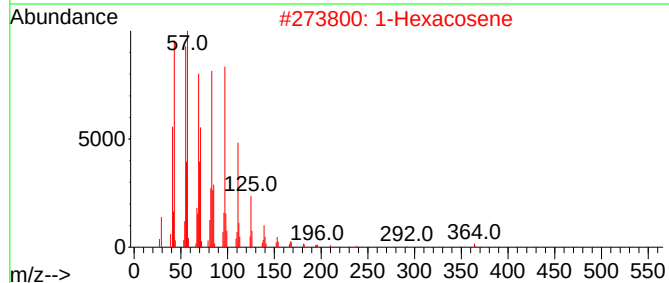
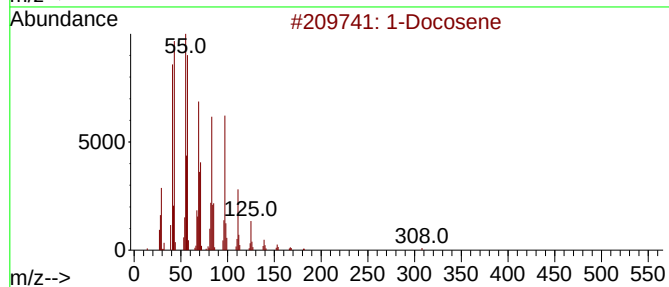
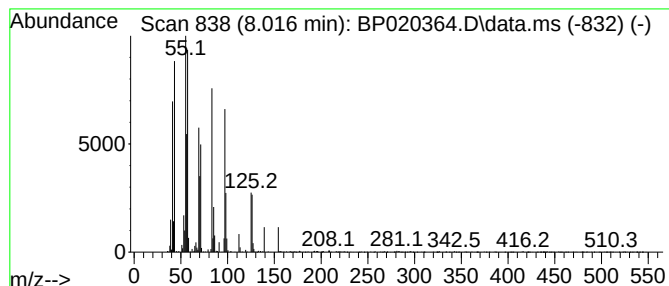
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.016	3.35 ng/ul	101234	1,4-Dichlorobenzene-d4	7.611	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	72
2	1-Hexacosene	364	C26H52	018835-33-1	64
3	Cyclopropane, 1,1,2,3-tetramethyl-	98	C7H14	074752-93-5	50
4	Cyclohexanone, 3,3,5,5-tetramethyl-	154	C10H18O	014376-79-5	43
5	Cyclohexanone, 3-ethyl-	126	C8H14O	022461-89-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

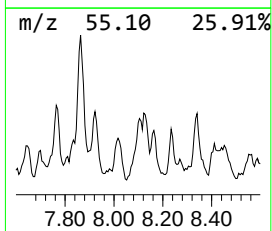
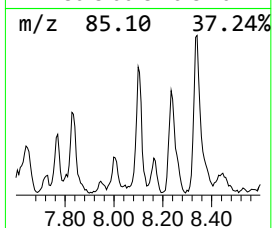
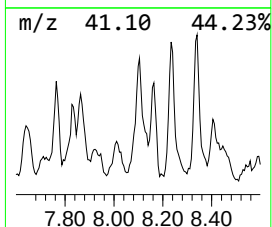
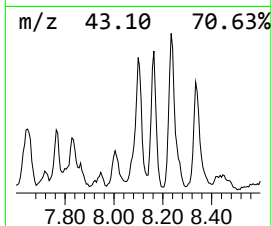
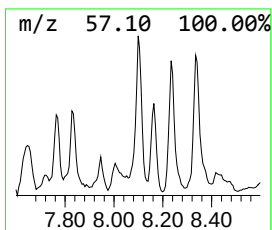
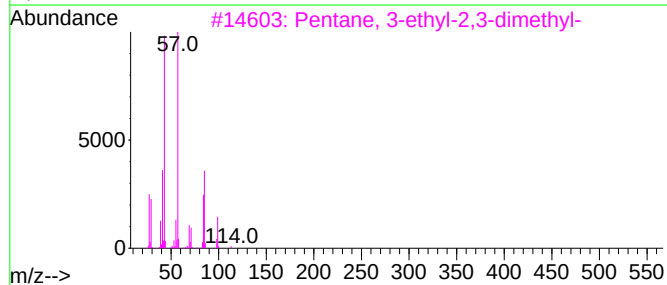
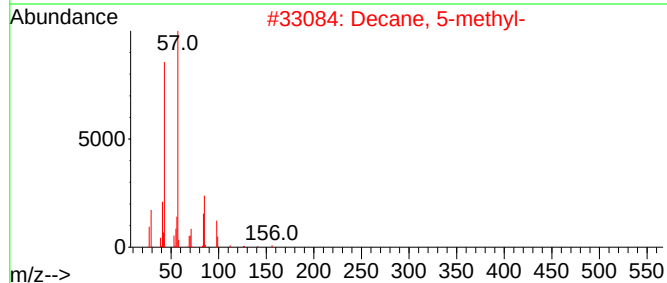
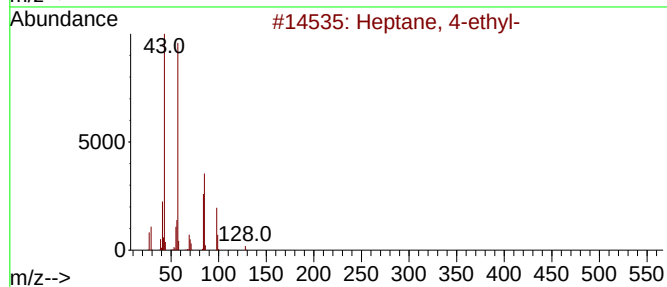
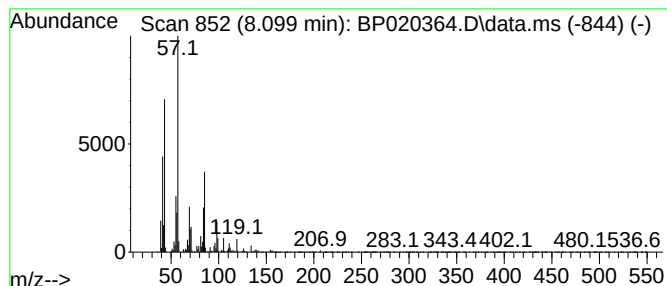
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.099	11.69 ng/ul	352734	1,4-Dichlorobenzene-d4	7.611	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 4-ethyl-	128	C9H20	002216-32-2	53
2	Decane, 5-methyl-	156	C11H24	013151-35-4	53
3	Pentane, 3-ethyl-2,3-dimethyl-	128	C9H20	016747-33-4	53
4	Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	43
5	Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
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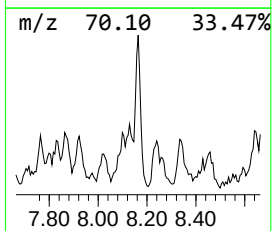
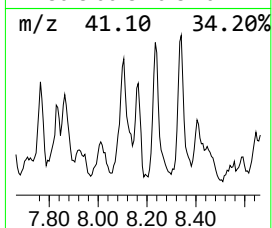
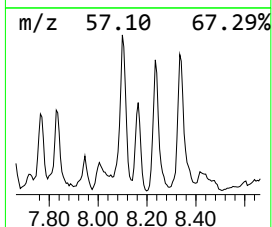
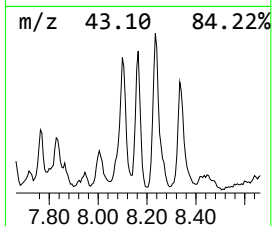
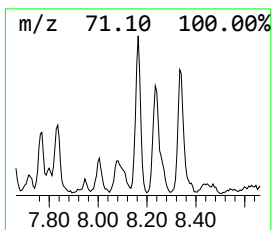
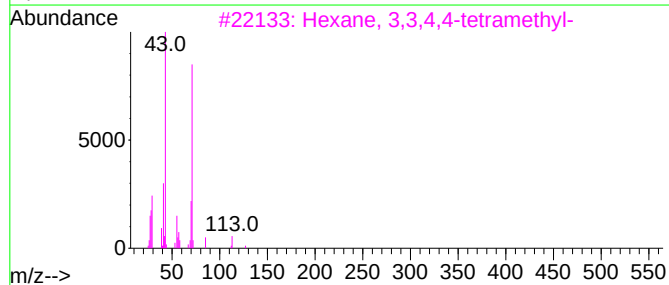
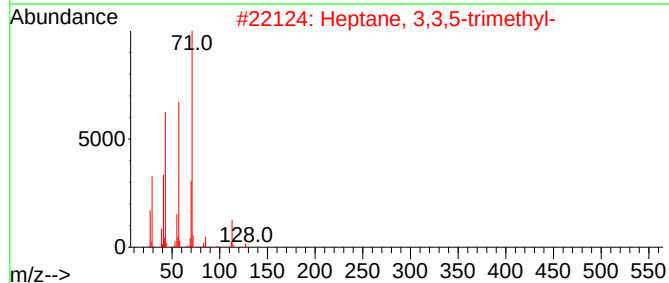
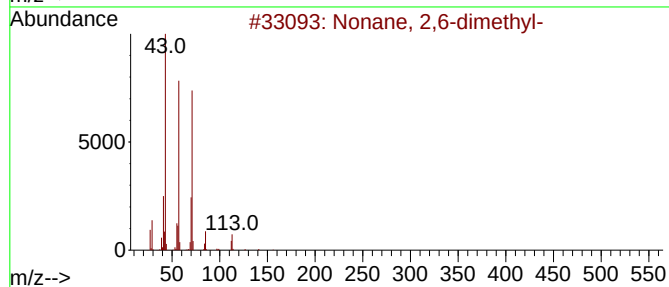
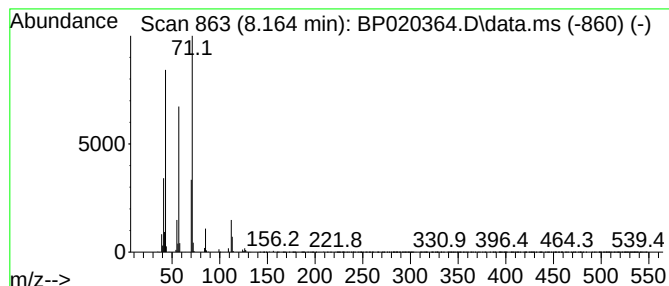
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.164	6.22 ng/ul	187832	1,4-Dichlorobenzene-d4	7.611

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	83
2		Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72
3		Hexane, 3,3,4,4-tetramethyl-	142	C10H22	005171-84-6	72
4		4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	64
5		Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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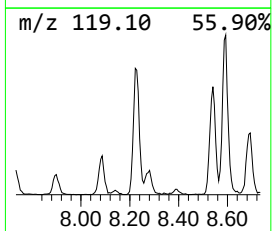
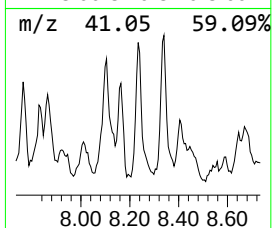
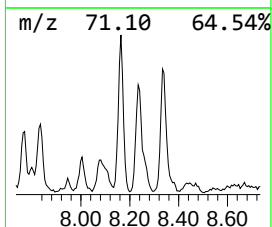
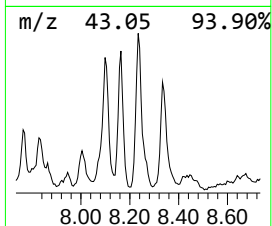
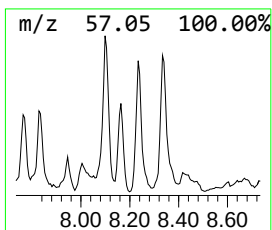
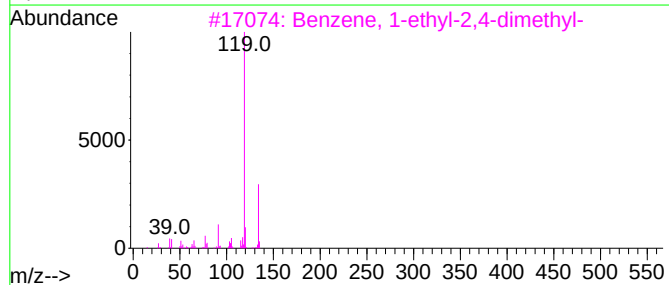
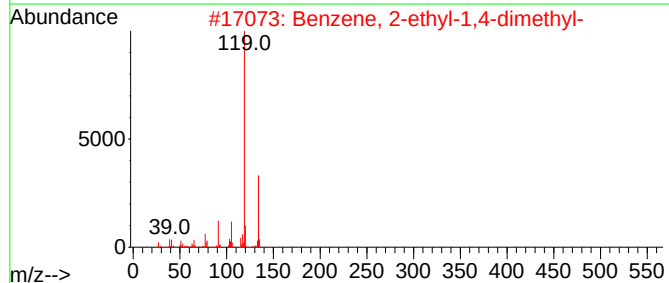
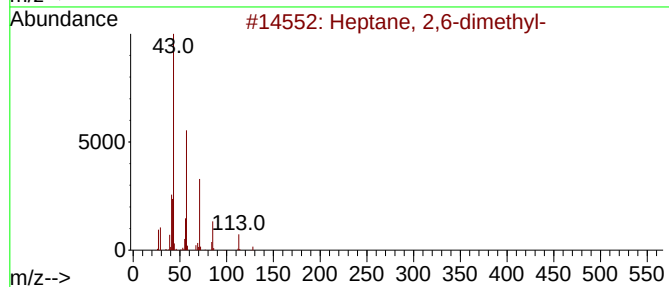
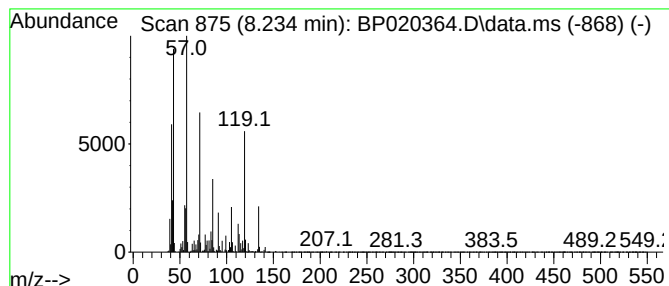
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	12.70 ng/ul	383096	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	49
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	35
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	35
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	35
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
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Acq On : 13 May 2024 22:13
Operator : MA/JU
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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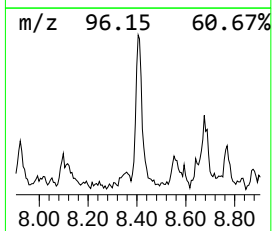
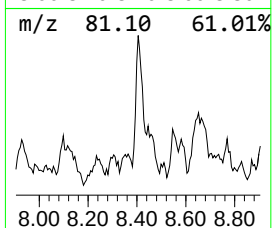
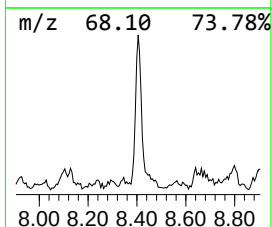
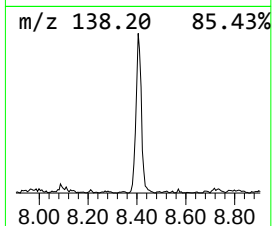
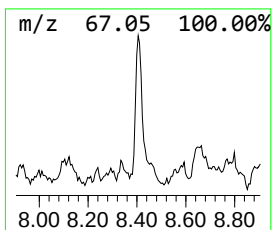
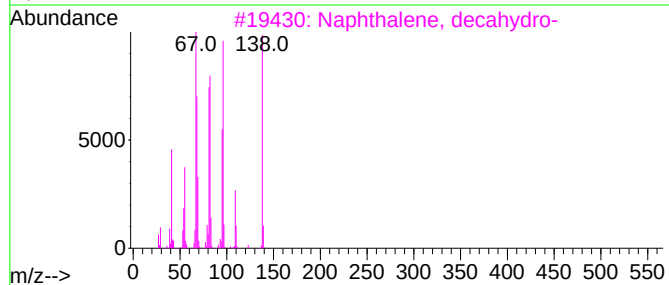
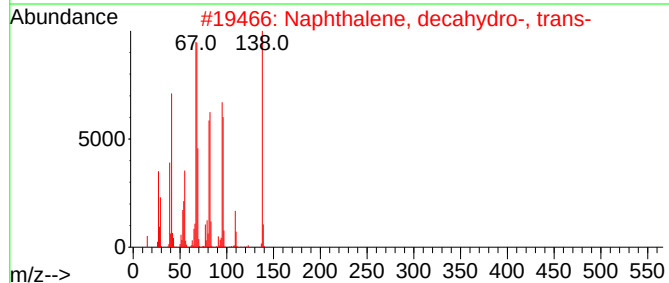
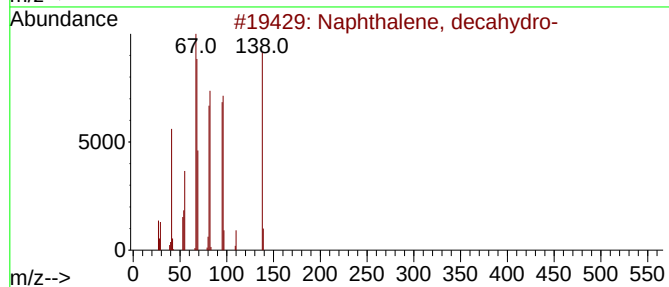
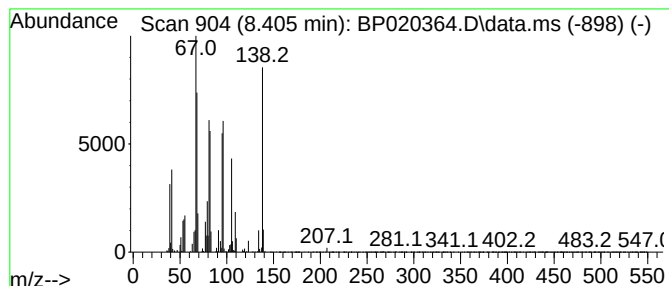
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 Naphthalene, decahydro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.405	7.16 ng/ul	216098	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-	138	C10H18	000091-17-8	93
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	93
3			Naphthalene, decahydro-	138	C10H18	000091-17-8	93
4			Naphthalene, decahydro-	138	C10H18	000091-17-8	93
5			Naphthalene, decahydro-	138	C10H18	000091-17-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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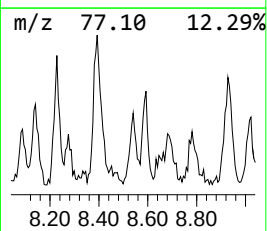
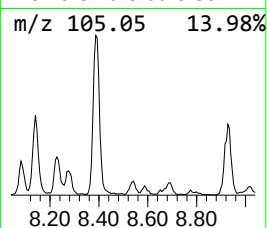
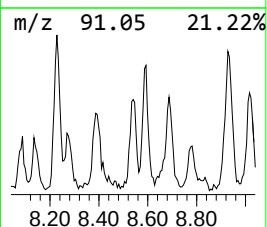
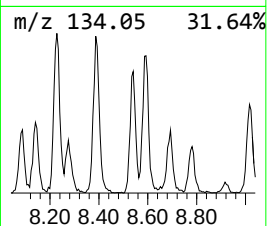
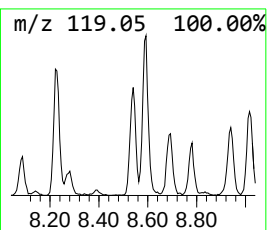
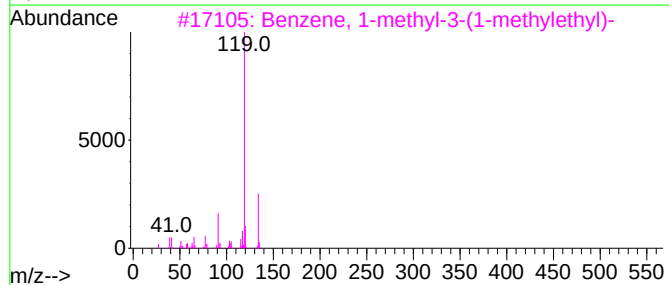
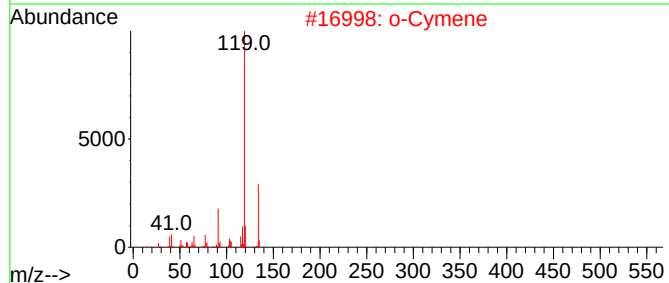
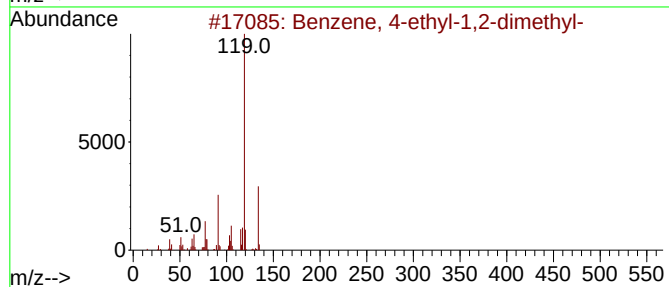
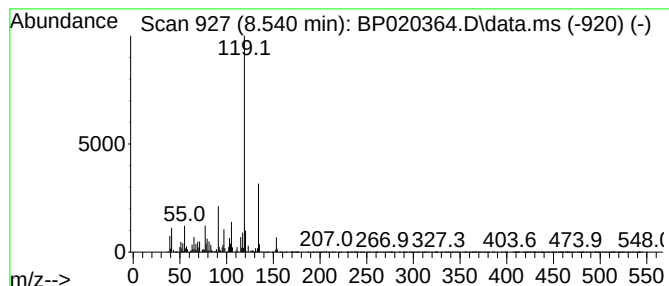
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.540	4.59 ng/ul	138550	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			o-Cymene	134	C10H14	000527-84-4	94
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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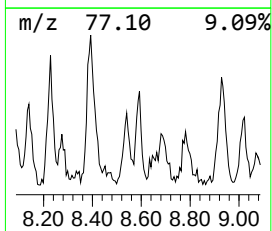
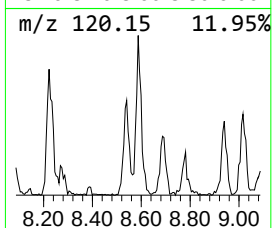
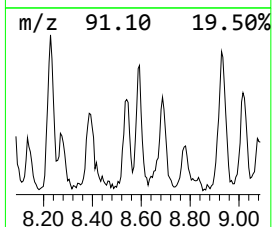
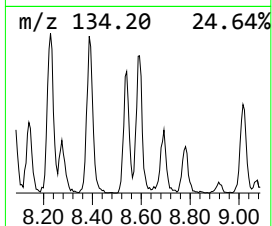
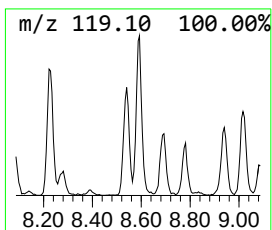
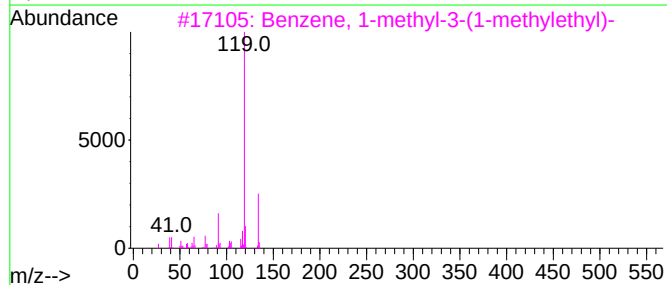
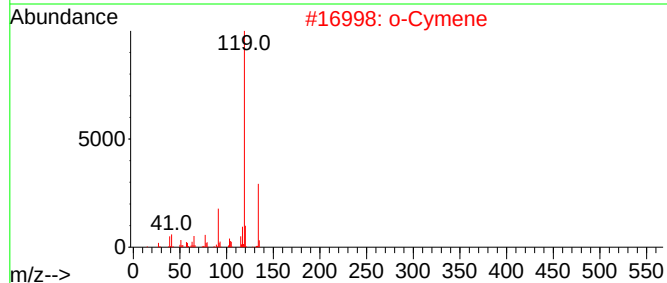
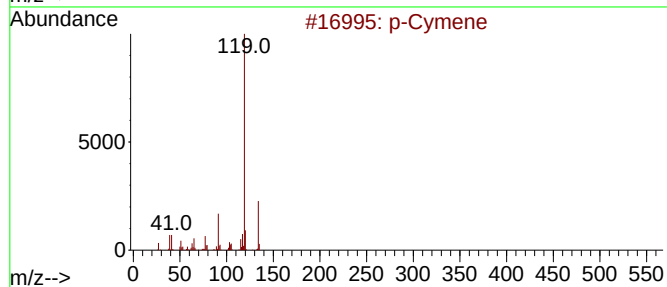
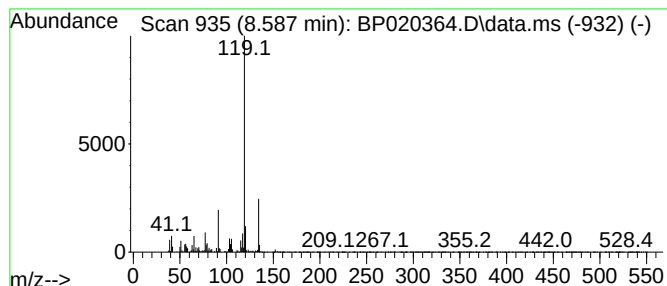
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 p-Cymene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.587	3.20 ng/ul	96706	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4			o-Cymene	134	C10H14	000527-84-4	97
5			o-Cymene	134	C10H14	000527-84-4	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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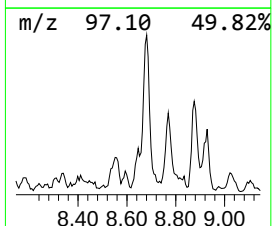
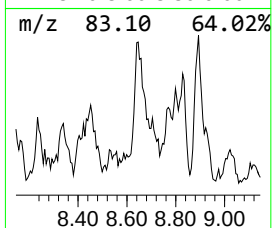
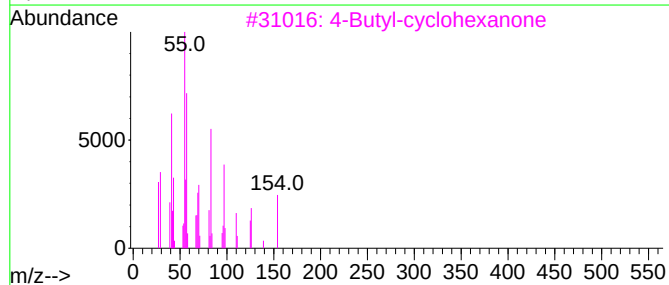
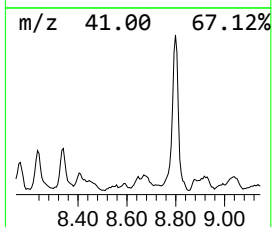
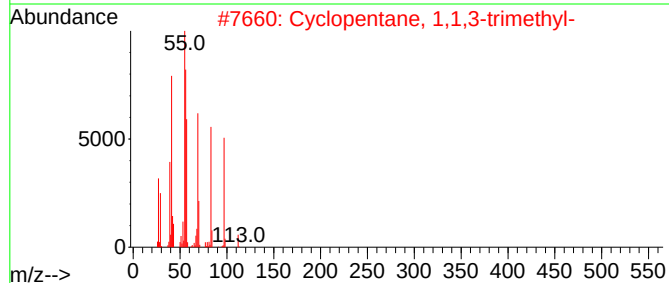
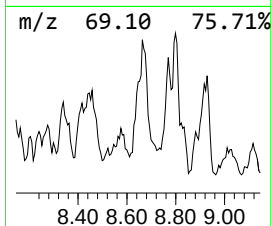
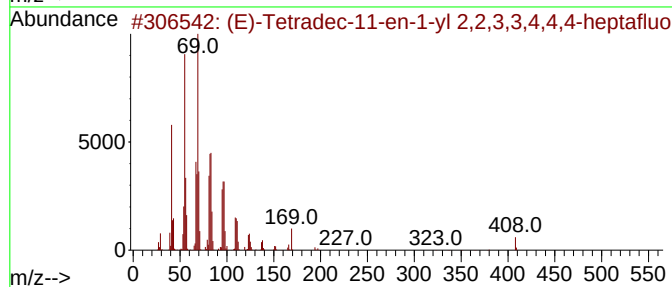
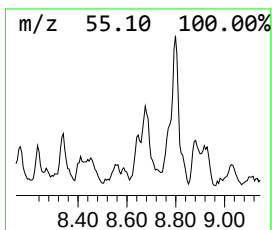
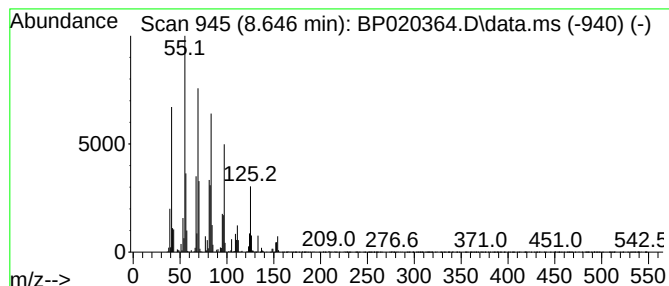
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 (E)-Tetradec-11-en-1-yl 2,2... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.646	2.67 ng/ul	80439	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-Tetradec-11-en-1-yl	2,2,3,3,...	408	C18H27F702	1000385-85-1	58	
2	Cyclopentane, 1,1,3-trimethyl-		112	C8H16	004516-69-2	58	
3	4-Butyl-cyclohexanone		154	C10H18O	061203-82-5	43	
4	Cyclopentane, 1,1,3-trimethyl-		112	C8H16	004516-69-2	42	
5	(2-Methylbutyl)cyclohexane		154	C11H22	054105-77-0	38	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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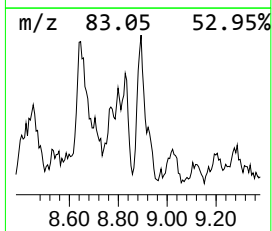
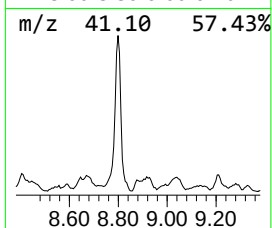
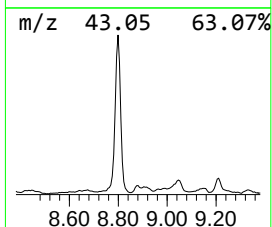
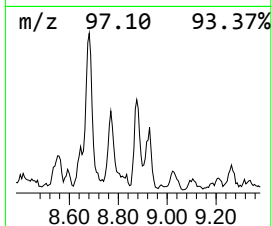
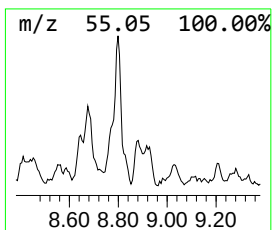
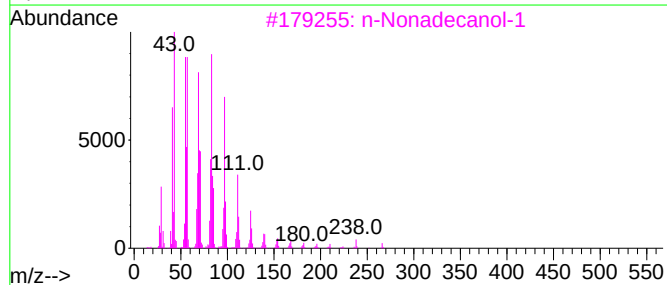
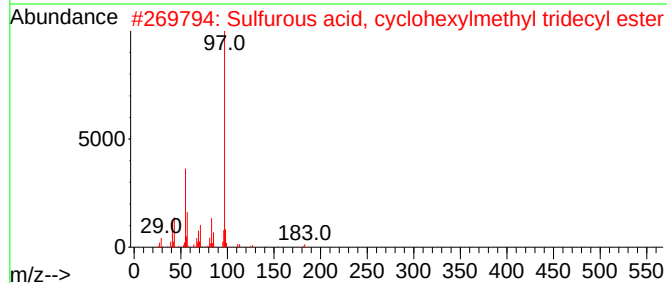
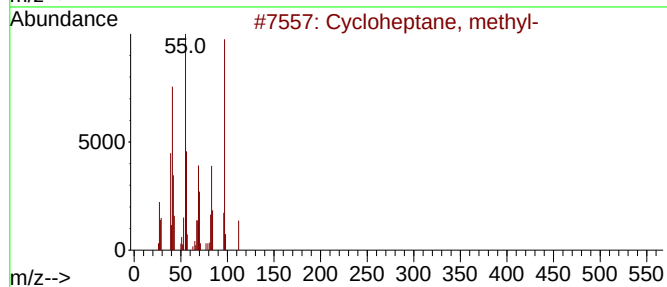
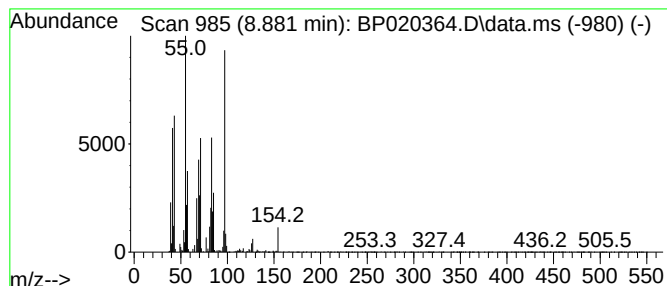
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 (DEL) Alkane: Cyclic8.881 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.881	3.52 ng/ul	106347	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloheptane, methyl-	112	C8H16	004126-78-7	47
2			Sulfurous acid, cyclohexylmethyl...	360	C20H40O3S	1000309-22-1	47
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	38
4			n-Pentadecanol	228	C15H32O	000629-76-5	38
5			1-Hexadecanol	242	C16H34O	036653-82-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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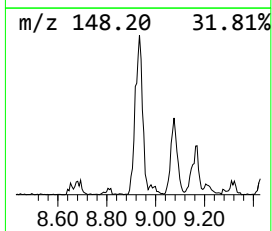
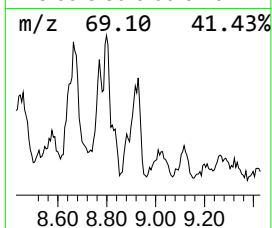
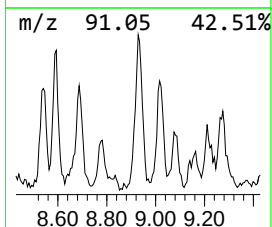
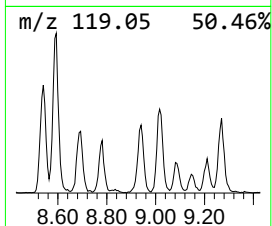
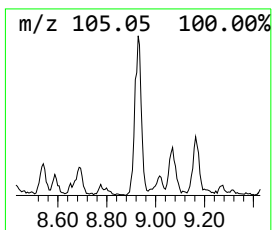
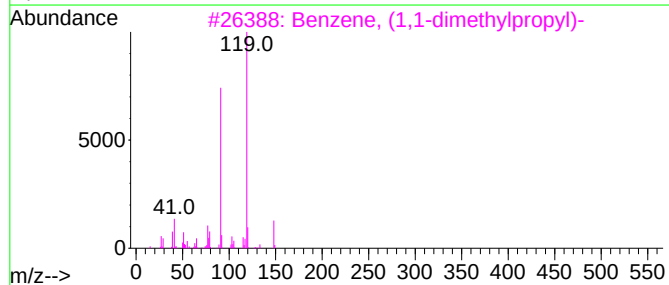
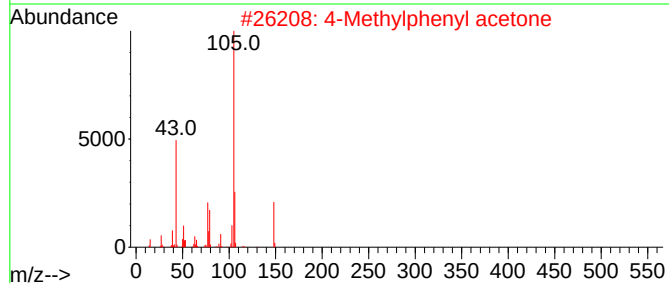
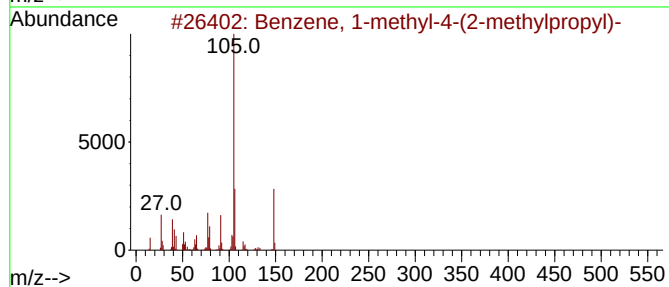
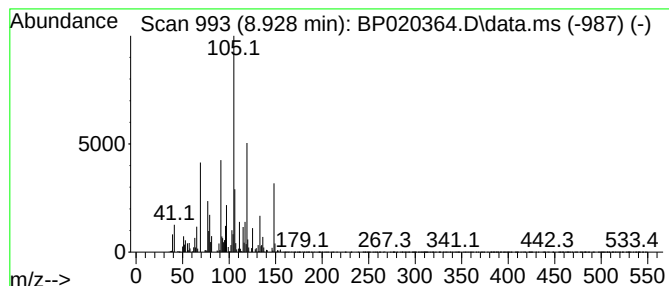
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	8.31 ng/ul	250863	1,4-Dichlorobenzene-d4	7.611

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	38
2			4-Methylphenyl acetone	148	C10H12O	002096-86-8	38
3			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	30
4			6,6-DIMETHYL-2-VINYLBICYCLO[3.1...	148	C11H16	000473-00-7	27
5			Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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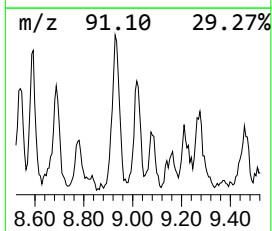
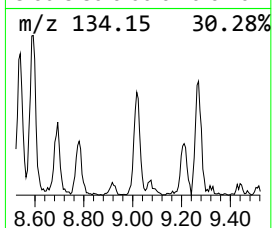
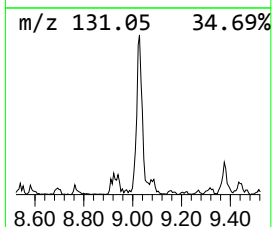
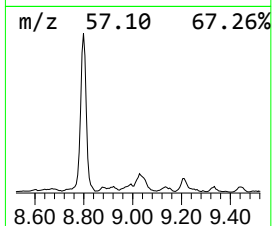
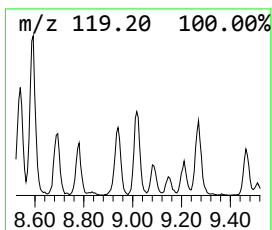
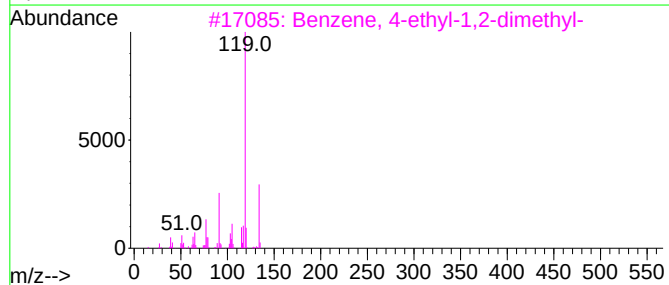
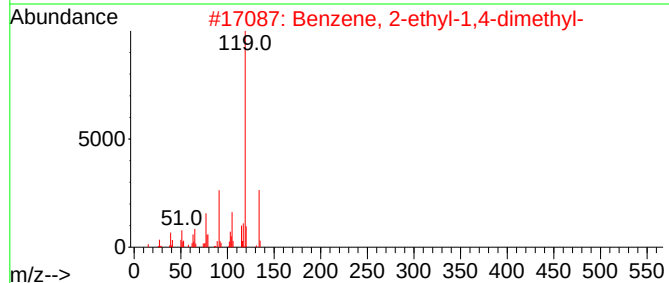
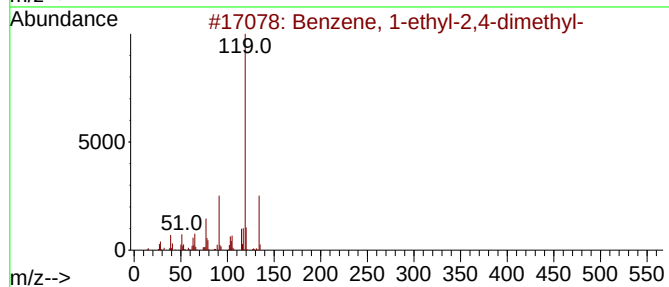
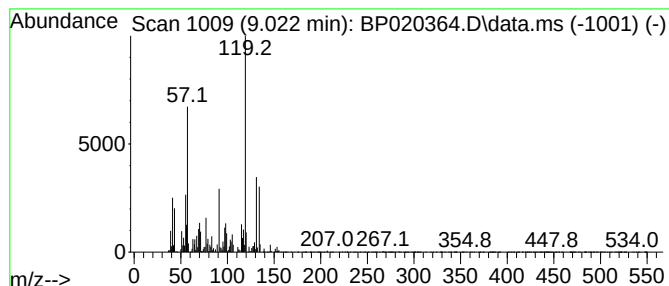
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.022	6.57 ng/ul	214723	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	86
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	86
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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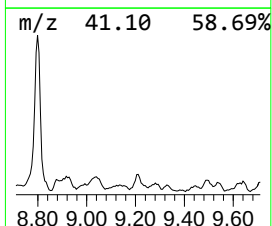
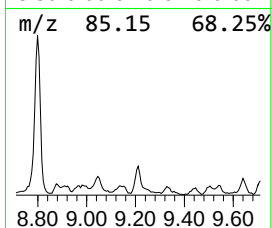
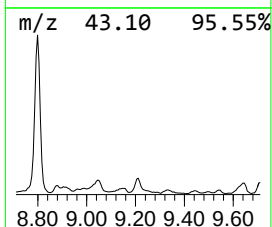
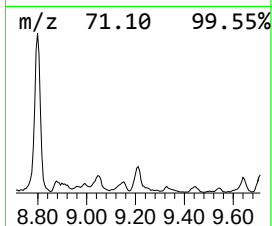
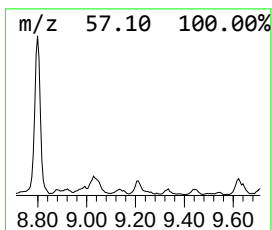
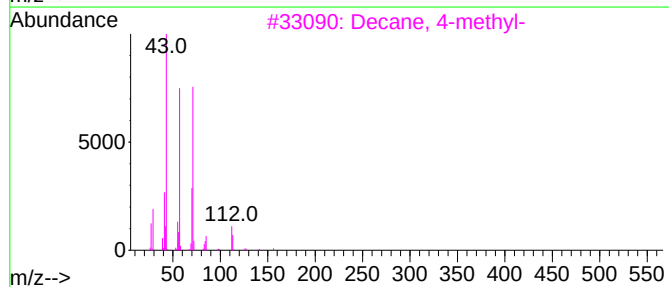
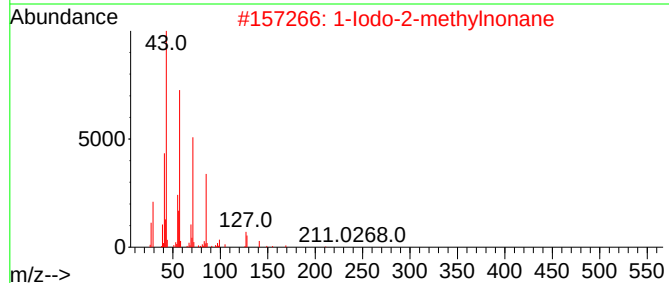
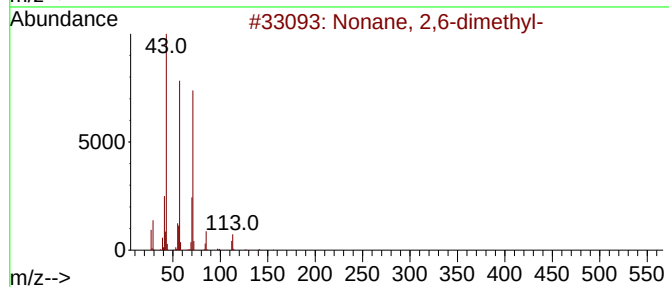
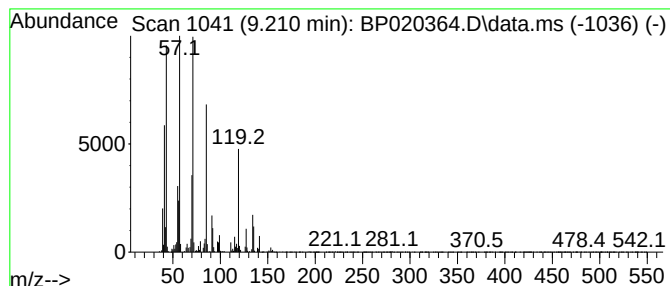
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.210	4.11 ng/ul	134437	Naphthalene-d8	10.387

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	43
2		1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	38
3		Decane, 4-methyl-	156	C11H24	002847-72-5	38
4		Octane, 4-methyl-	128	C9H20	002216-34-4	38
5		Nonane	128	C9H20	000111-84-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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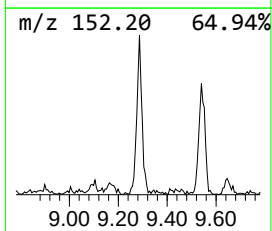
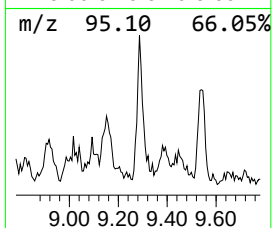
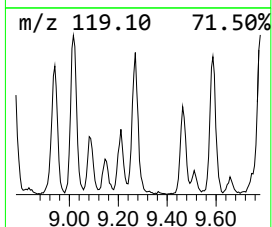
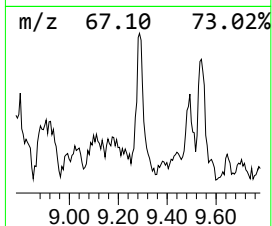
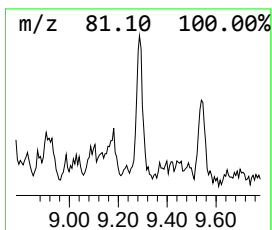
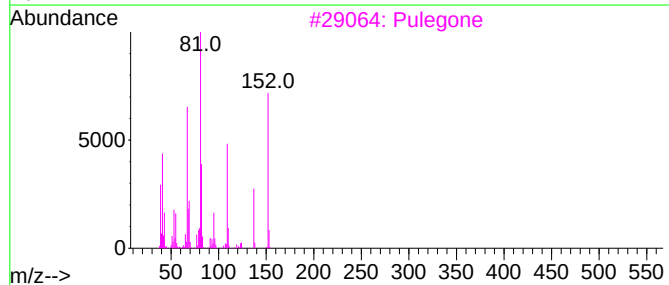
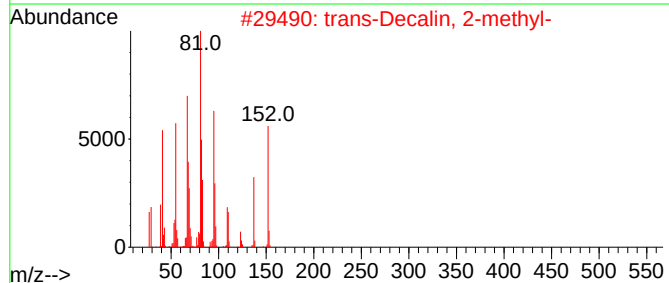
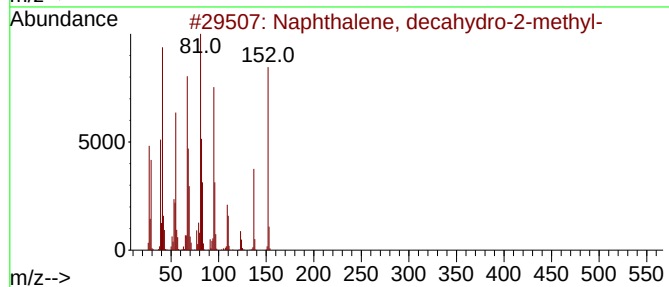
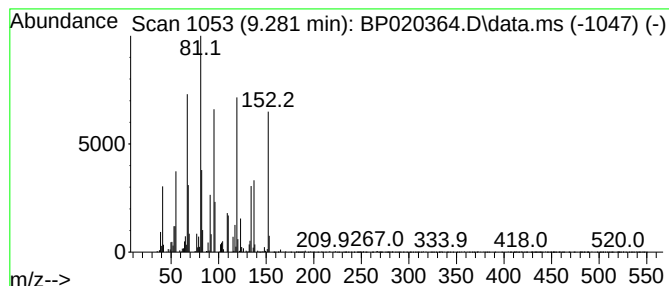
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 Naphthalene, decahydro-2-me... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.281	3.23 ng/ul	105577	Naphthalene-d8	10.387

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	89
2		trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	89
3		Pulegone	152	C10H16O	000089-82-7	60
4		Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	49
5		Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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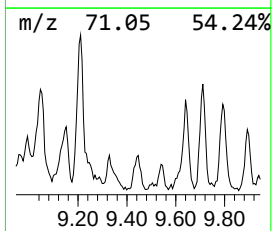
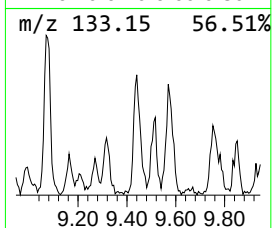
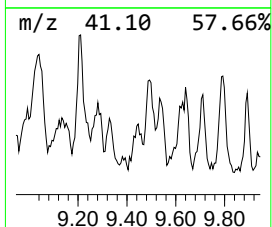
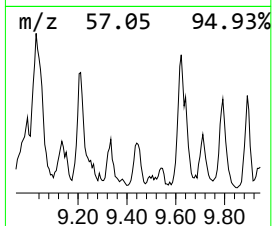
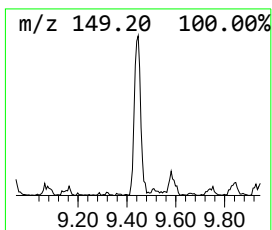
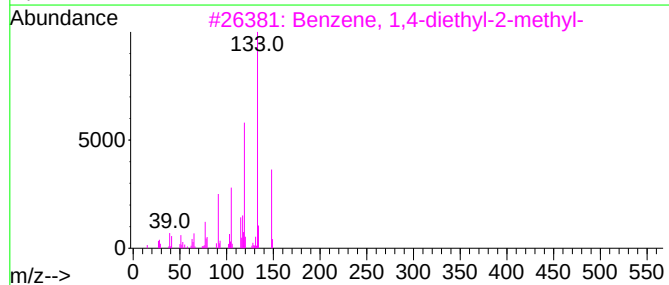
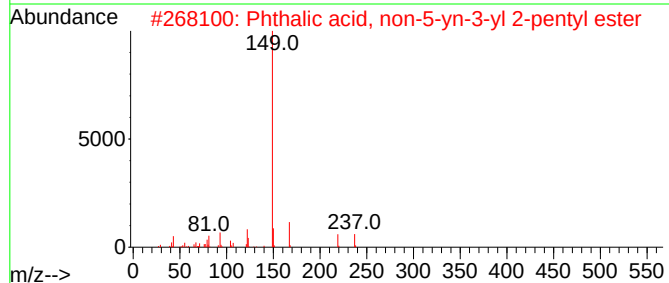
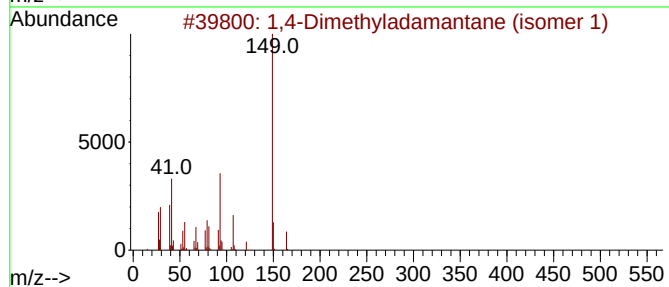
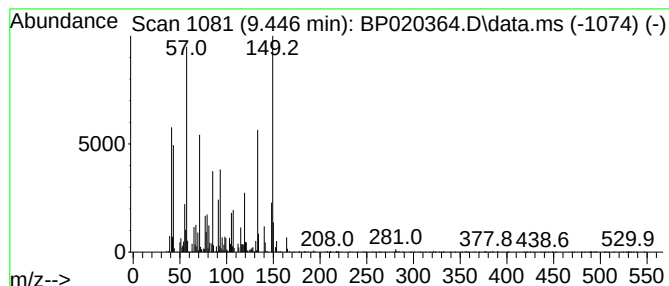
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 1,4-Dimethyladamantane (iso... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.446	3.18 ng/ul	103795	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dimethyladamantane (isomer 1)	164	C12H20	1000460-67-0	50
2			Phthalic acid, non-5-yn-3-yl 2-p...	358	C22H30O4	1000315-74-9	38
3			Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	25
4			m-Tolyl isothiocyanate	149	C8H7NS	000621-30-7	25
5			Benzaldehyde, 3-pentachloropheno...	412	C15H9Cl5O3	329222-74-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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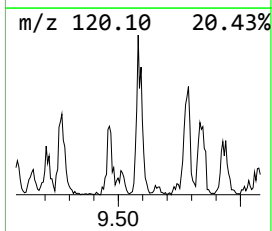
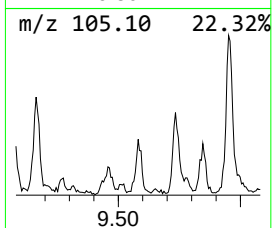
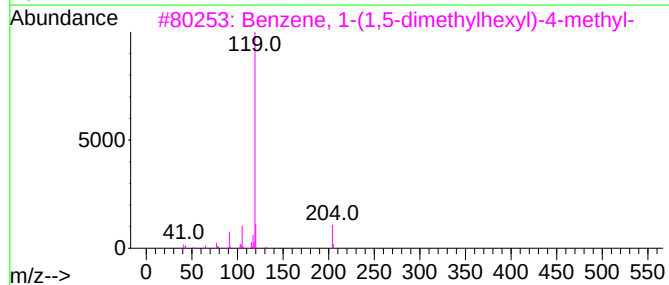
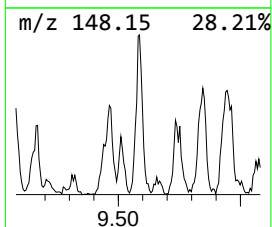
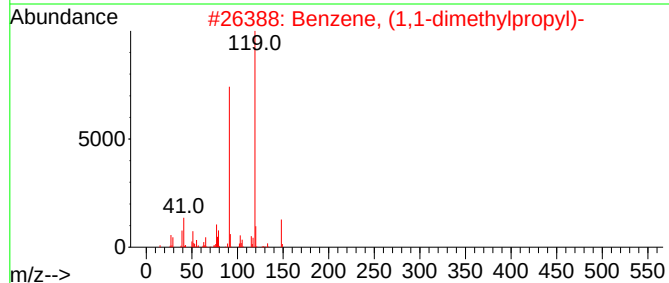
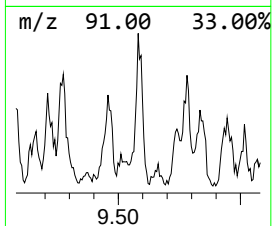
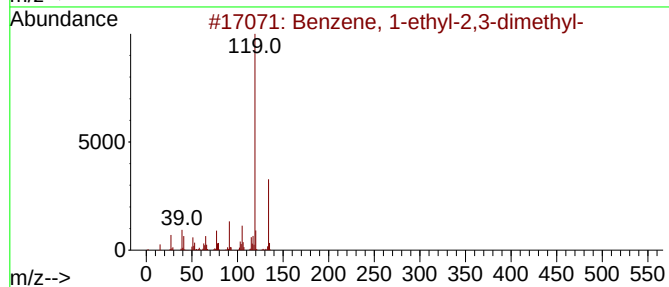
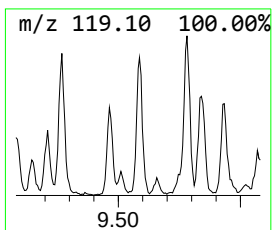
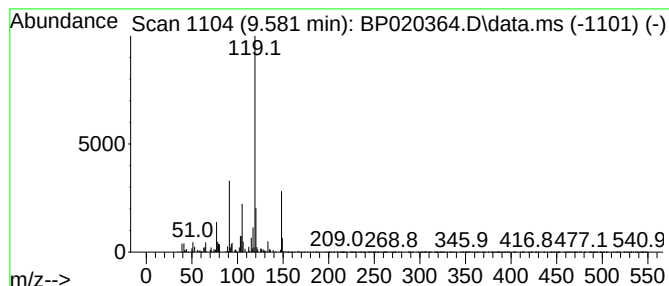
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.581	2.68 ng/ul	87560	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	59
2			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	59
3			Benzene, 1-(1,5-dimethylhexyl)-4...	204	C15H24	001461-02-5	59
4			2,5-Dimethylphenylacetic acid	164	C10H12O2	1000342-65-5	53
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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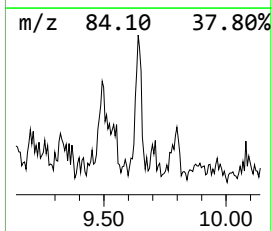
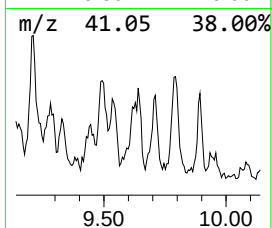
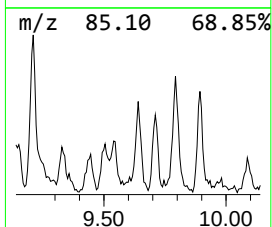
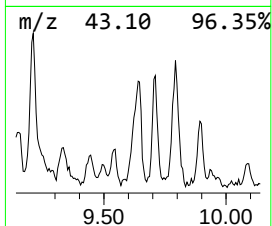
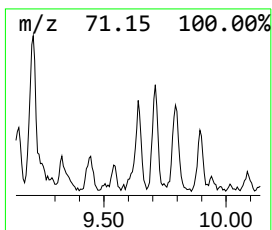
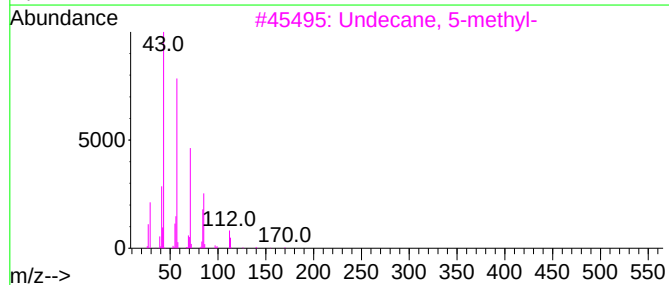
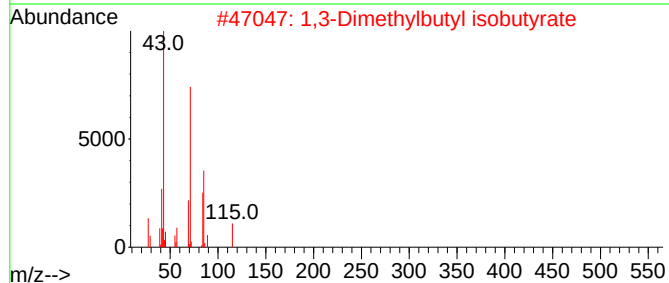
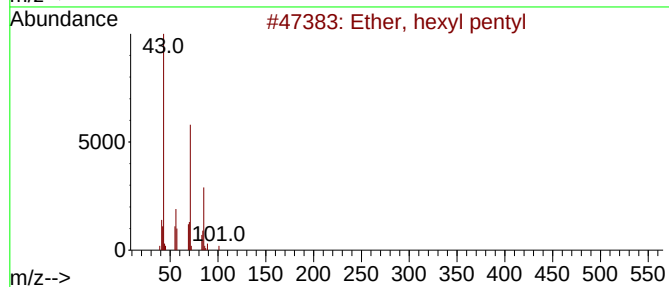
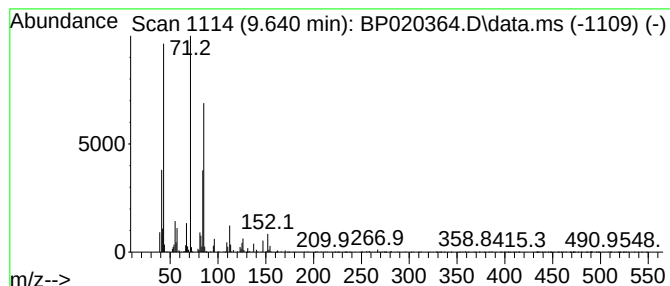
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 Ether, hexyl pentyl Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.640	3.16 ng/ul	103215	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ether, hexyl pentyl	172	C11H24O	032357-83-8	64
2			1,3-Dimethylbutyl isobutyrate	172	C10H20O2	1000429-31-6	50
3			Undecane, 5-methyl-	170	C12H26	001632-70-8	35
4			Butyric acid, 2,2-dimethyl-, vin...	142	C8H14O2	013170-00-8	35
5			Pentane, 3-bromo-	150	C5H11Br	001809-10-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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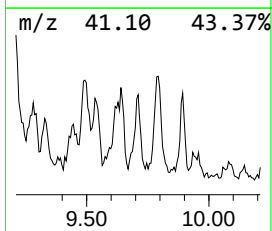
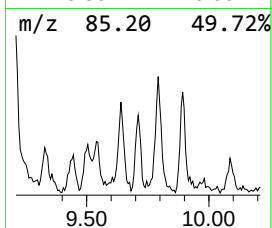
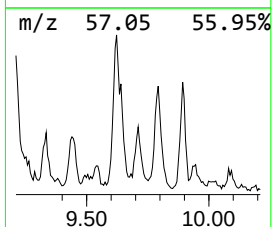
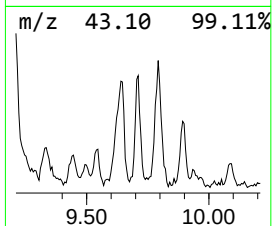
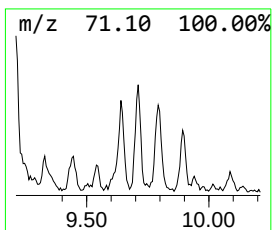
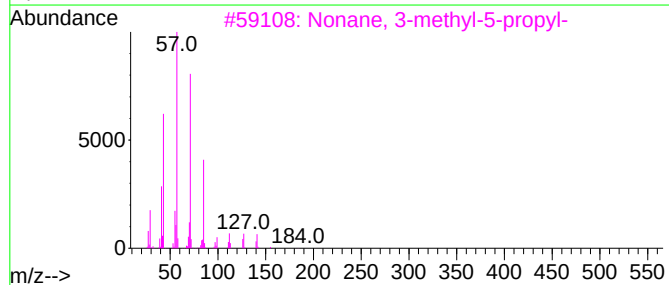
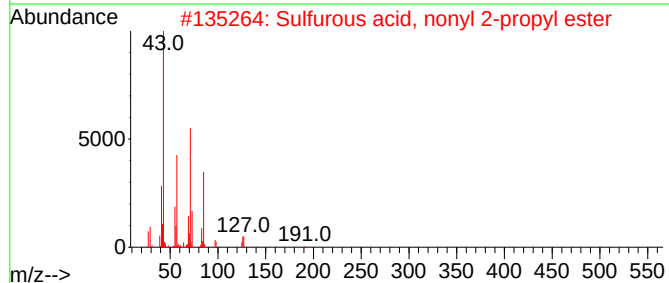
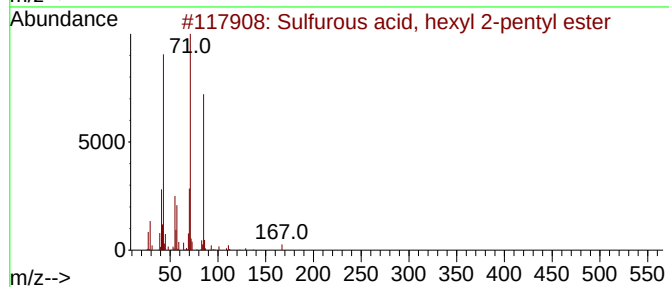
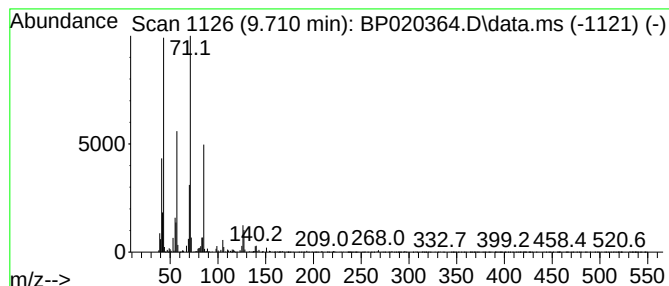
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 Sulfurous acid, hexyl 2-pen... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.710	3.16 ng/ul	103309	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	72
2			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	72
3			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	72
4			Octane, 2,6,6-trimethyl-	156	C11H24	054166-32-4	59
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
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Sample : P2369-13MEDL 5X
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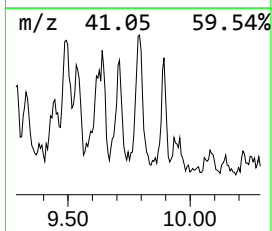
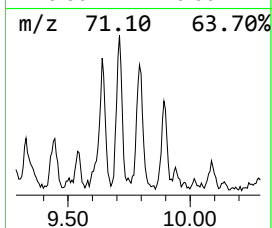
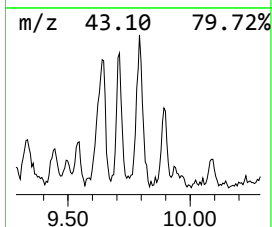
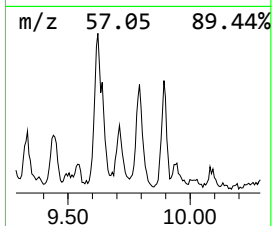
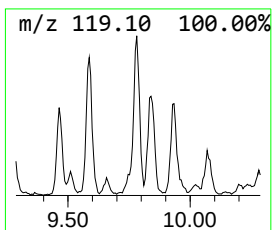
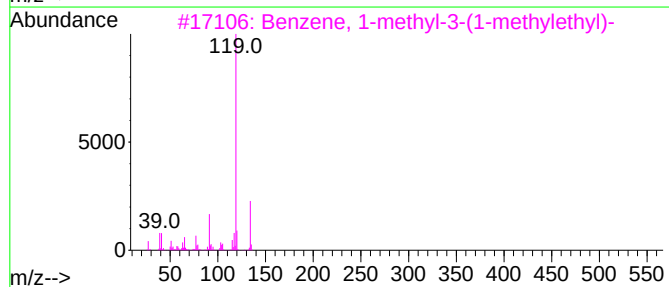
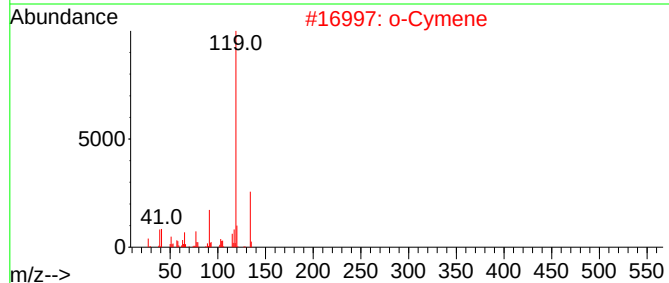
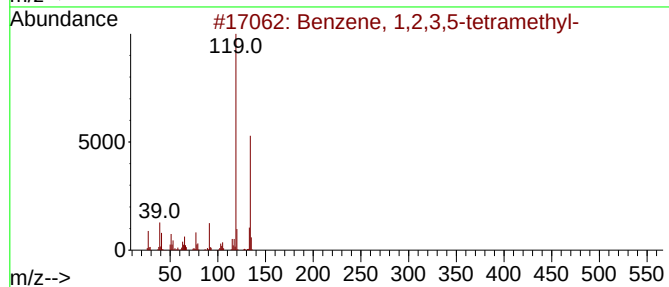
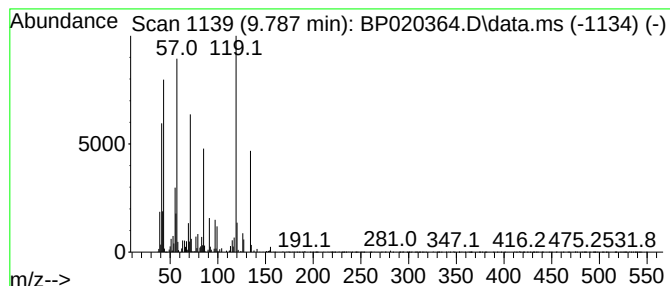
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.787	5.22 ng/ul	170723	Naphthalene-d8	10.387	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	50
2	o-Cymene	134	C10H14	000527-84-4	45
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	45
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	45
5	p-Cymene	134	C10H14	000099-87-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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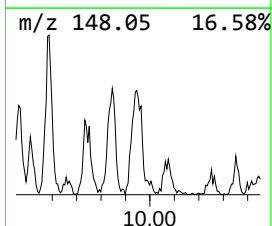
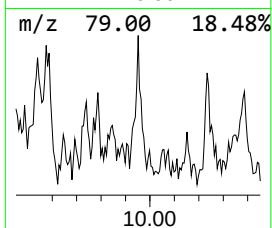
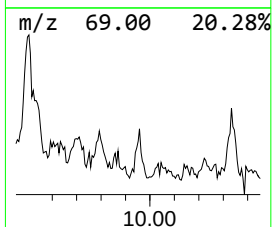
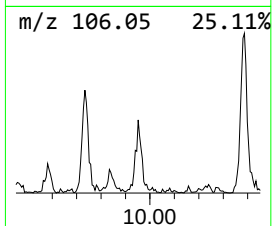
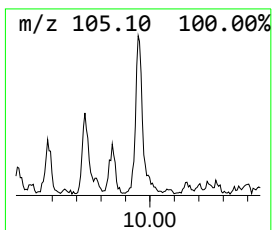
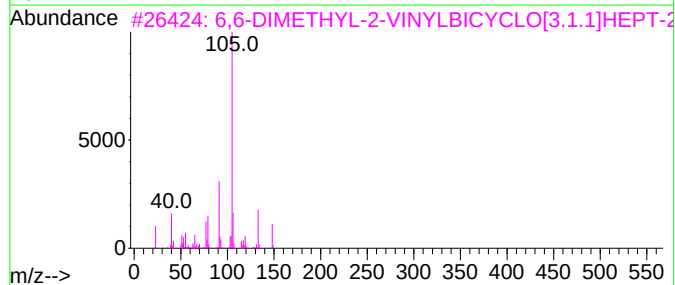
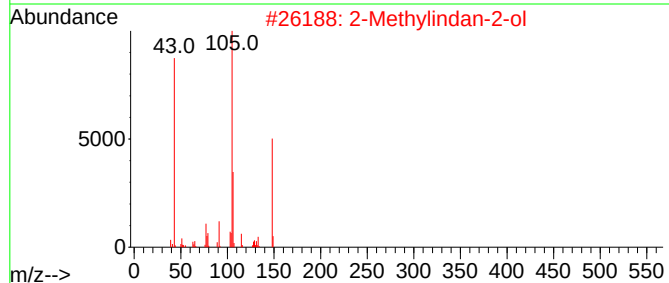
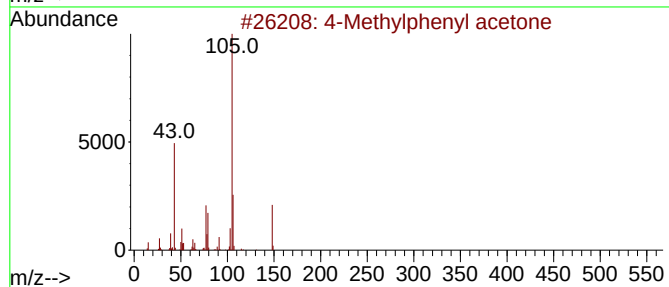
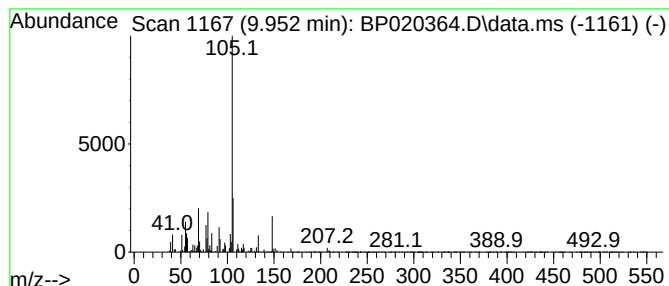
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 4-Methylphenyl acetone Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.952	2.55 ng/ul	83326	Naphthalene-d8	10.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylphenyl acetone	148	C10H12O	002096-86-8	58
2			2-Methylindan-2-ol	148	C10H12O	033223-84-6	58
3			6,6-DIMETHYL-2-VINYLBICYCLO[3.1.1]HEPT-2-ENE	148	C11H16	000473-00-7	52
4			Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	49
5			Benzene, (1,2-dimethylpropyl)-	148	C11H16	004481-30-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

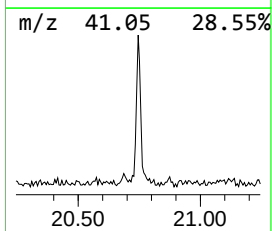
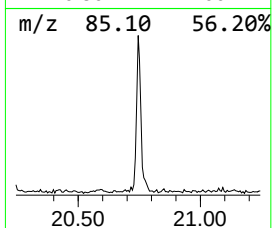
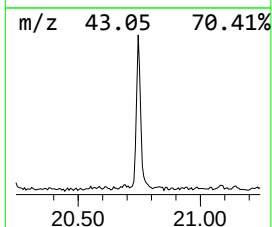
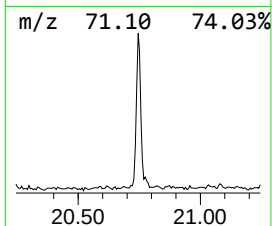
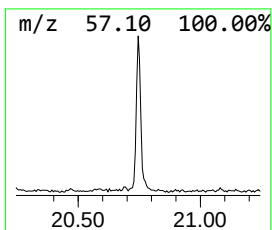
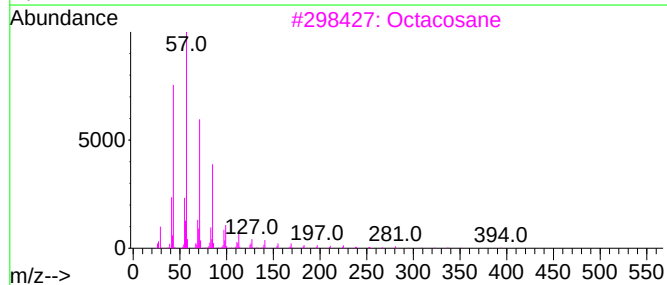
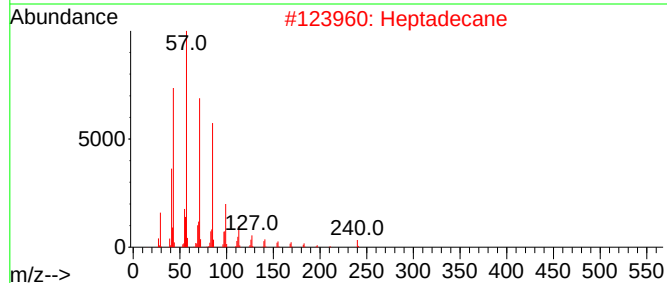
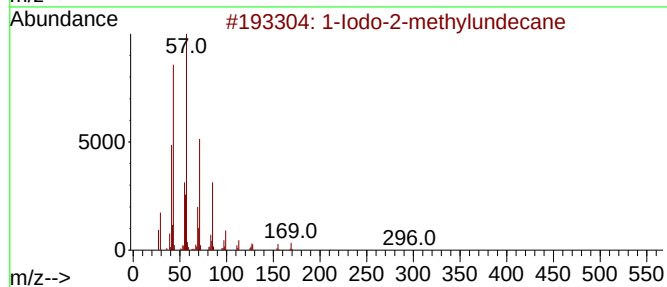
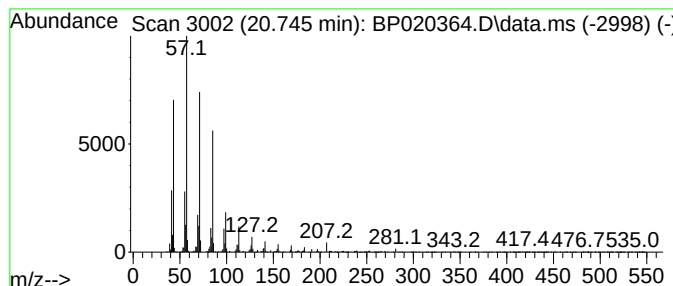
Instrument :
BNA_P
ClientSampleId :
A43N7MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 1-Iodo-2-methylundecane Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.745	2.81 ng/ul	195834	Chrysene-d12	21.610	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	93
2	Heptadecane	240	C17H36	000629-78-7	93
3	Octacosane	394	C28H58	000630-02-4	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43N7MEDL

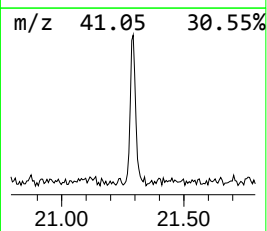
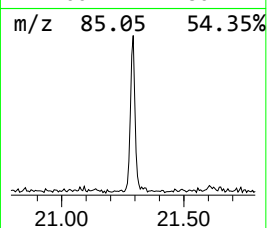
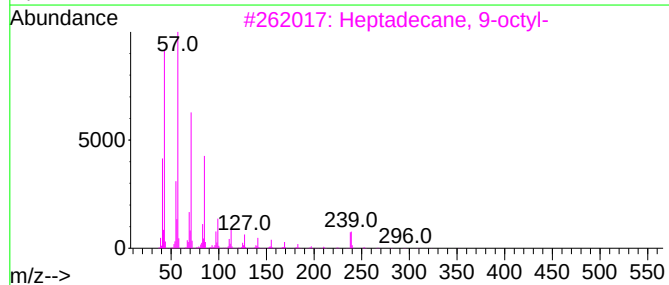
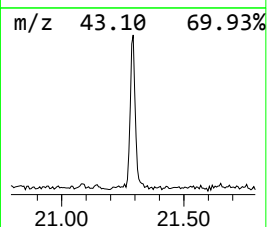
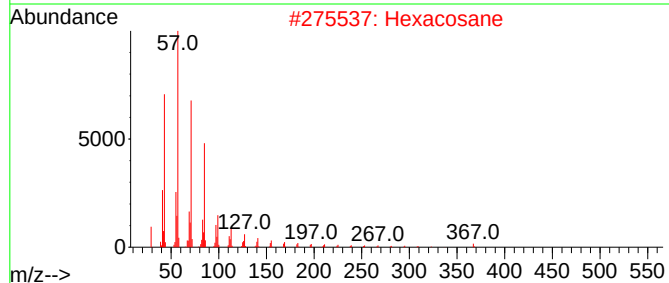
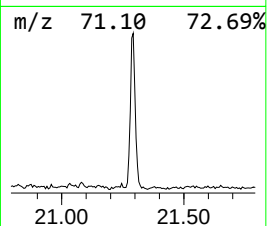
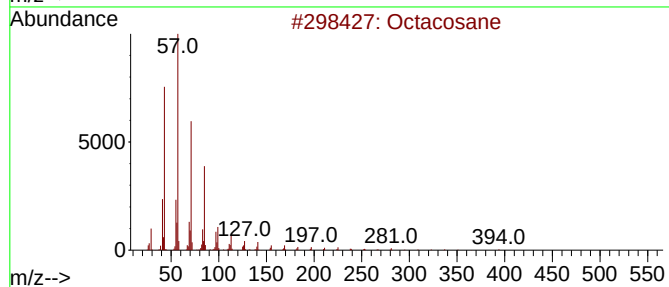
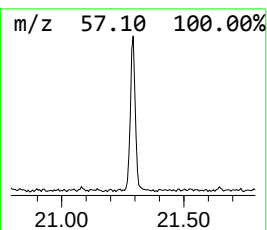
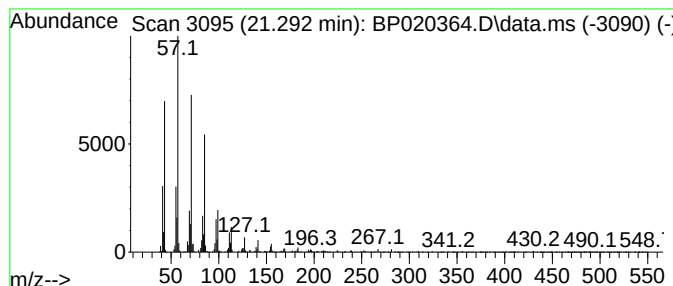
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.292	3.06 ng/ul	212855	Chrysene-d12	21.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020364.D
Acq On : 13 May 2024 22:13
Operator : MA/JU
Sample : P2369-13MEDL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

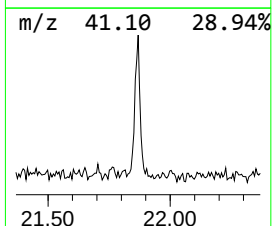
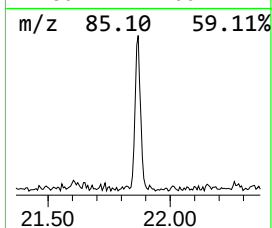
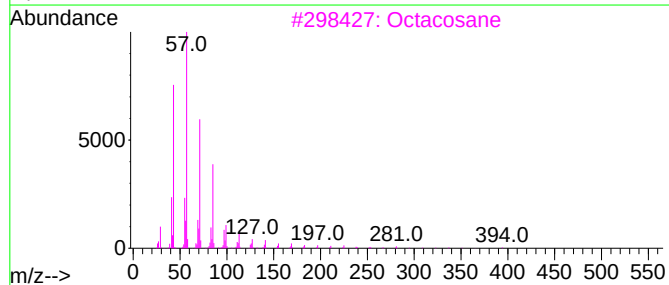
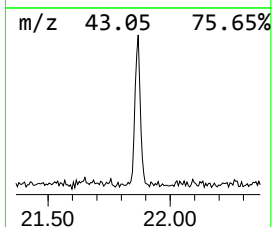
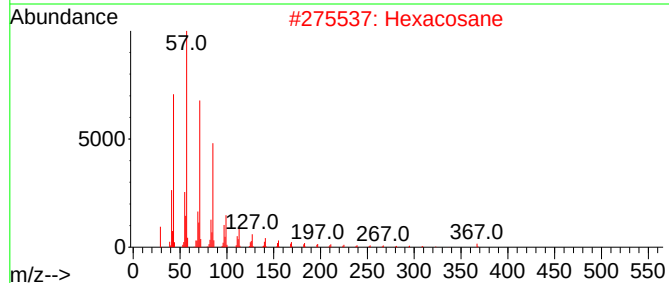
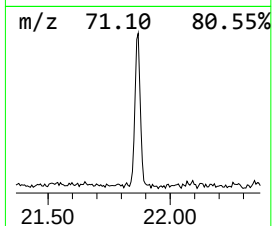
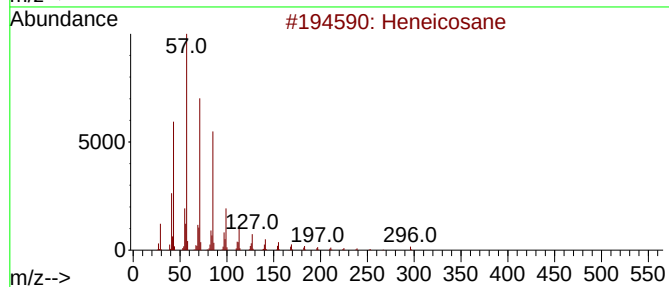
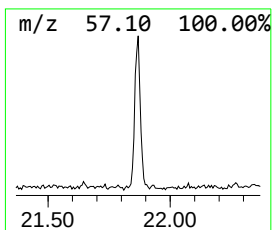
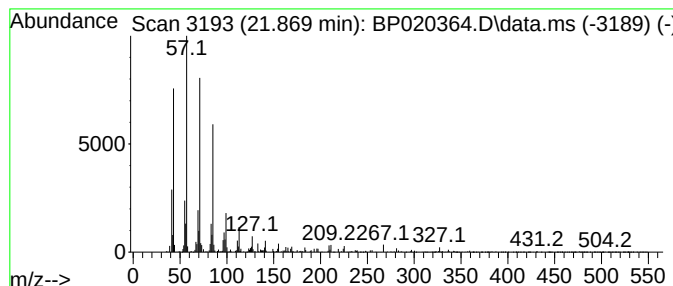
Instrument :
BNA_P
ClientSampleId :
A43N7MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.869	2.02 ng/ul	140468	Chrysene-d12		21.610	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
 Data File : BP020364.D
 Acq On : 13 May 2024 22:13
 Operator : MA/JU
 Sample : P2369-13MEDL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43N7MEDL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	5.887	3.3	ng/ul	98470	1	7.611	603525	20.0
(DEL) Alkane: S...	6.164	3.0	ng/ul	90262	1	7.611	603525	20.0
(DEL) Alkane: S...	6.269	6.9	ng/ul	208131	1	7.611	603525	20.0
(DEL) Alkane: S...	6.593	3.7	ng/ul	110529	1	7.611	603525	20.0
Carbonic acid, ...	6.646	2.8	ng/ul	84896	1	7.611	603525	20.0
(DEL) Alkane: S...	6.752	7.8	ng/ul	234882	1	7.611	603525	20.0
(DEL) Alkane: C...	6.917	2.5	ng/ul	74128	1	7.611	603525	20.0
Cyclohexane, 1-...	7.075	4.1	ng/ul	124411	1	7.611	603525	20.0
Benzene, 1-ethy...	7.258	2.2	ng/ul	67741	1	7.611	603525	20.0
Sulfurous acid,...	7.305	2.5	ng/ul	76190	1	7.611	603525	20.0
1-Decanol, 2-he...	7.411	4.2	ng/ul	125231	1	7.611	603525	20.0
Decan-2-yl isob...	7.499	8.0	ng/ul	242334	1	7.611	603525	20.0
(DEL) Alkane: S...	7.558	13.3	ng/ul	402590	1	7.611	603525	20.0
Benzene, 1,2,4-...	7.711	6.0	ng/ul	182548	1	7.611	603525	20.0
(DEL) Alkane: S...	7.764	5.0	ng/ul	150575	1	7.611	603525	20.0
(DEL) Alkane: S...	7.828	4.1	ng/ul	122568	1	7.611	603525	20.0
(DEL) Alkane: C...	7.864	5.4	ng/ul	161786	1	7.611	603525	20.0
(DEL) Alkane: C...	7.922	3.0	ng/ul	90729	1	7.611	603525	20.0
(DEL) Alkane: S...	8.016	3.4	ng/ul	101234	1	7.611	603525	20.0
(DEL) Alkane: S...	8.099	11.7	ng/ul	352734	1	7.611	603525	20.0
(DEL) Alkane: S...	8.164	6.2	ng/ul	187832	1	7.611	603525	20.0
(DEL) Alkane: S...	8.234	12.7	ng/ul	383096	1	7.611	603525	20.0
Naphthalene, de...	8.405	7.2	ng/ul	216098	1	7.611	603525	20.0
Benzene, 4-ethy...	8.540	4.6	ng/ul	138550	1	7.611	603525	20.0
p-Cymene	8.587	3.2	ng/ul	96706	1	7.611	603525	20.0
(E)-Tetradec-11...	8.646	2.7	ng/ul	80439	1	7.611	603525	20.0
(DEL) Alkane: C...	8.881	3.5	ng/ul	106347	1	7.611	603525	20.0
unknown-01	8.928	8.3	ng/ul	250863	1	7.611	603525	20.0
Benzene, 1-ethy...	9.022	6.6	ng/ul	214723	2	10.387	653753	20.0
(DEL) Alkane: S...	9.210	4.1	ng/ul	134437	2	10.387	653753	20.0
Naphthalene, de...	9.281	3.2	ng/ul	105577	2	10.387	653753	20.0
1,4-Dimethylada...	9.446	3.2	ng/ul	103795	2	10.387	653753	20.0
Benzene, 1-ethy...	9.581	2.7	ng/ul	87560	2	10.387	653753	20.0
Ether, hexyl pe...	9.640	3.2	ng/ul	103215	2	10.387	653753	20.0
Sulfurous acid,...	9.710	3.2	ng/ul	103309	2	10.387	653753	20.0
Benzene, 1,2,3,...	9.787	5.2	ng/ul	170723	2	10.387	653753	20.0
4-Methylphenyl ...	9.952	2.5	ng/ul	83326	2	10.387	653753	20.0
1-Iodo-2-methyl...	20.745	2.8	ng/ul	195834	5	21.610	1392130	20.0
(DEL) Alkane: S...	21.292	3.1	ng/ul	212855	5	21.610	1392130	20.0
(DEL) Alkane: S...	21.869	2.0	ng/ul	140468	5	21.610	1392130	20.0