

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
 Data File : BP015053.D
 Acq On : 10 May 2023 21:05
 Operator : CG/JU
 Sample : 02591-21
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREJ6

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-BP050323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015053.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.452	8	11	24	rVB	28754	62353	2.35%	0.133%
2	4.734	216	229	239	rBV	83633	269201	10.14%	0.574%
3	7.304	655	666	679	rBV2	379061	818562	30.83%	1.745%
4	7.469	688	694	703	rBV	338775	530821	19.99%	1.131%
5	7.675	723	729	738	rVB	408996	675651	25.45%	1.440%
6	8.063	788	795	802	rBV	151944	257723	9.71%	0.549%
7	8.151	802	810	816	rVV	963032	1537822	57.92%	3.278%
8	8.216	816	821	848	rVB	368125	1000067	37.67%	2.132%
9	8.869	925	932	945	rVB	451839	809596	30.49%	1.726%
10	8.987	945	952	960	rBV	674031	1084433	40.85%	2.311%
11	9.122	967	975	989	rVB2	46783	125416	4.72%	0.267%
12	9.328	1004	1010	1017	rBV	405299	686213	25.85%	1.463%
13	9.463	1024	1033	1039	rBV	142876	290290	10.93%	0.619%
14	9.528	1039	1044	1047	rVV2	42316	77149	2.91%	0.164%
15	9.581	1047	1053	1066	rVB	250382	484534	18.25%	1.033%
16	10.057	1128	1134	1145	rBV	327252	555271	20.91%	1.184%
17	10.157	1145	1151	1158	rVV3	46880	86928	3.27%	0.185%
18	10.245	1158	1166	1174	rVV	158706	338810	12.76%	0.722%
19	10.328	1174	1180	1184	rVB6	32847	63687	2.40%	0.136%
20	10.416	1190	1195	1200	rVB	85060	125963	4.74%	0.268%
21	10.492	1200	1208	1213	rBV4	152258	392573	14.79%	0.837%
22	10.604	1221	1227	1233	rVB	482761	793676	29.89%	1.692%
23	10.816	1257	1263	1268	rVV	368516	548078	20.64%	1.168%
24	10.875	1268	1273	1277	rVV6	29989	65289	2.46%	0.139%
25	10.939	1278	1284	1286	rVV3	124169	220058	8.29%	0.469%
26	10.986	1286	1292	1298	rVV	1298291	2231615	84.06%	4.756%
27	11.057	1298	1304	1309	rVV2	256806	568524	21.41%	1.212%
28	11.122	1309	1315	1331	rVV	321391	627150	23.62%	1.337%
29	11.251	1331	1337	1344	rVB8	21060	46193	1.74%	0.098%
30	11.522	1378	1383	1392	rBV2	58840	129871	4.89%	0.277%
31	11.639	1395	1403	1407	rBV4	41794	82711	3.12%	0.176%
32	11.704	1408	1414	1420	rVB	123269	227907	8.58%	0.486%
33	11.775	1420	1426	1432	rBV	167042	300456	11.32%	0.640%
34	11.828	1432	1435	1440	rVB4	47217	73184	2.76%	0.156%
35	11.892	1440	1446	1451	rBV3	94346	159076	5.99%	0.339%
36	11.945	1451	1455	1457	rBV	48427	70731	2.66%	0.151%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
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37	12.016	1463	1467	1474	rVB5	71392	130126	4.90%	0.277%
38	12.086	1474	1479	1486	rVB3	77362	139937	5.27%	0.298%
39	12.169	1489	1493	1499	rBV5	49715	88580	3.34%	0.189%
40	12.228	1499	1503	1508	rVB3	32833	50803	1.91%	0.108%

41	12.286	1508	1513	1515	rBV4	32768	55647	2.10%	0.119%
42	12.316	1515	1518	1523	rVV3	37586	72369	2.73%	0.154%
43	12.375	1525	1528	1534	rVV	51176	94708	3.57%	0.202%
44	12.445	1534	1540	1548	rVV	287600	449084	16.92%	0.957%
45	12.516	1549	1552	1558	rVV6	31889	59622	2.25%	0.127%

46	12.604	1559	1567	1580	rVB	839961	1635235	61.59%	3.485%
47	12.739	1586	1590	1600	rVB6	41402	104149	3.92%	0.222%
48	12.845	1601	1608	1612	rBV6	66270	144793	5.45%	0.309%
49	12.998	1629	1634	1640	rVB8	37553	76878	2.90%	0.164%
50	13.063	1642	1645	1648	rVV4	28355	40881	1.54%	0.087%

51	13.110	1648	1653	1661	rVB4	120262	237177	8.93%	0.506%
52	13.222	1667	1672	1674	rVV2	66024	106881	4.03%	0.228%
53	13.280	1678	1682	1689	rVV6	100802	245104	9.23%	0.522%
54	13.339	1690	1692	1697	rVB4	53032	63055	2.38%	0.134%
55	13.398	1697	1702	1709	rVB7	38127	90783	3.42%	0.193%

56	13.516	1717	1722	1725	rBV3	61294	91620	3.45%	0.195%
57	13.598	1730	1736	1738	rVV3	97622	207176	7.80%	0.442%
58	13.633	1738	1742	1748	rVV2	295952	438177	16.50%	0.934%
59	13.686	1748	1751	1754	rVV3	35519	56756	2.14%	0.121%
60	13.739	1756	1760	1769	rVB2	74275	133062	5.01%	0.284%

61	14.075	1811	1817	1820	rBV7	32970	76624	2.89%	0.163%
62	14.133	1821	1827	1831	rBV	970571	1373370	51.73%	2.927%
63	14.192	1833	1837	1842	rVB	759725	999836	37.66%	2.131%
64	14.310	1852	1857	1862	rBV6	90863	176780	6.66%	0.377%
65	14.439	1874	1879	1883	rBV	511736	663220	24.98%	1.414%

66	14.486	1883	1887	1896	rVB2	1015475	1553979	58.53%	3.312%
67	14.569	1897	1901	1904	rBV3	59509	74671	2.81%	0.159%
68	14.680	1914	1920	1927	rBV3	162213	331411	12.48%	0.706%
69	14.792	1934	1939	1945	rVB	1621036	2427760	91.44%	5.175%
70	14.857	1945	1950	1958	rBV	121728	199063	7.50%	0.424%

71	14.986	1968	1972	1977	rBV	283757	459580	17.31%	0.980%
72	15.186	2004	2006	2010	rVB	103430	131195	4.94%	0.280%
73	15.227	2011	2013	2017	rVB3	41342	39322	1.48%	0.084%
74	15.292	2020	2024	2029	rVB3	74819	111058	4.18%	0.237%
75	15.357	2032	2035	2041	rVB2	80554	116644	4.39%	0.249%

76	15.422	2042	2046	2056	rVB9	81389	204136	7.69%	0.435%
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77	15.557	2066	2069	2076	rVB5	70882	115226	4.34%	0.246%
78	15.622	2076	2080	2085	rBV3	135455	205359	7.74%	0.438%
79	15.786	2102	2108	2113	rBV	1148568	1695140	63.85%	3.613%
80	15.898	2124	2127	2135	rVB3	129686	212997	8.02%	0.454%
81	16.010	2141	2146	2152	rBV4	117482	287568	10.83%	0.613%
82	16.145	2167	2169	2172	rVB	119711	113600	4.28%	0.242%
83	16.245	2183	2186	2190	rVB2	83625	102891	3.88%	0.219%
84	16.486	2223	2227	2228	rBV	150742	178024	6.71%	0.379%
85	16.745	2268	2271	2277	rVB4	85581	147561	5.56%	0.315%
86	16.827	2281	2285	2289	rBV2	130565	225732	8.50%	0.481%
87	16.874	2289	2293	2296	rBV	254955	396443	14.93%	0.845%
88	17.016	2314	2317	2322	rVB	280784	352546	13.28%	0.751%
89	17.198	2342	2348	2358	rBV	1214984	1922096	72.40%	4.097%
90	17.280	2359	2362	2363	rBV	138219	133169	5.02%	0.284%
91	17.551	2402	2408	2412	rBV	1877770	2654909	100.00%	5.659%
92	17.592	2412	2415	2420	rVV	418123	674576	25.41%	1.438%
93	17.651	2420	2425	2432	rVB2	1065272	1734571	65.33%	3.697%
94	18.033	2485	2490	2493	rBV4	249176	428283	16.13%	0.913%
95	18.292	2530	2534	2538	rBV2	292168	533974	20.11%	1.138%
96	18.404	2549	2553	2558	rBV	887861	1246263	46.94%	2.656%
97	18.692	2598	2602	2605	rBV	401429	675534	25.44%	1.440%
98	18.927	2639	2642	2646	rBV3	407722	605077	22.79%	1.290%
99	19.245	2692	2696	2700	rBV2	933899	1284982	48.40%	2.739%
100	19.292	2701	2704	2708	rBV	542083	822274	30.97%	1.753%

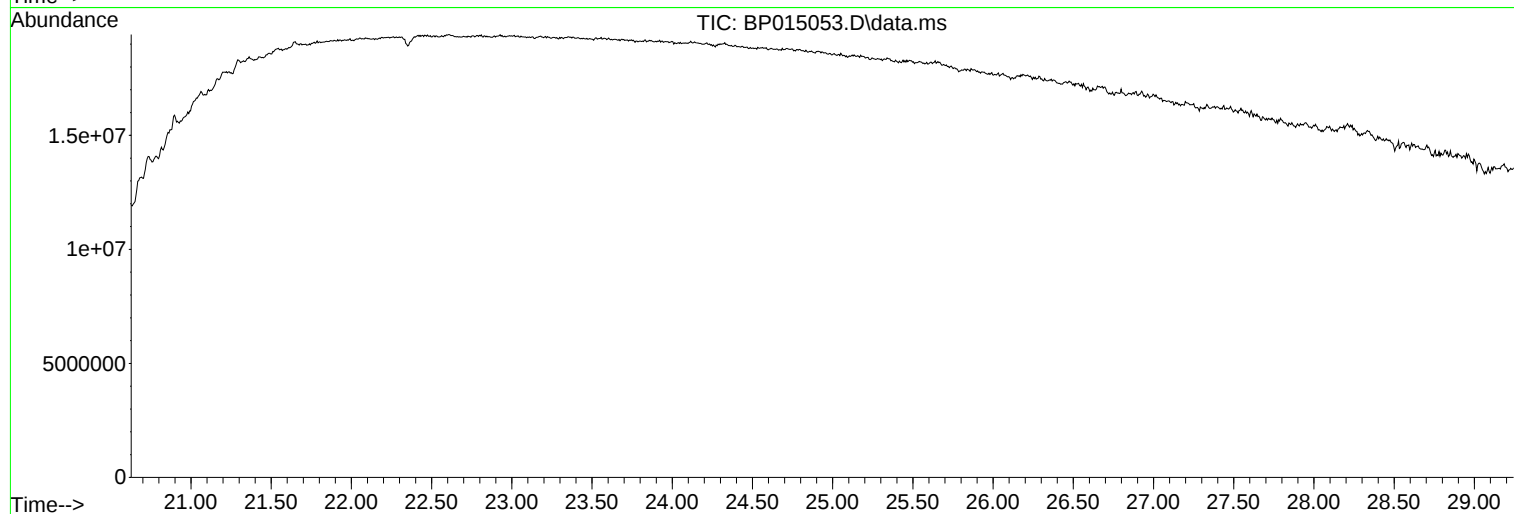
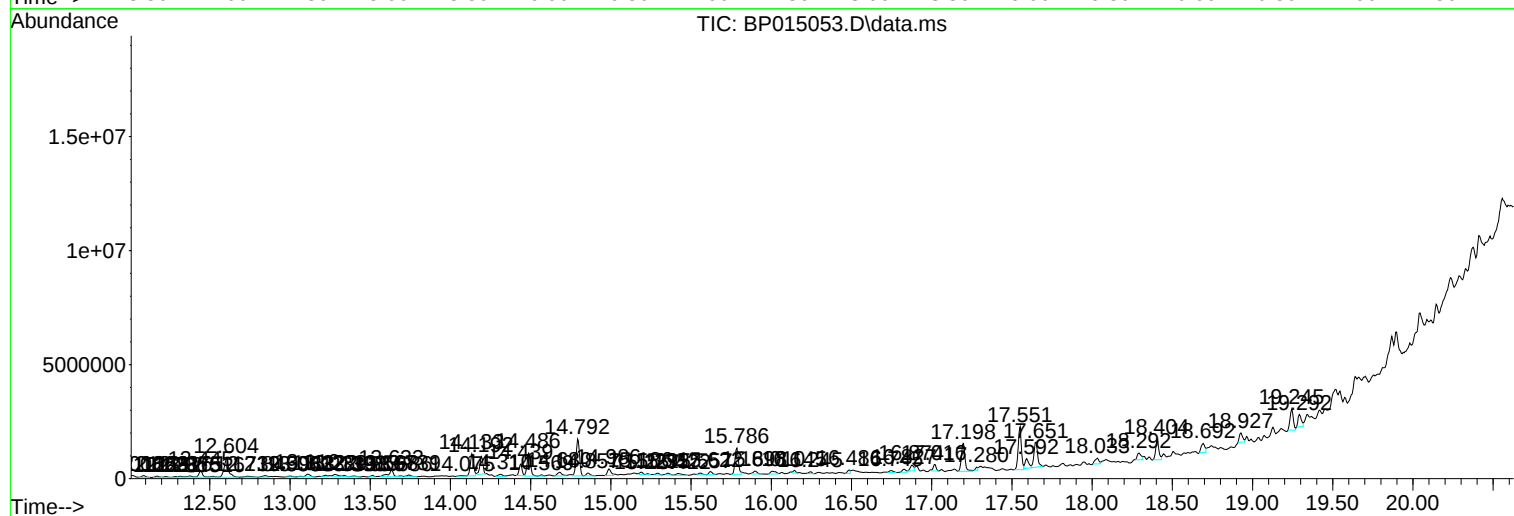
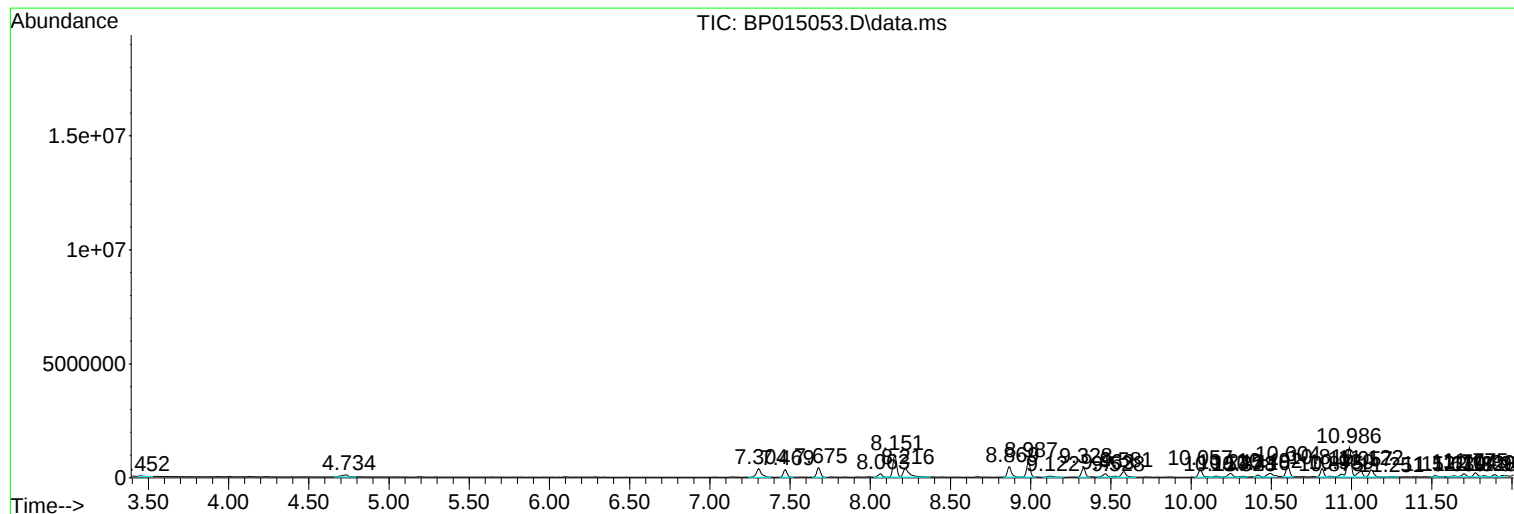
Sum of corrected areas: 46917629

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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



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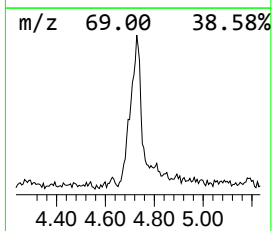
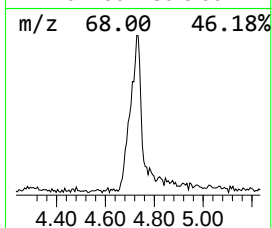
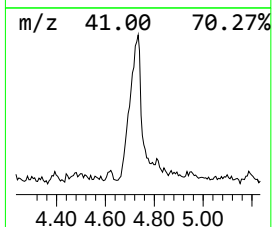
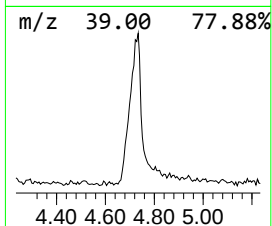
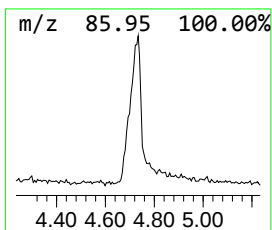
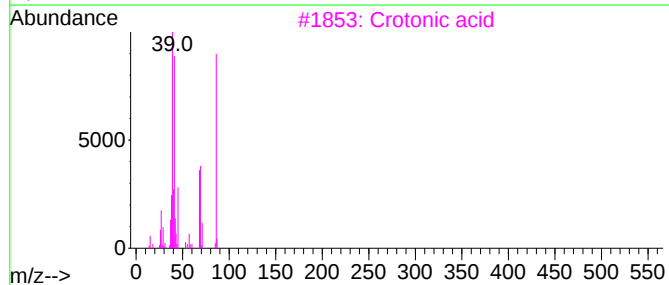
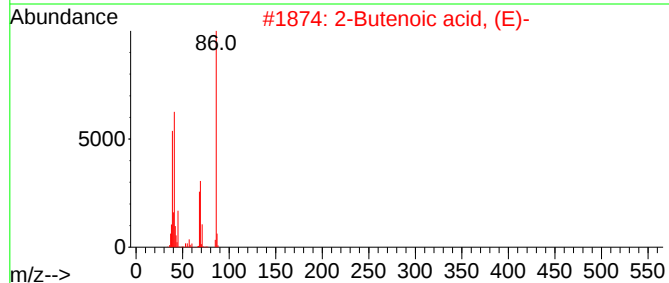
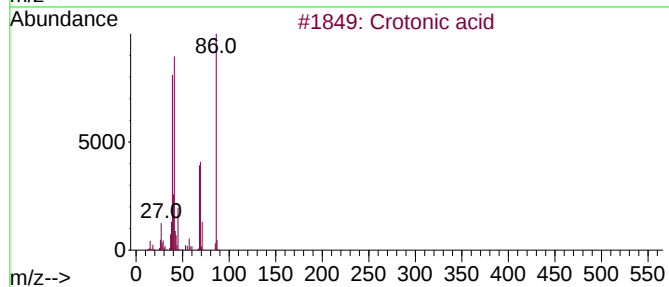
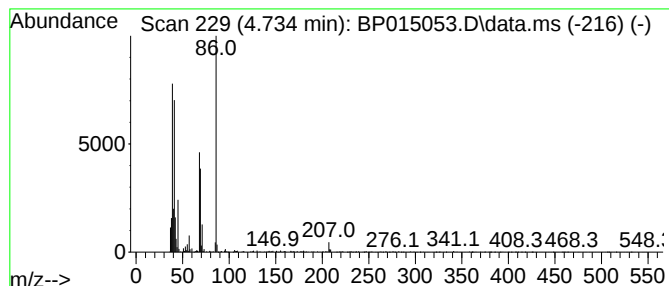
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Crotonic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.734	3.50 ng/ul	269201	1,4-Dichlorobenzene-d4	8.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Crotonic acid	86	C4H6O2	003724-65-0	94
2			2-Butenoic acid, (E)-	86	C4H6O2	000107-93-7	94
3			Crotonic acid	86	C4H6O2	003724-65-0	94
4			Isocrotonic acid	86	C4H6O2	000503-64-0	93
5			2-Butenoic acid, (E)-	86	C4H6O2	000107-93-7	93



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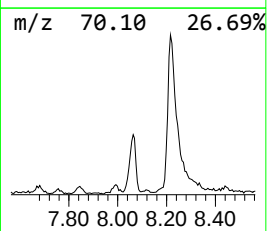
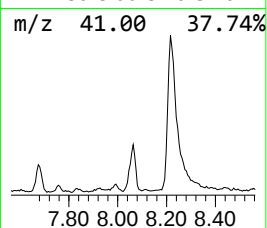
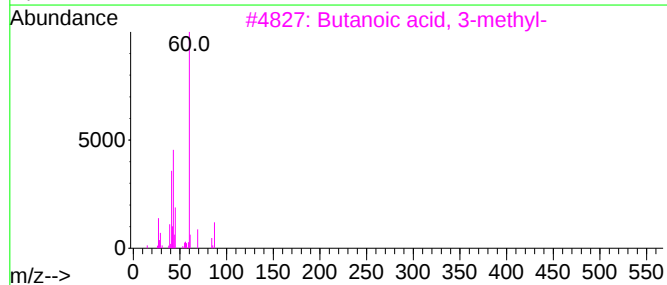
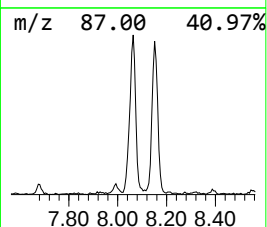
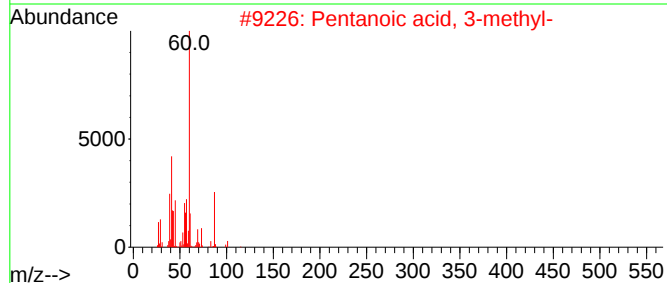
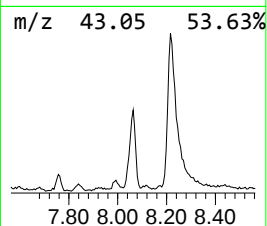
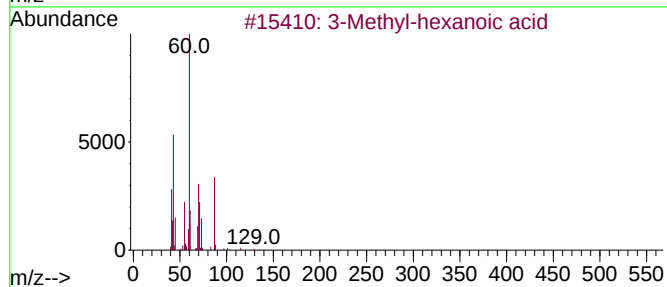
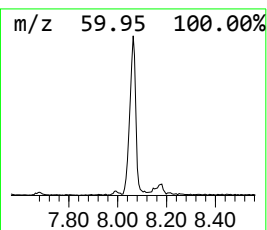
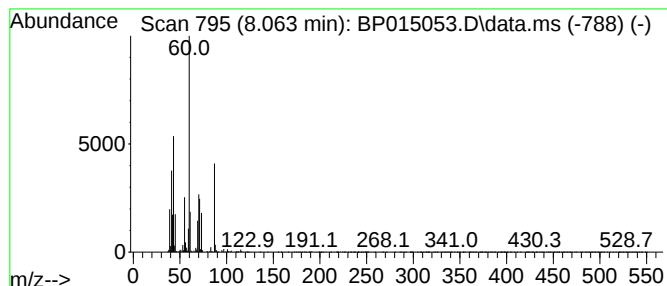
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 3-Methyl-hexanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.063	3.35 ng/ul	257723	1,4-Dichlorobenzene-d4	8.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methyl-hexanoic acid	130	C7H14O2	1000365-65-4	91
2			Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	40
3			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	38
4			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	38
5			Heptanoic acid	130	C7H14O2	000111-14-8	37



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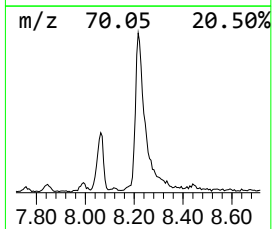
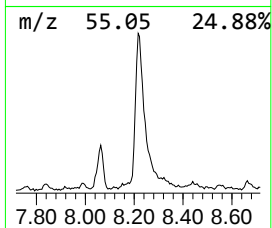
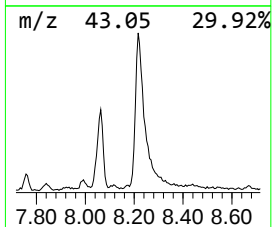
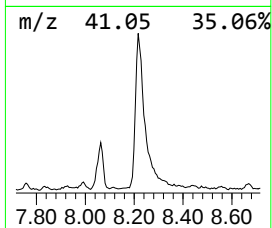
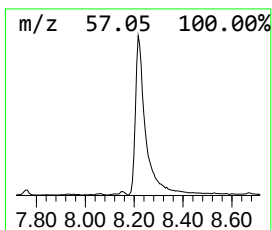
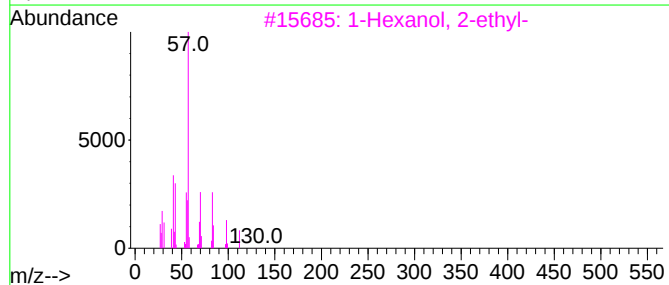
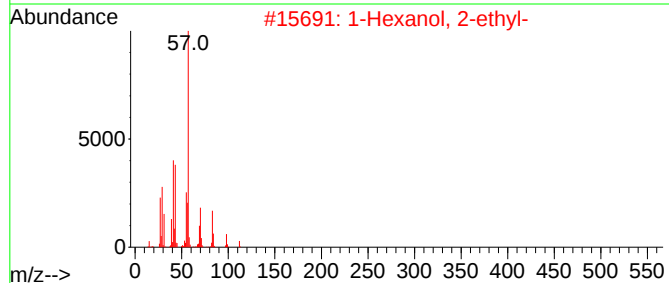
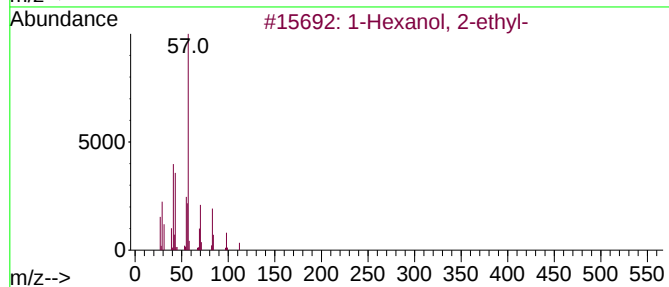
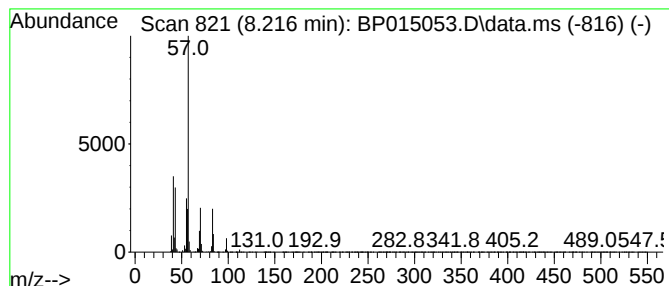
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.216	13.01 ng/ul	1000070	1,4-Dichlorobenzene-d4	8.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	74
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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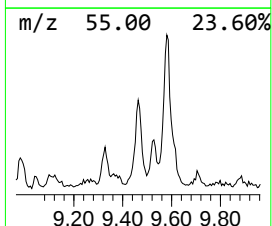
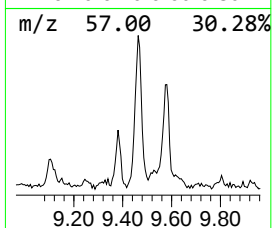
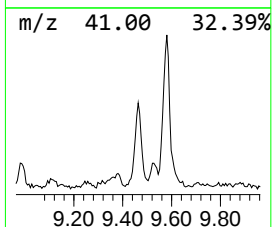
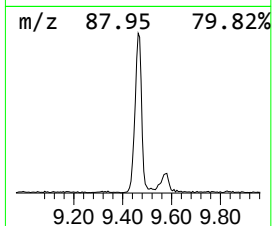
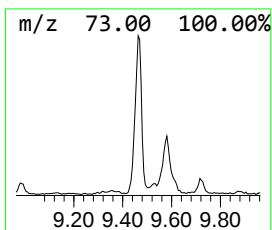
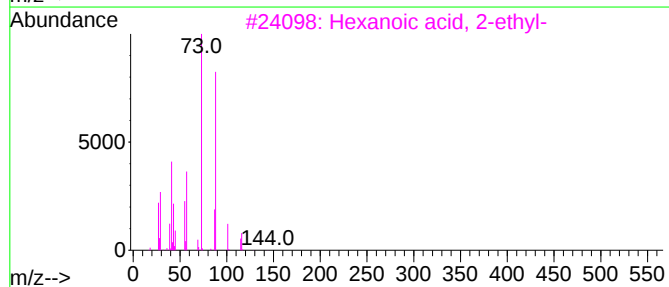
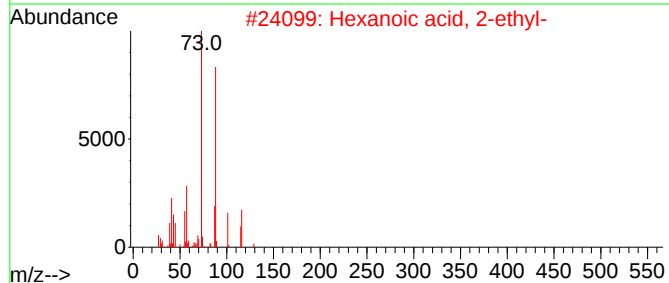
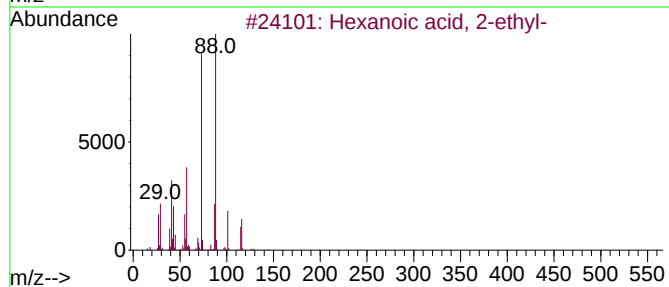
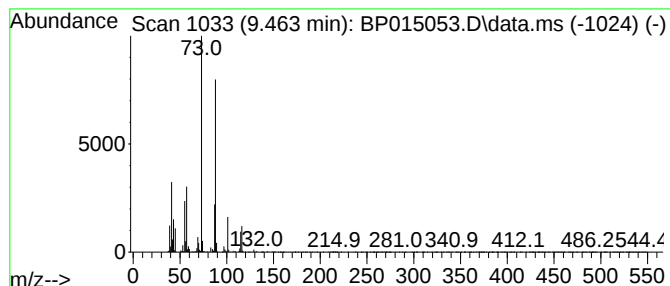
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.463	3.78 ng/ul	290290	1,4-Dichlorobenzene-d4	8.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	95
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72
5			Acetic acid, diethyl-	116	C6H12O2	000088-09-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
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Sample : 02591-21
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ClientSampleId :
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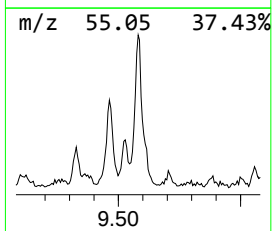
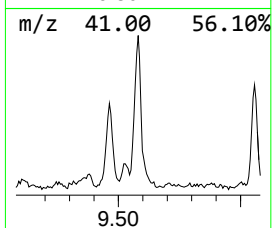
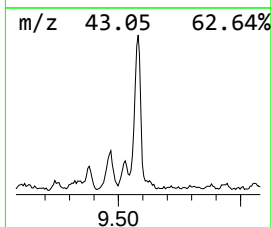
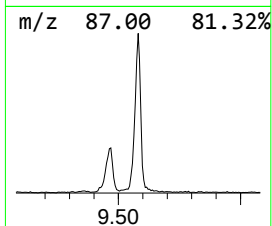
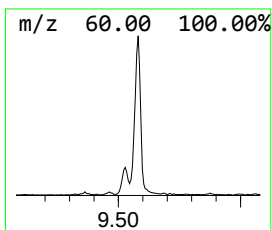
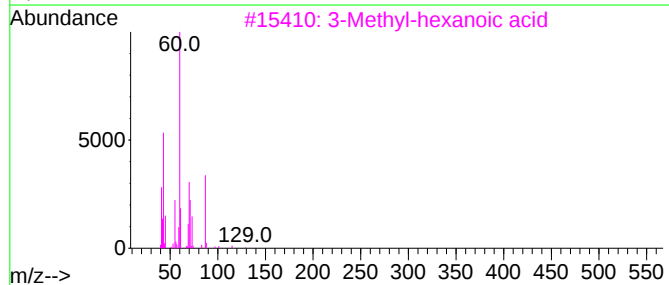
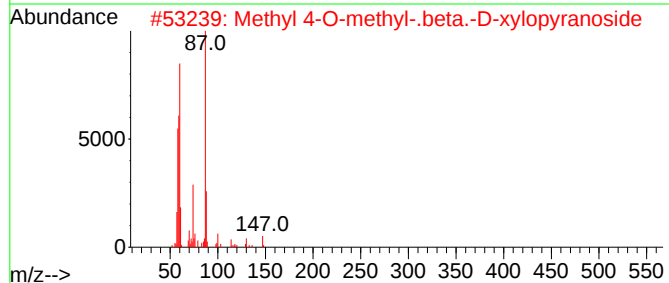
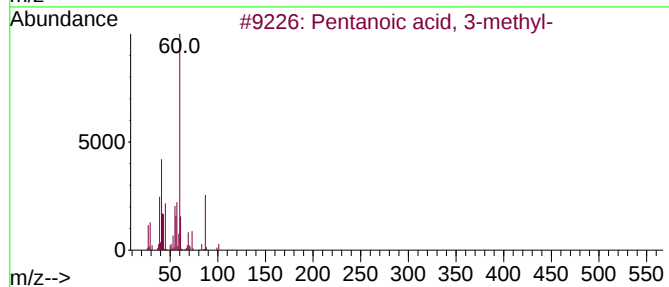
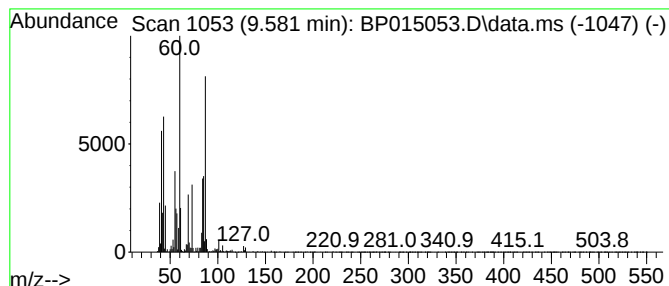
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.581	4.34 ng/ul	484534	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	47
2			Methyl 4-O-methyl-.beta.-D-xylop...	178	C7H14O5	058918-59-5	40
3			3-Methyl-hexanoic acid	130	C7H14O2	1000365-65-4	37
4			3-Methyloctanoic acid	158	C9H18O2	006061-10-5	32
5			Butanoic acid	88	C4H8O2	000107-92-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
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Misc :
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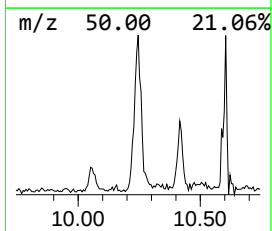
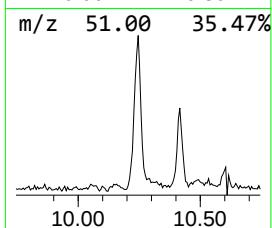
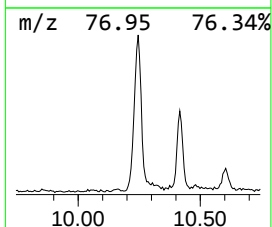
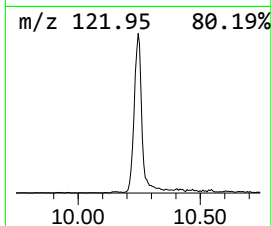
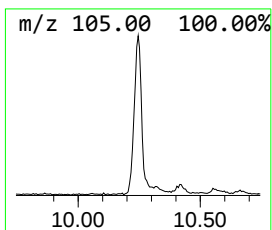
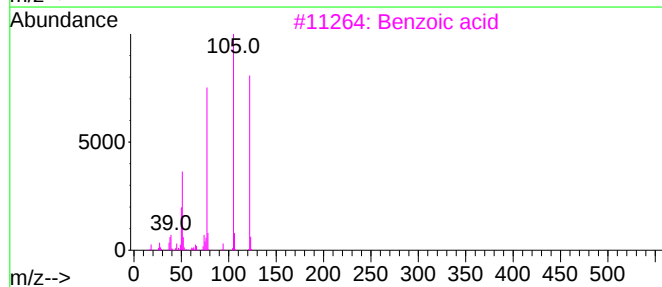
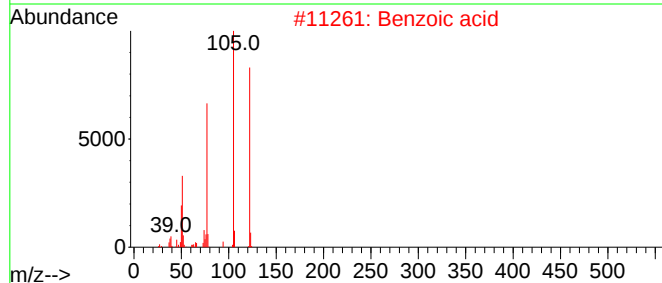
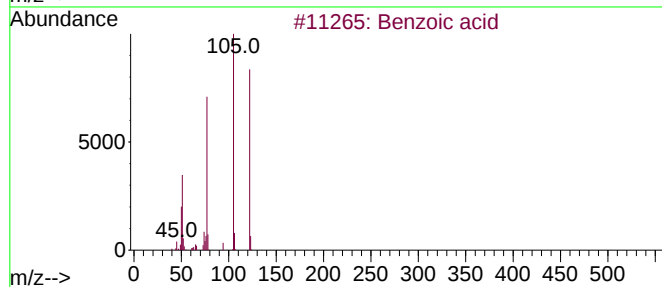
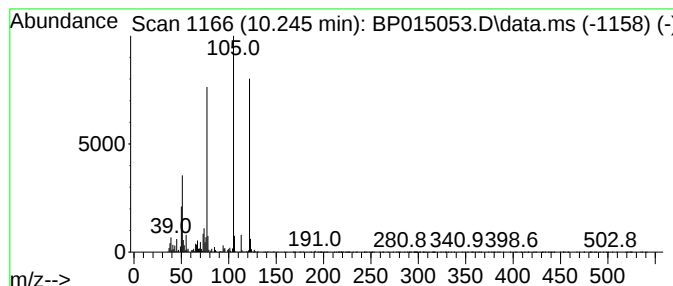
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.245	3.04 ng/ul	338810	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	97
2			Benzoic acid	122	C7H6O2	000065-85-0	96
3			Benzoic acid	122	C7H6O2	000065-85-0	94
4			Benzoic acid	122	C7H6O2	000065-85-0	93
5			Benzoic acid, silver(1+) salt	228	C7H5AgO2	000532-31-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
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ALS Vial : 14 Sample Multiplier: 1

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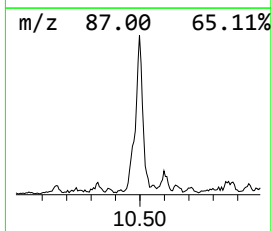
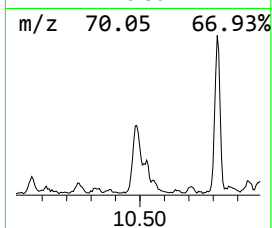
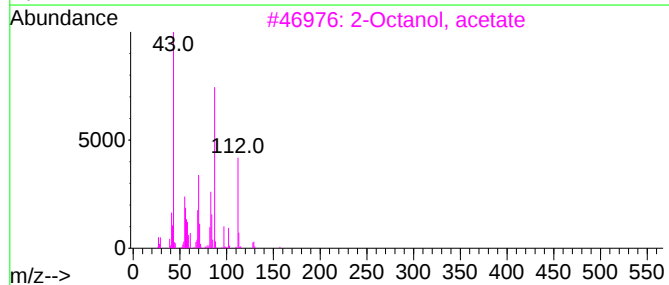
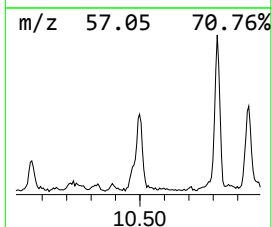
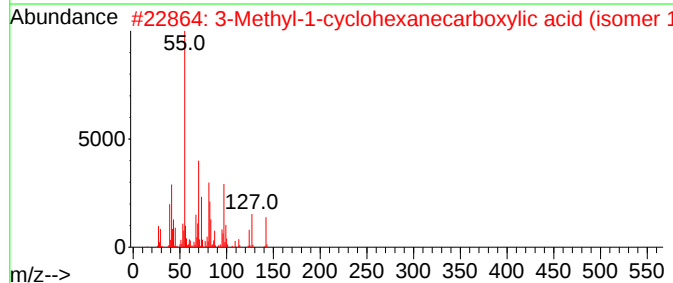
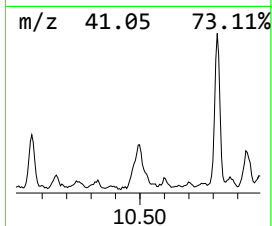
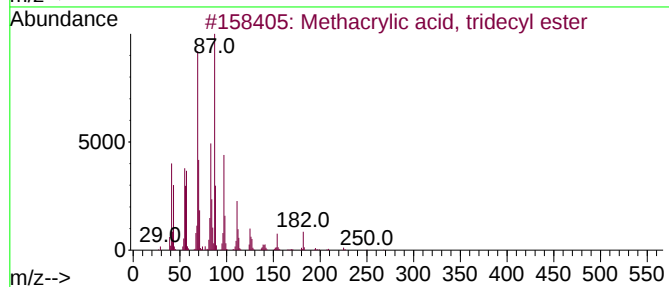
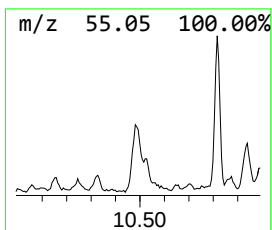
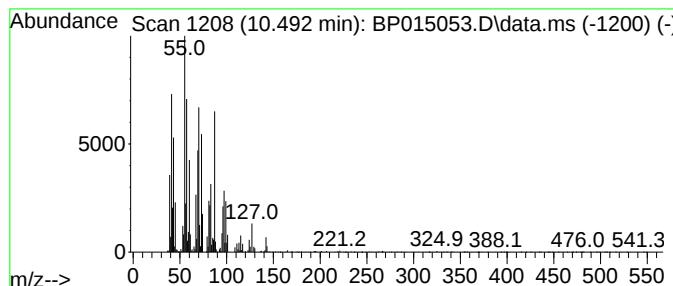
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.492	3.52 ng/ul	392573	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylic acid, tridecyl ester	268	C17H32O2	1000340-28-9	38
2			3-Methyl-1-cyclohexanecarboxylic...	142	C8H14O2	1000453-57-0	38
3			2-Octanol, acetate	172	C10H20O2	002051-50-5	27
4			2-Octenoic acid, (E)-	142	C8H14O2	001871-67-6	27
5			Hex-5-enoic acid, ethyl ester	142	C8H14O2	1000302-89-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
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ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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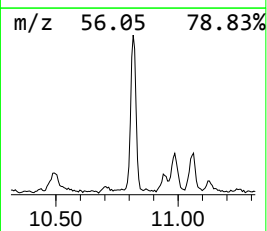
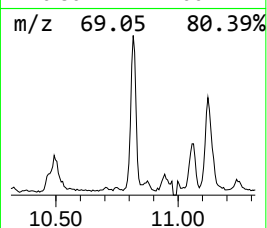
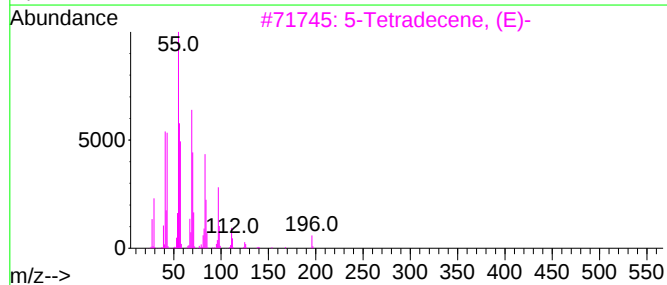
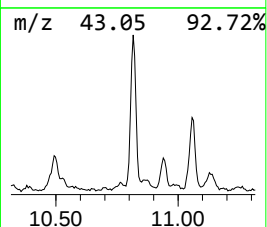
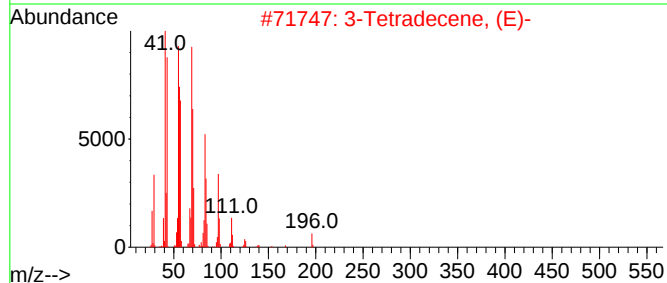
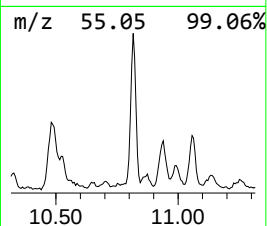
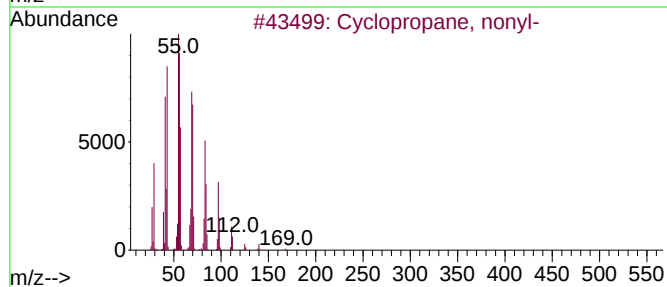
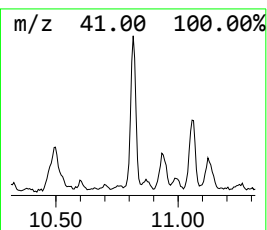
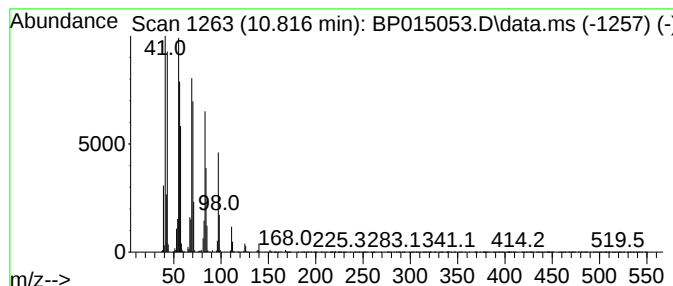
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic10.816 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.816	4.91 ng/ul	548078	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, nonyl-	168	C12H24	074663-85-7	93
2			3-Tetradecene, (E)-	196	C14H28	041446-68-8	91
3			5-Tetradecene, (E)-	196	C14H28	041446-66-6	91
4			1-Decene	140	C10H20	000872-05-9	90
5			3-Tetradecene, (Z)-	196	C14H28	041446-67-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
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Acq On : 10 May 2023 21:05
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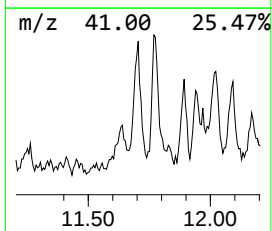
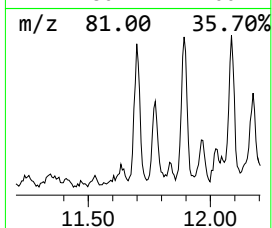
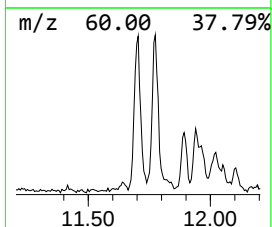
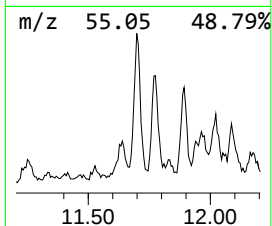
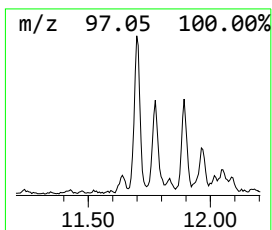
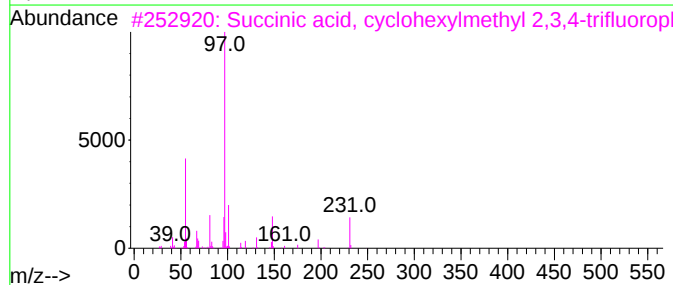
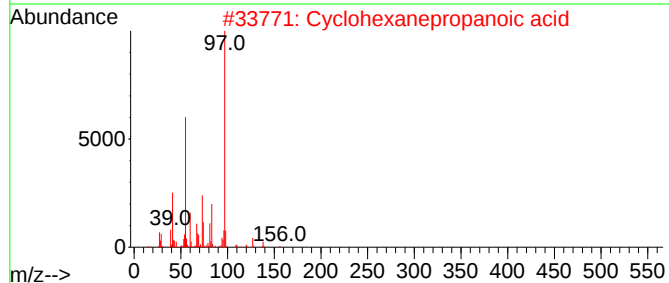
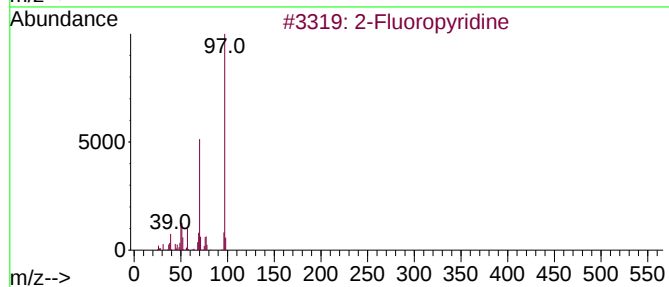
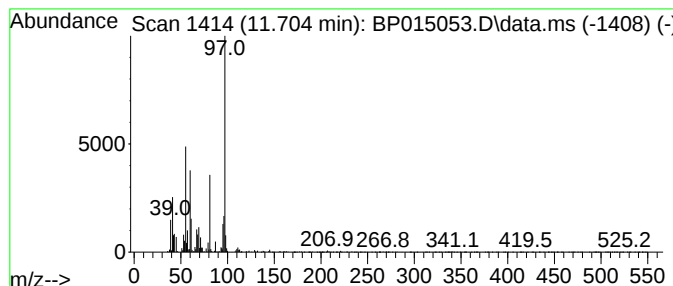
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-Fluoropyridine Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.704	2.04 ng/ul	227907	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Fluoropyridine	97	C5H4FN	000372-48-5	50
2			Cyclohexanepropanoic acid	156	C9H16O2	000701-97-3	50
3			Succinic acid, cyclohexylmethyl ...	344	C17H19F3O4	1000390-76-4	50
4			Cyclohexane, 1-bromo-2-methyl-	176	C7H13Br	006294-39-9	50
5			4-Methylcyclohexaneacetic acid (...)	156	C9H16O2	1000453-59-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
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Acq On : 10 May 2023 21:05
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Sample : 02591-21
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ALS Vial : 14 Sample Multiplier: 1

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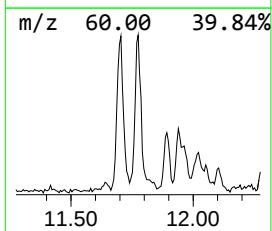
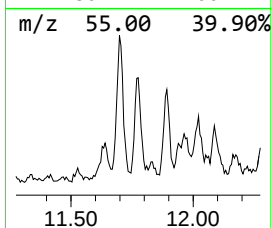
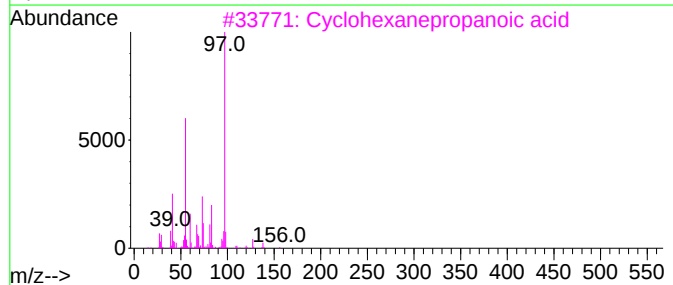
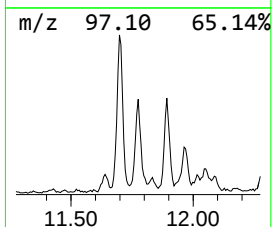
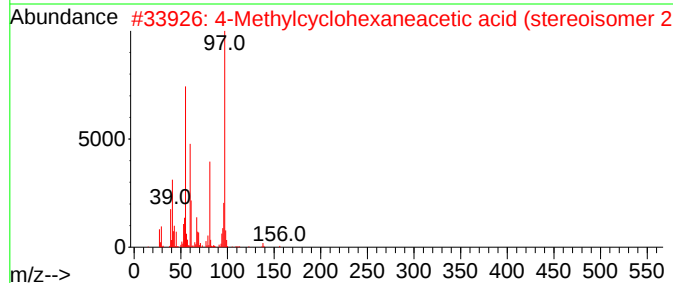
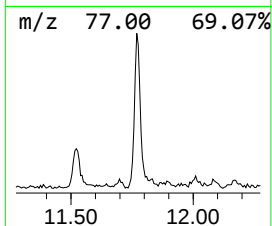
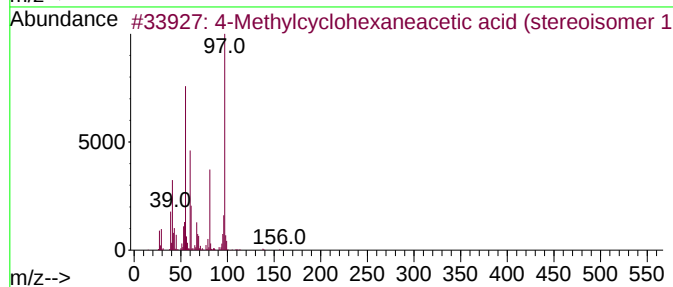
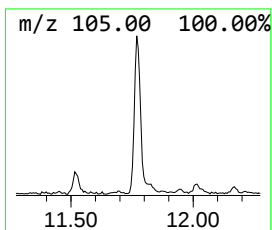
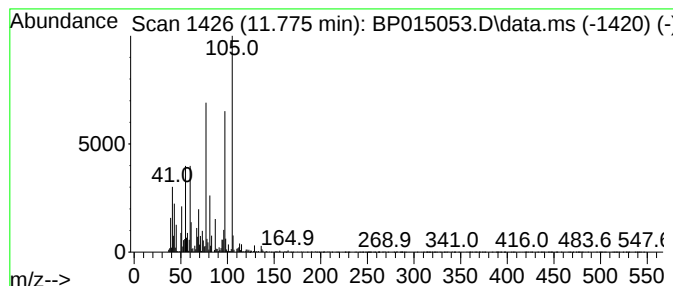
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-03 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.775	2.69 ng/ul	300456	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylcyclohexaneacetic acid (...)	156	C9H16O2	1000453-59-3	27
2			4-Methylcyclohexaneacetic acid (...)	156	C9H16O2	1000453-59-2	22
3			Cyclohexanepropanoic acid	156	C9H16O2	000701-97-3	14
4			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	10
5			4H-1,2,4-Triazole, 4-ethyl-	97	C4H7N3	043183-55-7	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

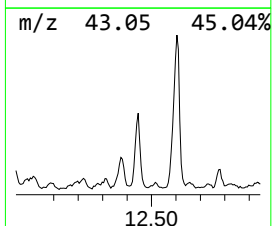
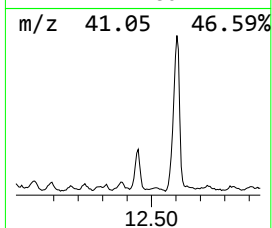
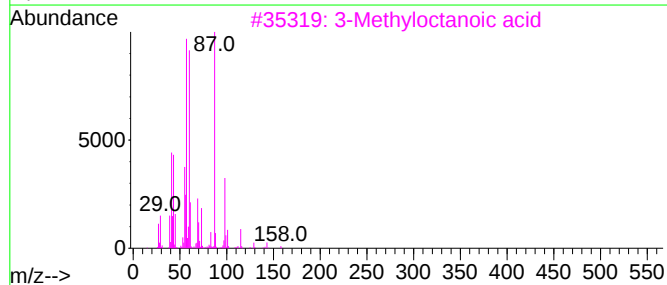
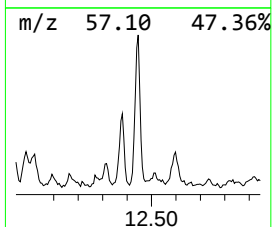
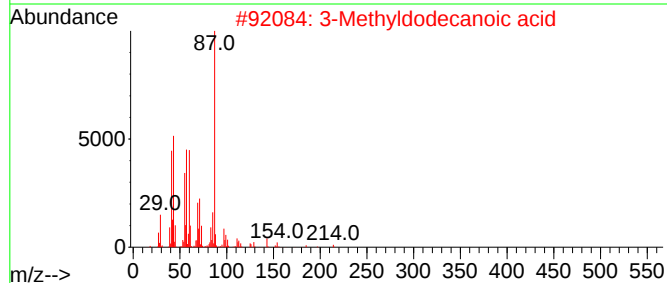
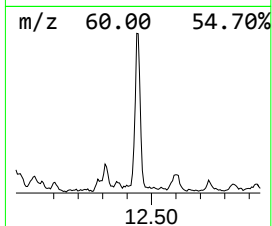
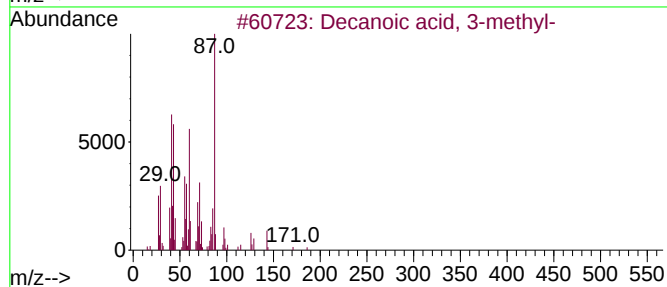
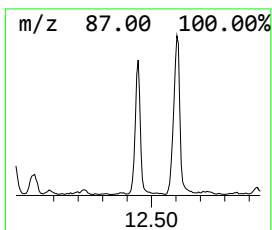
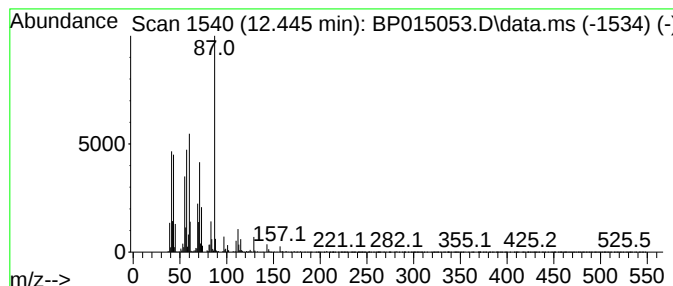
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Decanoic acid, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.445	4.02 ng/ul	449084	Naphthalene-d8	10.986

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decanoic acid, 3-methyl-	186	C11H22O2	060308-82-9	59
2			3-Methyldodecanoic acid	214	C13H26O2	013490-36-3	50
3			3-Methyloctanoic acid	158	C9H18O2	006061-10-5	50
4			5-Methylhexanoic acid	130	C7H14O2	000628-46-6	43
5			Diethylene glycol 2-ethylhexyl e...	260	C14H28O4	1000506-26-1	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

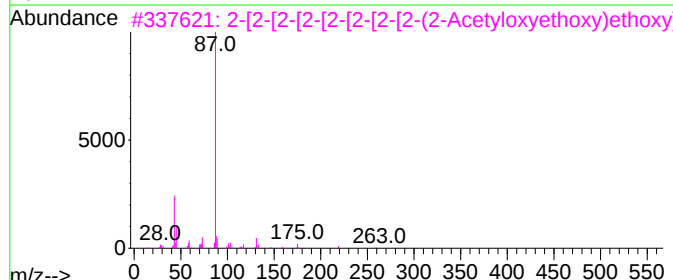
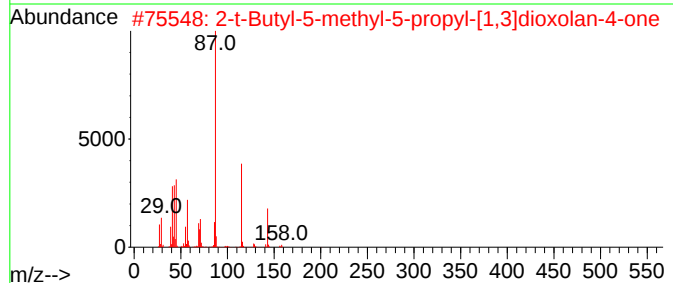
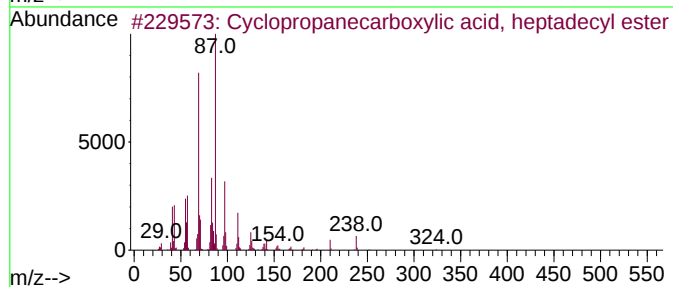
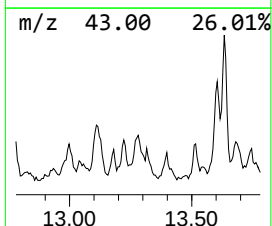
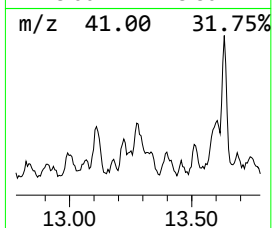
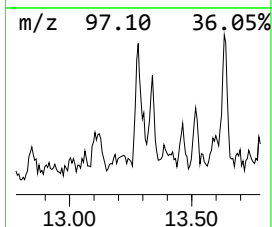
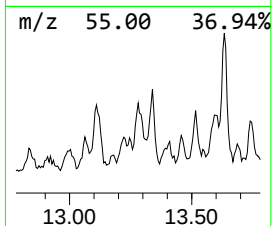
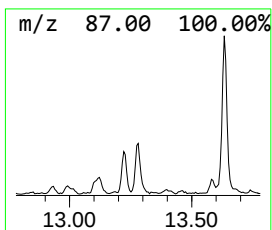
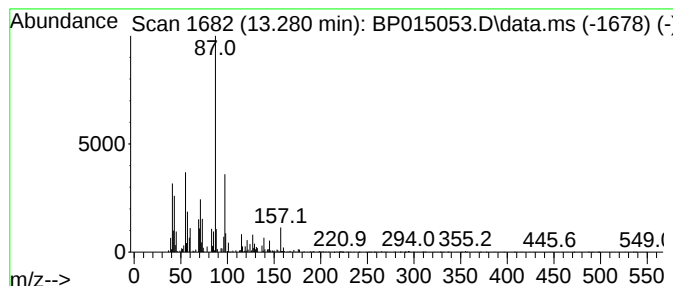
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Cyclopropanecarboxylic acid... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.281	2.02 ng/ul	245104	Acenaphthene-d10	14.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropanecarboxylic acid, hep...	324	C21H40O2	1000282-78-5	53
2			2-t-Butyl-5-methyl-5-propyl-[1,3...	200	C11H20O3	1000191-45-7	43
3			2-[2-[2-[2-[2-[2-(2-Acetyl...	498	C22H42O12	1010351-90-5	43
4			1,3-Dioxolane-2-propanol, 2-meth...	188	C9H16O4	029021-95-2	43
5			2,2-Dimethyl-5-(2-methyl-[1,3]di...	230	C12H22O4	1000185-52-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

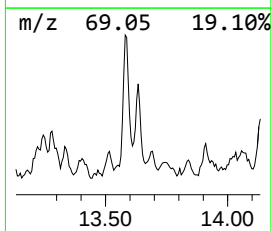
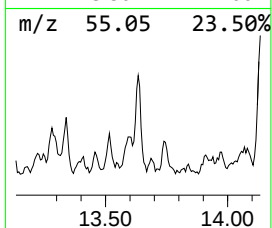
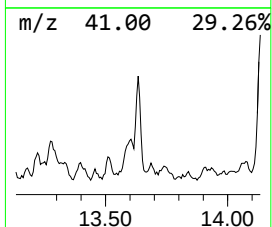
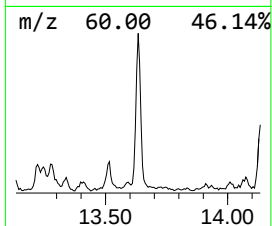
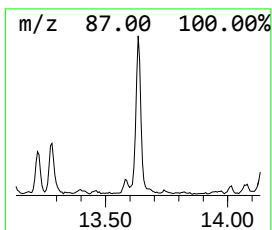
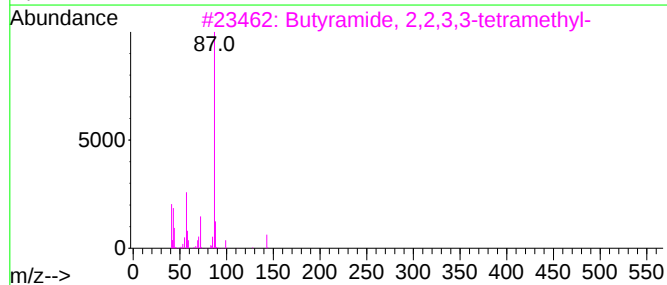
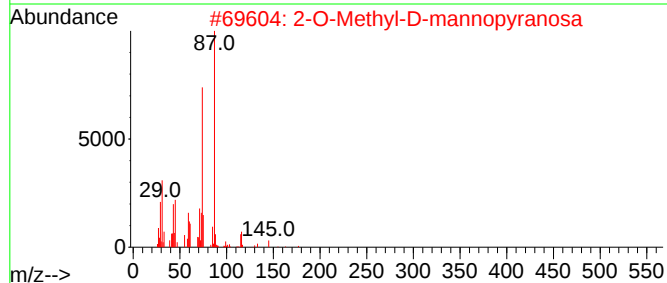
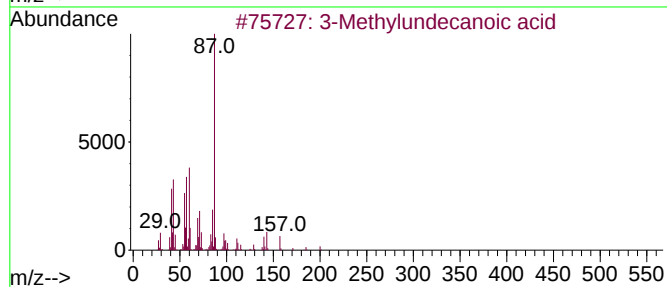
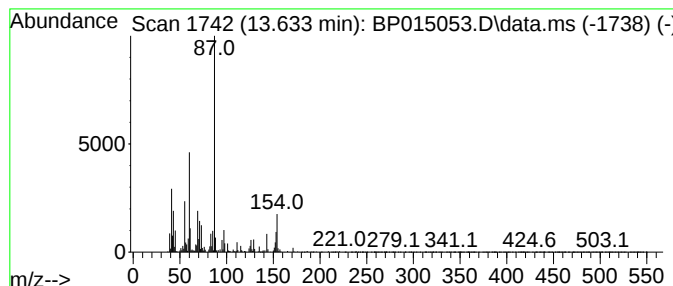
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 3-Methylundecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.633	3.61 ng/ul	438177	Acenaphthene-d10	14.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methylundecanoic acid	200	C12H24O2	065781-38-6	64
2			2-O-Methyl-D-mannopyranosa	194	C7H14O6	036864-61-6	38
3			Butyramide, 2,2,3,3-tetramethyl-	143	C8H17NO	098486-62-5	38
4			3-Methyloctanoic acid	158	C9H18O2	006061-10-5	38
5			.alpha.-D-Xylofuranoside, methyl...	178	C7H14O5	032469-86-6	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

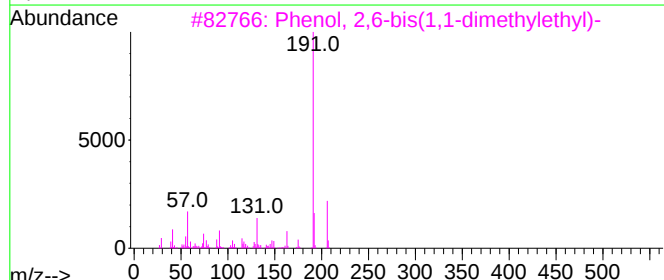
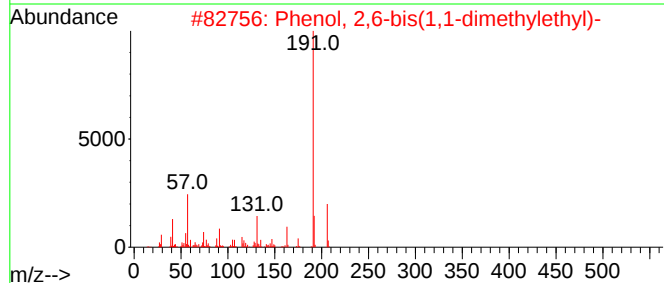
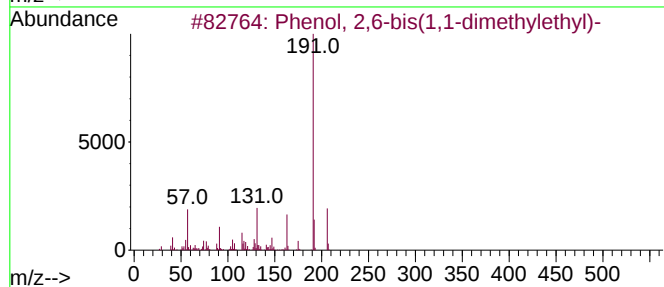
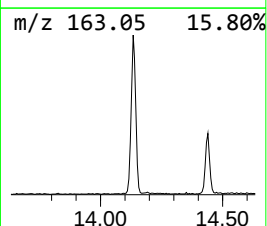
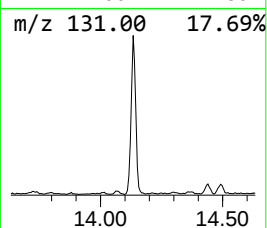
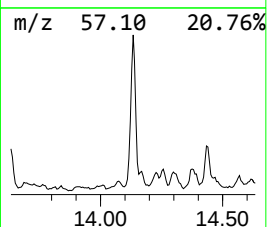
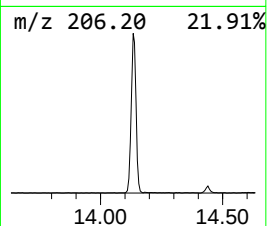
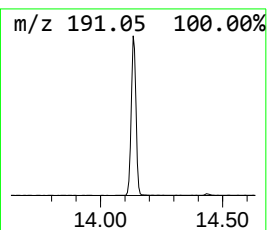
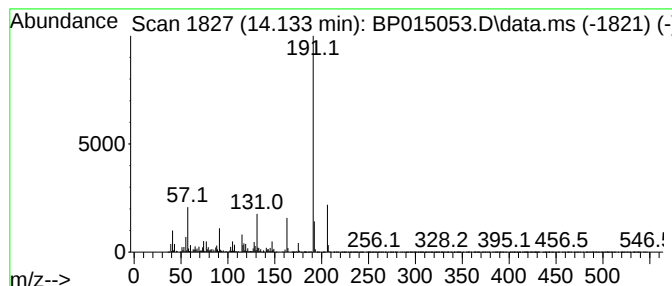
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.133	11.31 ng/ul	1373370	Acenaphthene-d10	14.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	98
2			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	94
3			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	93
4			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	93
5			Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

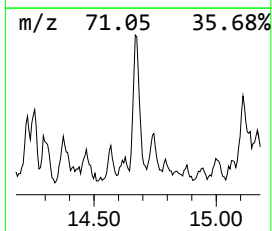
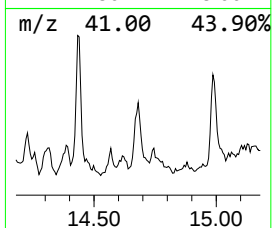
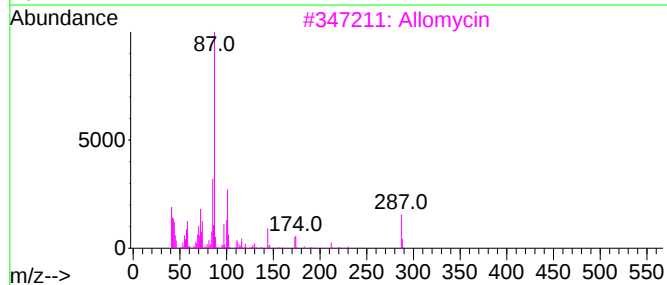
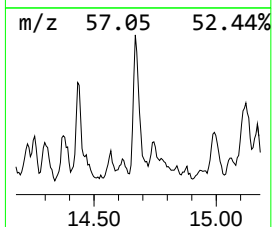
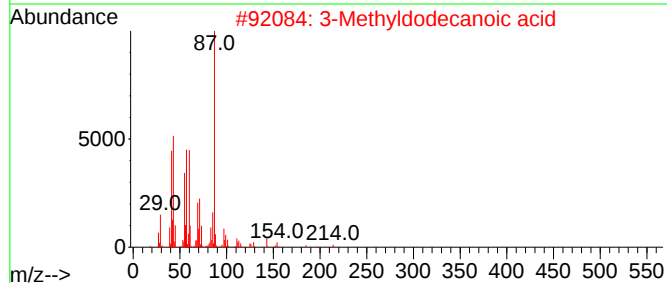
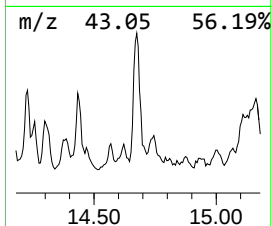
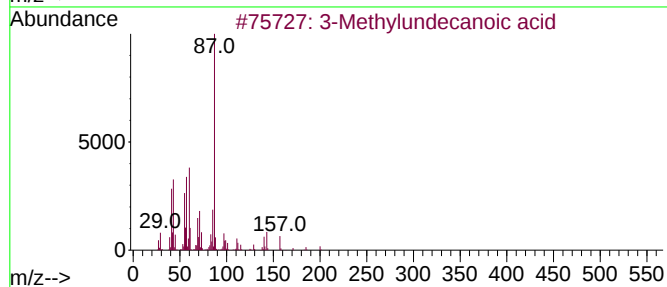
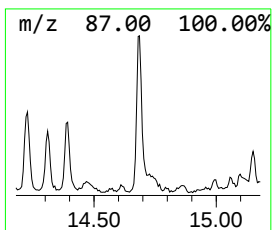
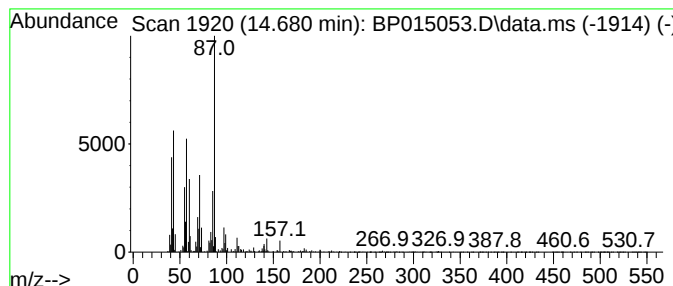
Instrument :
BNA_P
ClientSampleId :
JREJ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 3-Methyldodecanoic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.680	2.73 ng/ul	331411	Acenaphthene-d10		14.792	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Methylundecanoic acid		200	C12H24O2	065781-38-6	95
2	3-Methyldodecanoic acid		214	C13H26O2	013490-36-3	64
3	Allomycin		618	C29H42N6O9	017650-86-1	50
4	4-Nonanol, 4-methyl-		158	C10H22O	023418-38-4	43
5	2,2-Dimethyl-5-hexen-3-ol		128	C8H16O	019550-89-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

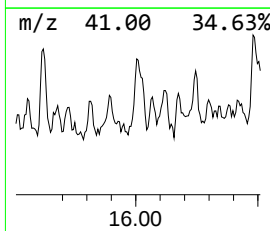
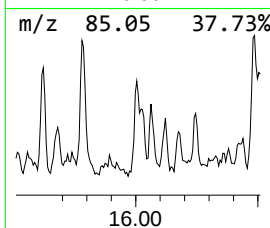
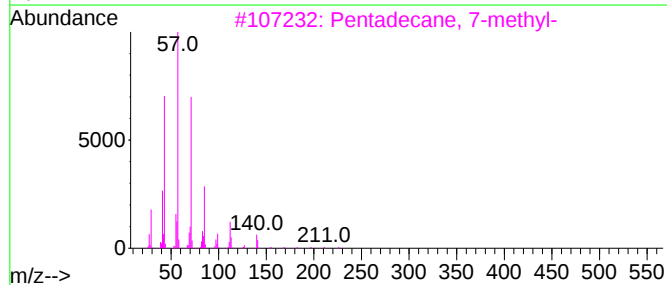
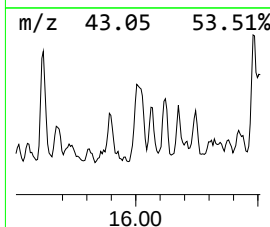
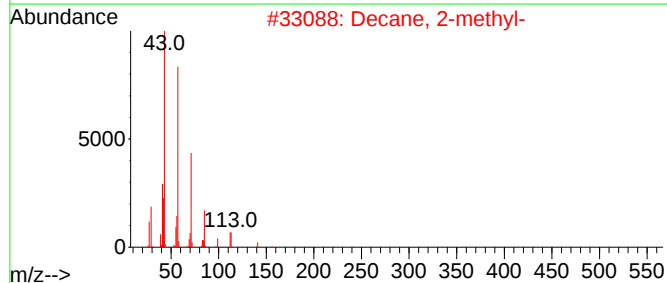
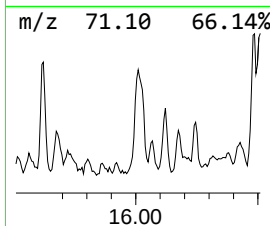
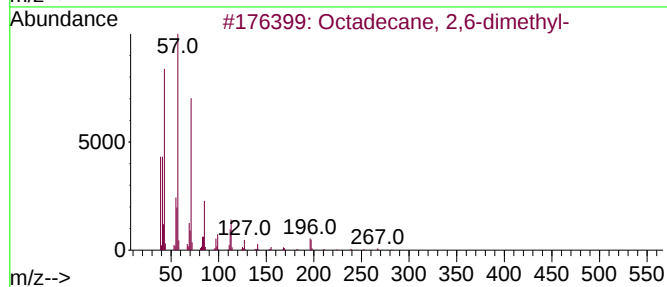
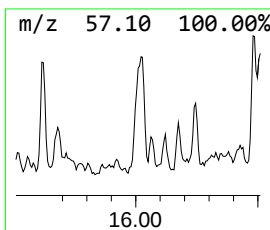
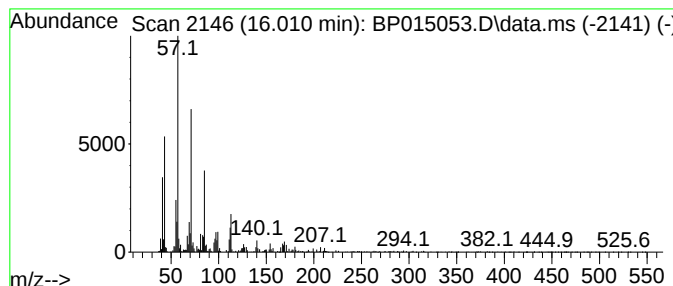
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ClientSampleId :
JREJ6

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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 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.010	2.37 ng/ul	287568	Acenaphthene-d10	14.792	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	90
2	Decane, 2-methyl-	156	C11H24	006975-98-0	87
3	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	87
4	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	83
5	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

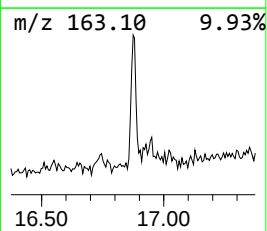
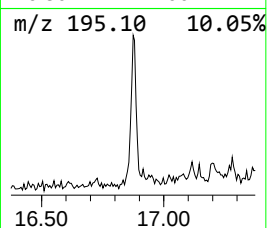
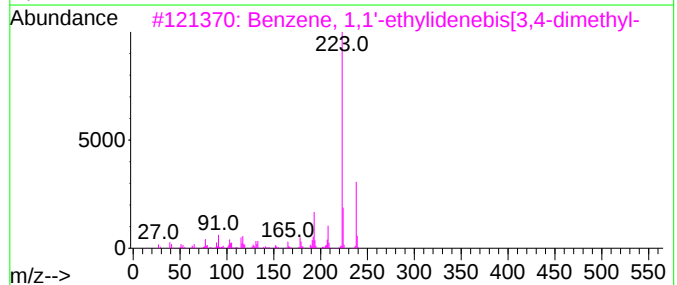
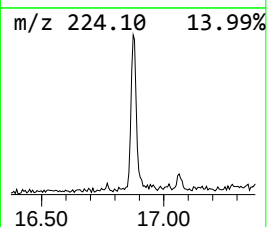
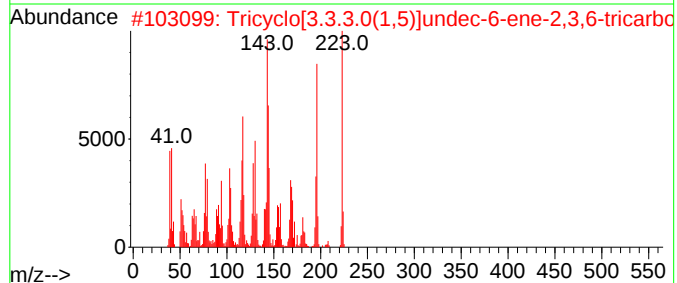
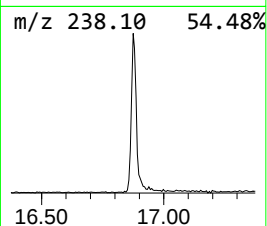
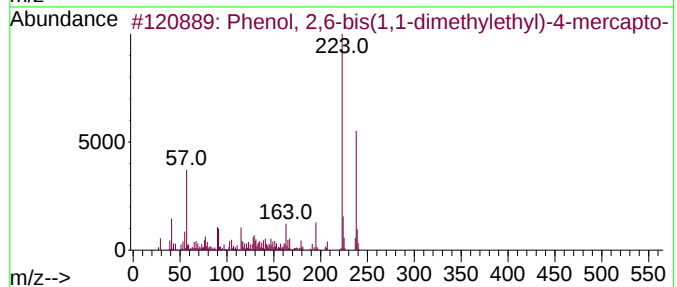
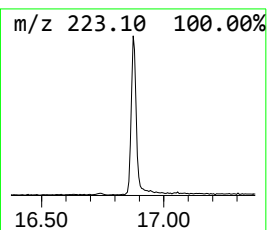
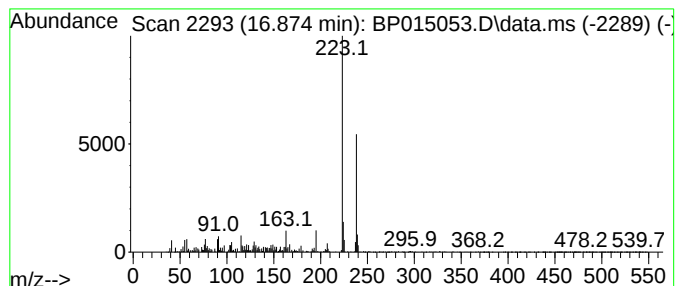
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Tricyclo[3.3.3.0(1,5)]undec... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.874	2.99 ng/ul	396443	Phenanthrene-d10	17.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-bis(1,1-dimethylethy...	238	C14H22OS	000950-59-4	91
2			Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	86
3			Benzene, 1,1'-ethyldienebis[3,4-...	238	C18H22	001742-14-9	72
4			4-tert-Butyl-benzophenone	238	C17H18O	022679-54-5	72
5			3H-Pyrrolo[3,2-f]quinoline, 1,2,...	238	C16H18N2	1000302-73-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

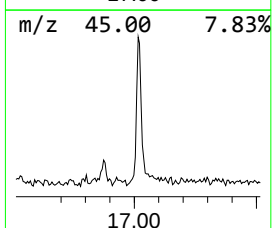
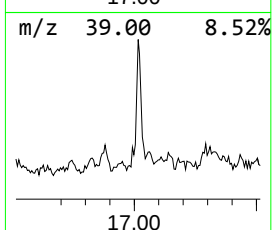
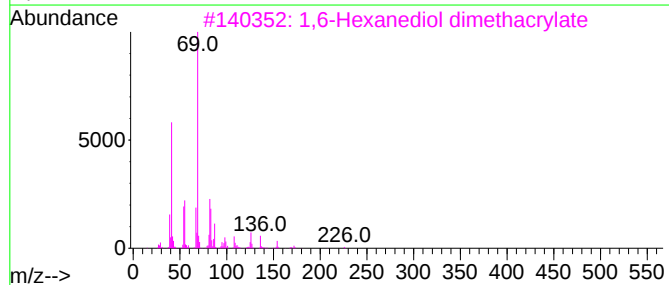
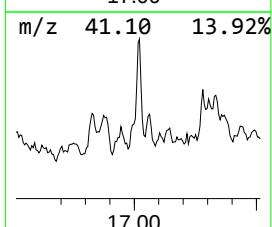
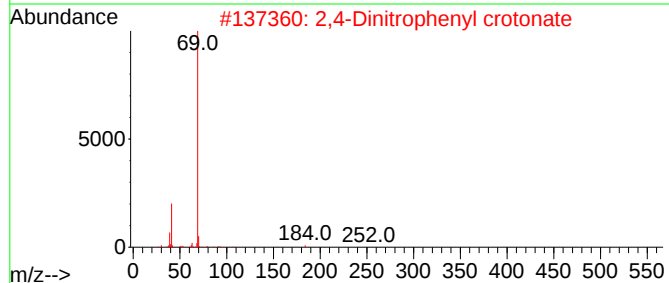
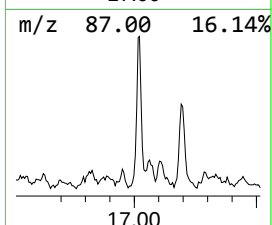
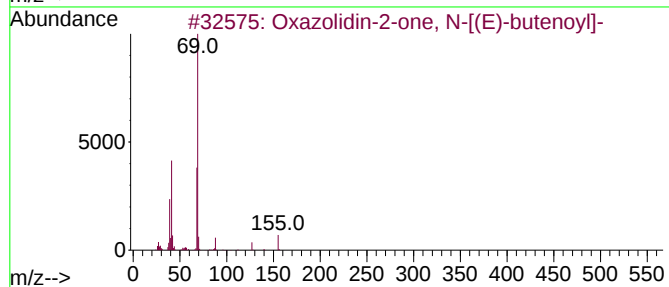
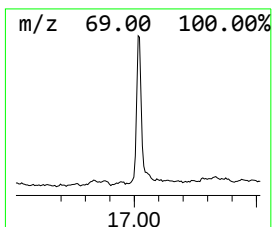
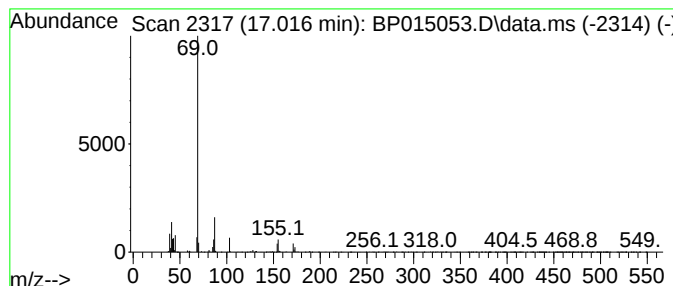
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TIC Integration Parameters: LSCINT.P

Peak Number 18 unknown-04

Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.016	2.66 ng/ul	352546	Phenanthrene-d10	17.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1010156-45-5	42
2			2,4-Dinitrophenyl crotonate	252	C10H8N2O6	069817-88-5	9
3			1,6-Hexanediol dimethacrylate	254	C14H22O4	006606-59-3	9
4			Cyclopropanecarboxylic acid, 2-m...	170	C10H18O2	1000354-65-8	9
5			Cyclopropanecarboxylic acid, buty...	142	C8H14O2	054947-39-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
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Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

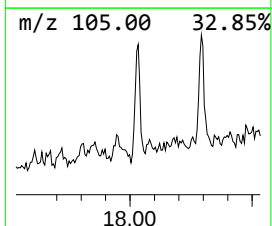
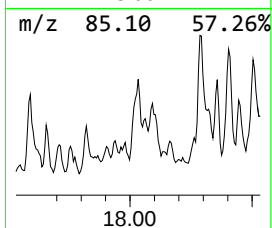
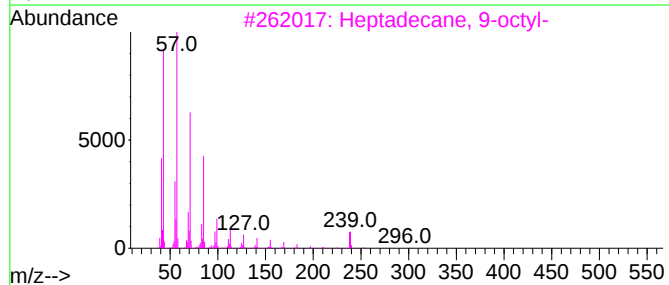
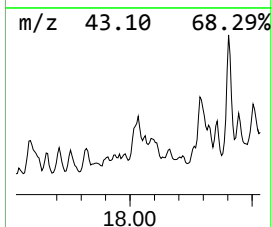
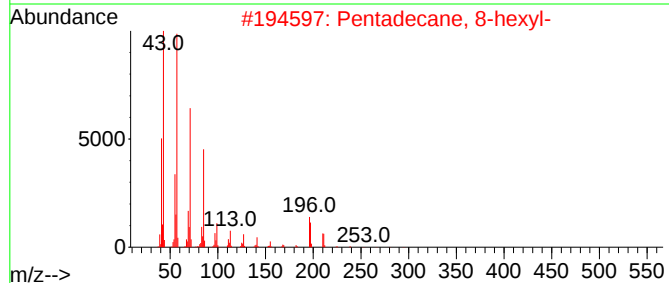
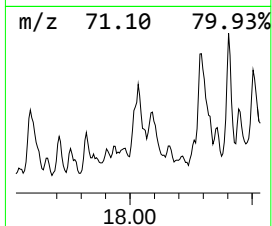
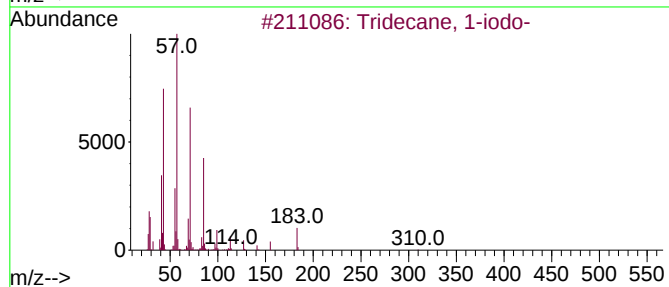
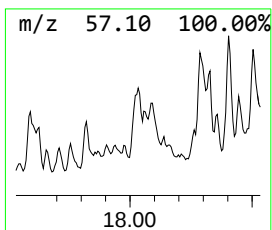
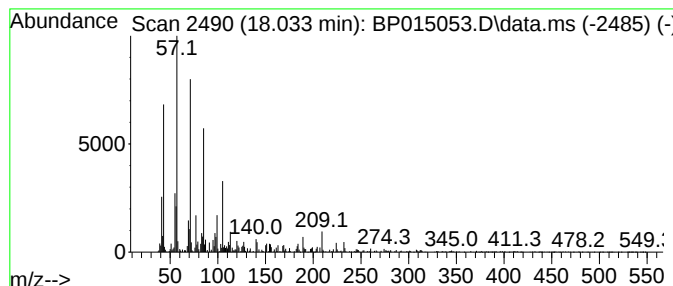
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ClientSampleId :
JREJ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 19 Tridecane, 1-iodo- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.033	3.23 ng/ul	428283	Phenanthrene-d10	17.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
2	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	89
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	76
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	76
5	Tridecane, 7-propyl-	226	C16H34	055045-09-5	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
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Acq On : 10 May 2023 21:05
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ALS Vial : 14 Sample Multiplier: 1

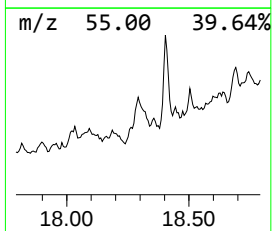
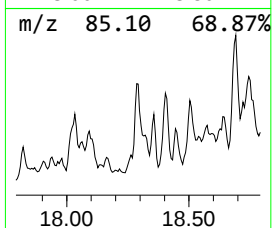
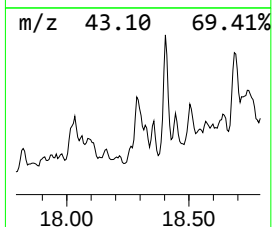
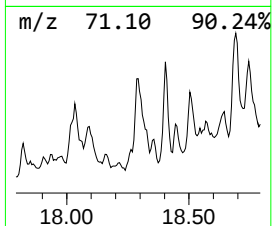
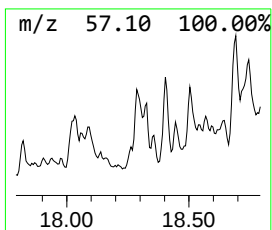
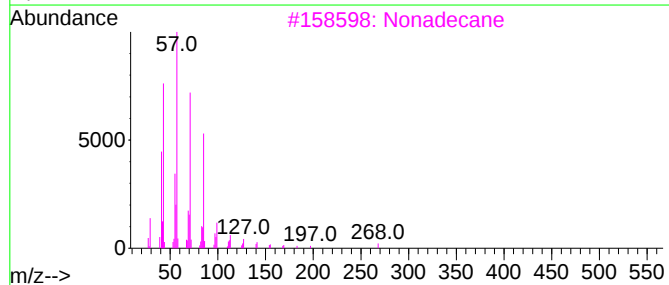
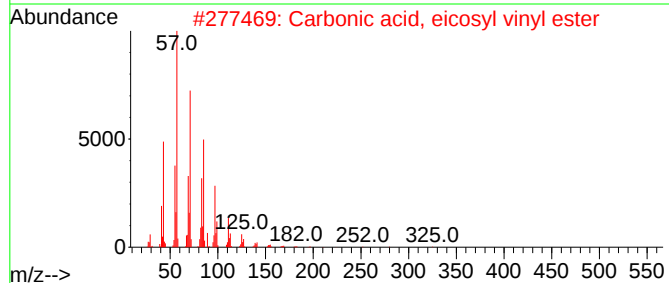
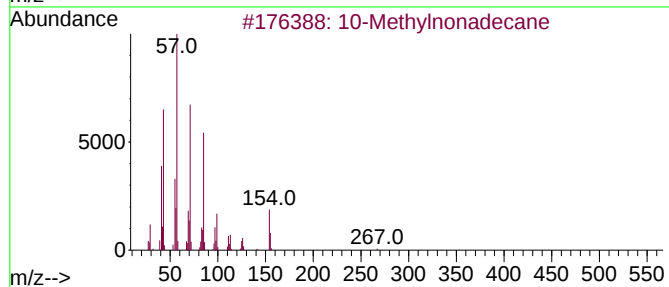
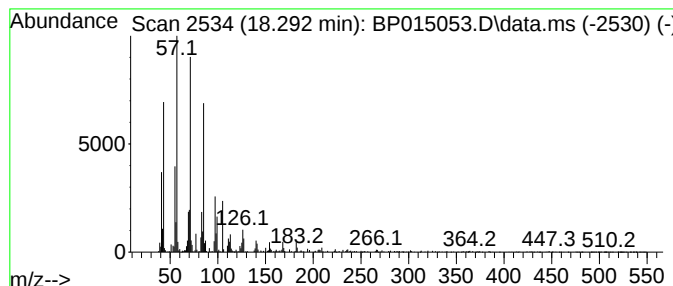
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.292	4.02 ng/ul	533974	Phenanthrene-d10	17.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane	282	C20H42	056862-62-5	83
2	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	80
3	Nonadecane	268	C19H40	000629-92-5	74
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	70
5	Hexacosane	366	C26H54	000630-01-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
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Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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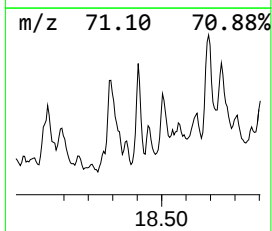
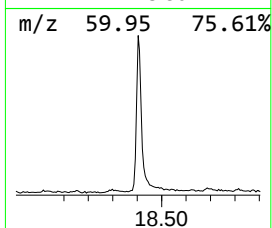
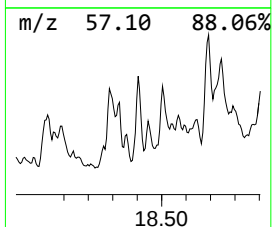
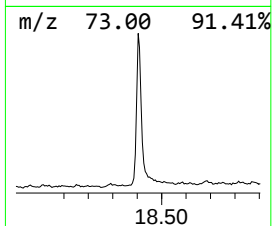
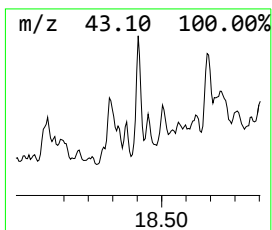
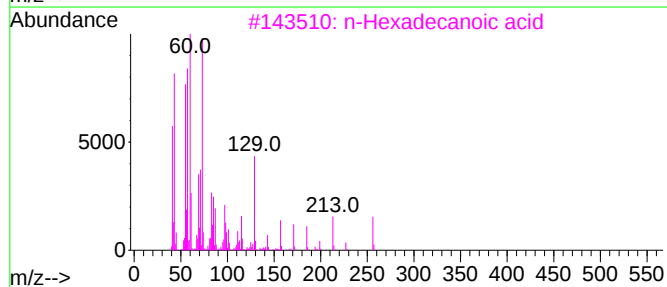
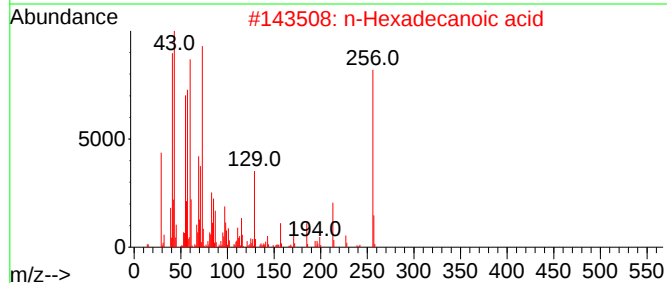
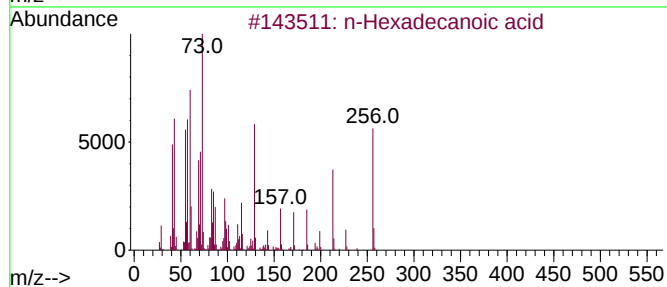
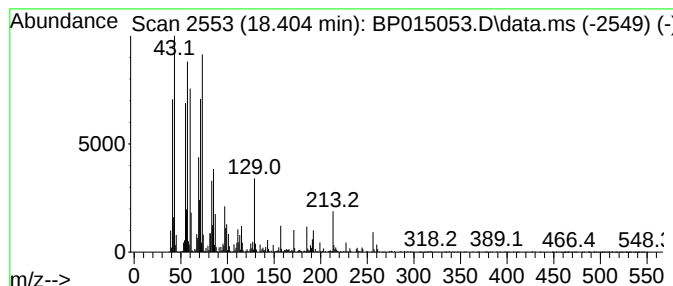
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	9.39 ng/ul	1246260	Phenanthrene-d10	17.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
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Acq On : 10 May 2023 21:05
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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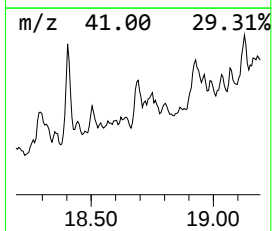
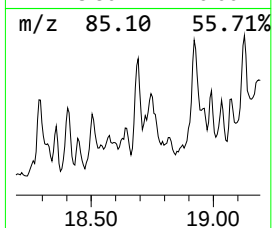
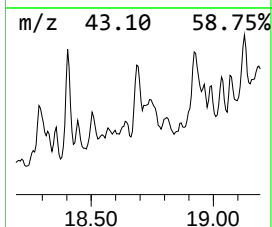
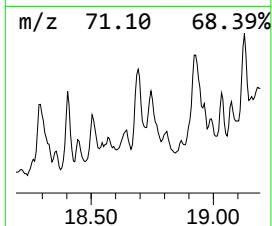
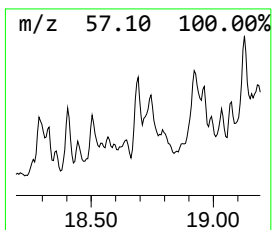
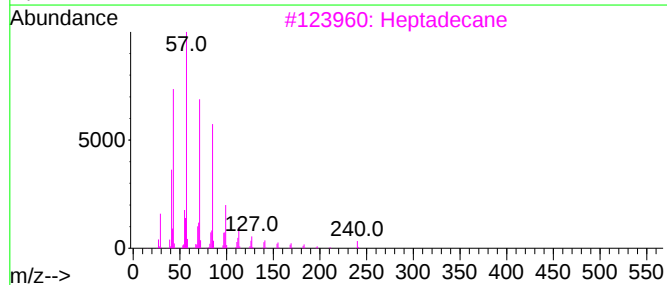
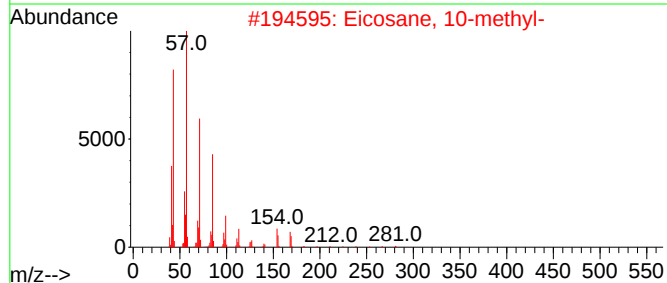
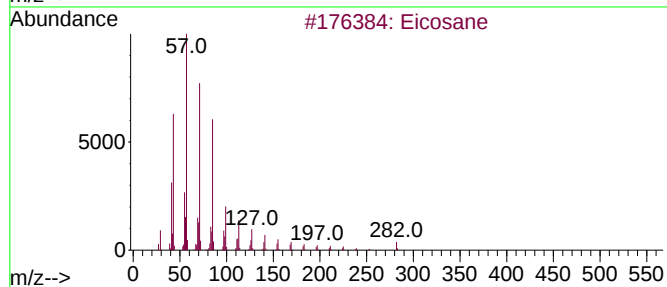
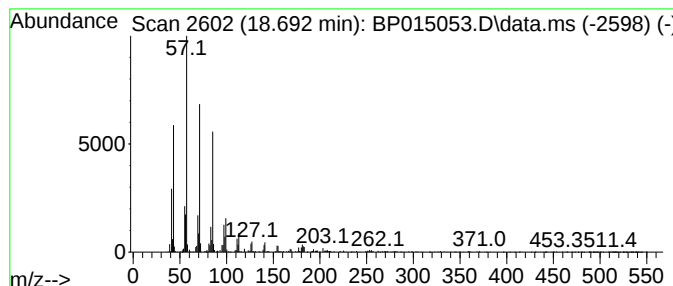
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.692	5.09 ng/ul	675534	Phenanthrene-d10	17.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	93
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
3		Heptadecane	240	C17H36	000629-78-7	87
4		Tetracosane	338	C24H50	000646-31-1	86
5		Octadecane	254	C18H38	000593-45-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JREJ6

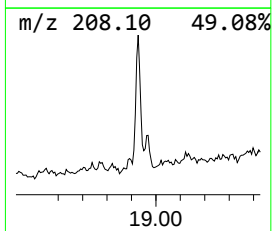
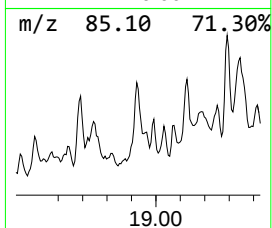
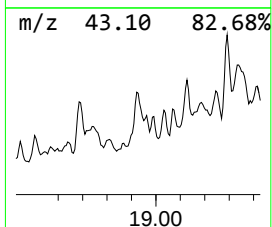
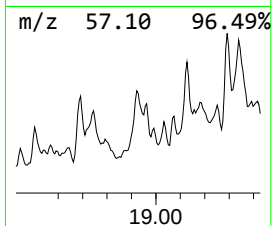
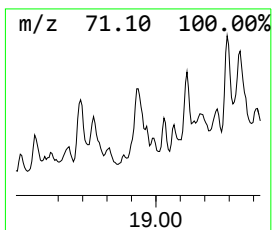
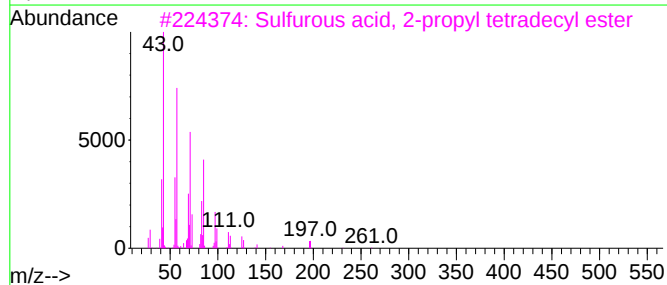
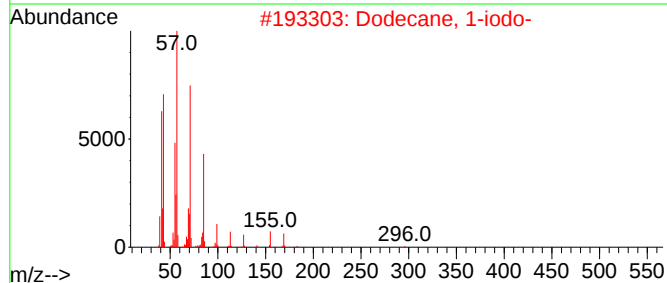
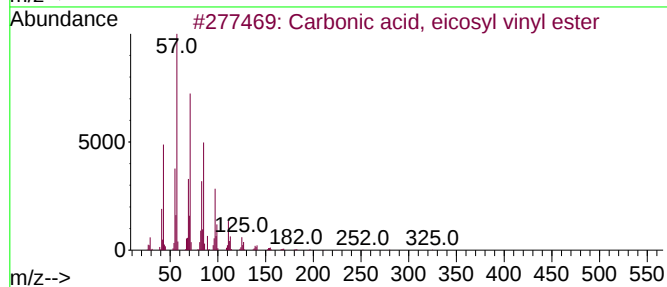
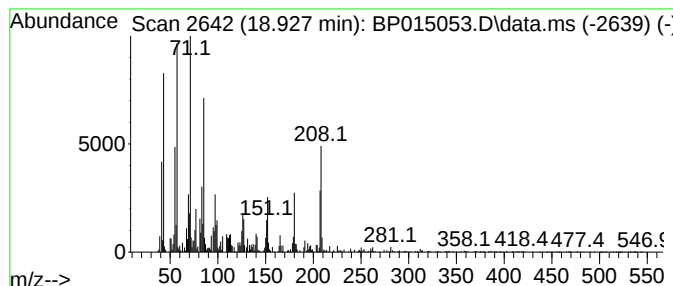
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.927	4.56 ng/ul	605077	Phenanthrene-d10	17.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	49
2			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	46
3			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	46
4			1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	46
5			Tridecane, 6-propyl-	226	C16H34	055045-10-8	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

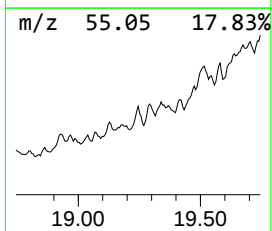
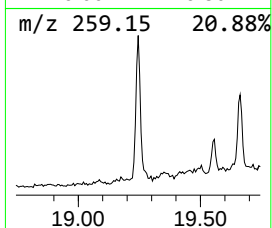
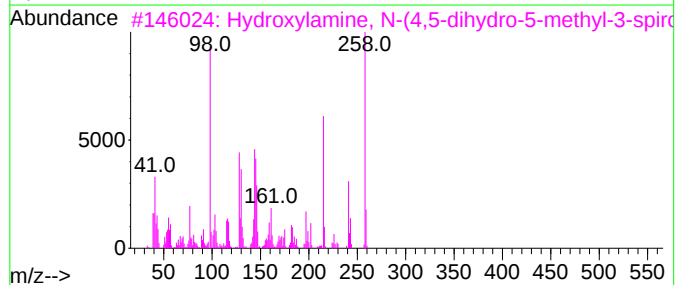
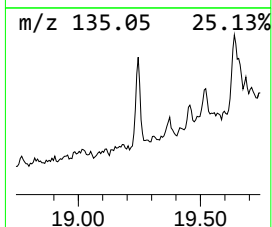
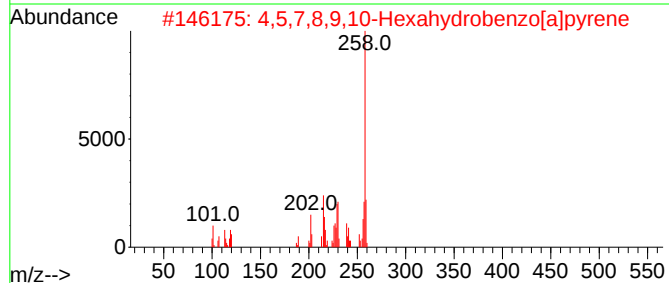
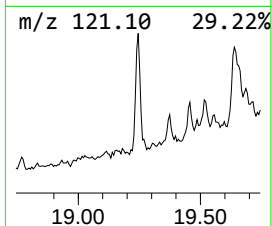
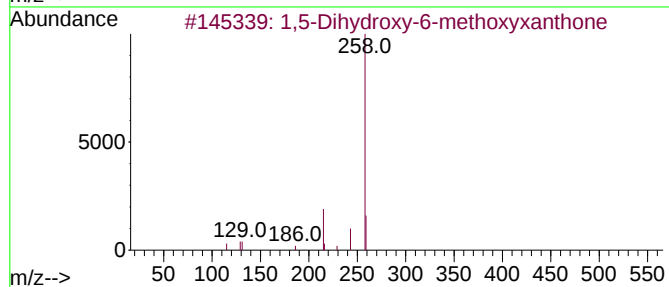
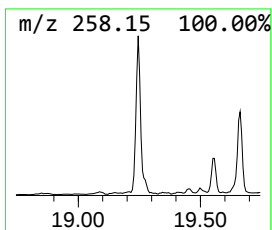
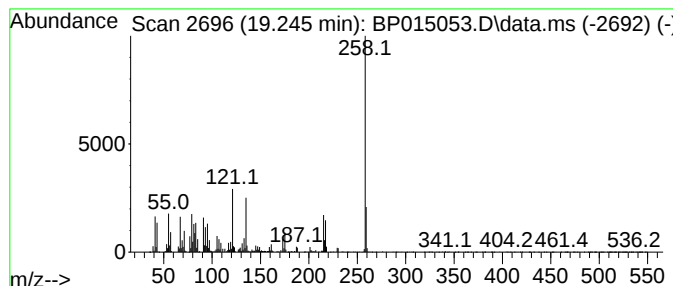
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 24 1,5-Dihydroxy-6-methoxyxant... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.245	9.68 ng/ul	1284980	Phenanthrene-d10	17.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,5-Dihydroxy-6-methoxyxanthone	258	C14H10O5	181307-35-7	58
2			4,5,7,8,9,10-Hexahydrobenzo[a]py...	258	C20H18	073712-75-1	53
3			Hydroxylamine, N-(4,5-dihydro-5-...	258	C16H22N2O	1000274-03-9	52
4			6-Methoxy-benzo[c]phenanthrene	258	C19H14O	004176-37-8	50
5			1,1'-Binaphthalene, 3,3',4,4'-te...	258	C20H18	005405-96-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015053.D
Acq On : 10 May 2023 21:05
Operator : CG/JU
Sample : 02591-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREJ6

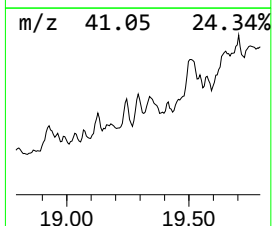
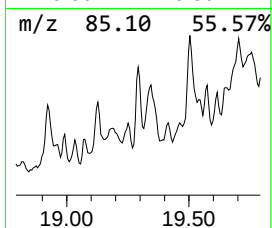
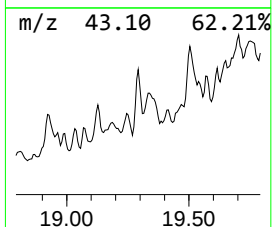
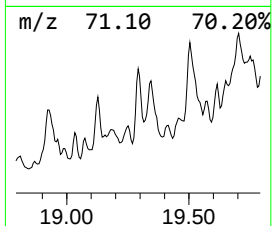
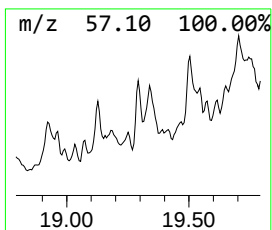
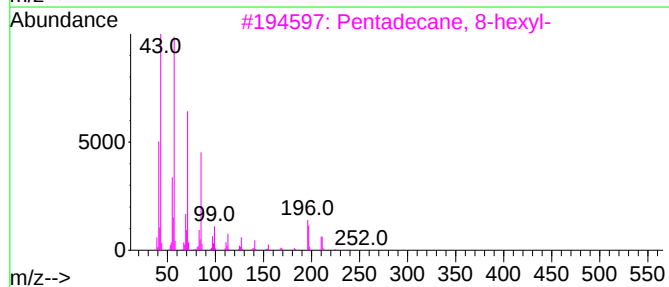
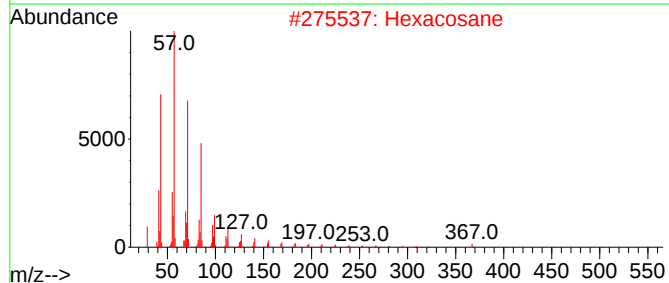
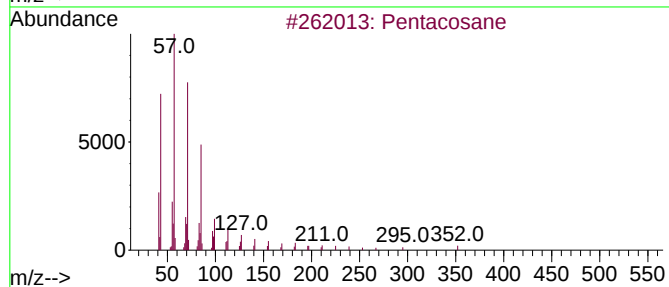
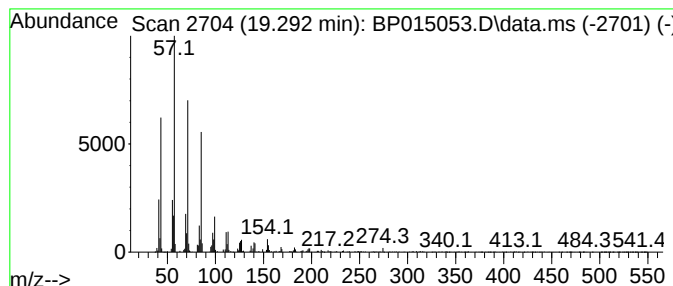
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.292	6.19 ng/ul	822274	Phenanthrene-d10	17.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentacosane	352	C25H52	000629-99-2	95
2		Hexacosane	366	C26H54	000630-01-3	90
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	89
5		Hexatriacontane	507	C36H74	000630-06-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
 Data File : BP015053.D
 Acq On : 10 May 2023 21:05
 Operator : CG/JU
 Sample : 02591-21
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREJ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.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
Crotonic acid	4.734	3.5	ng/ul	269201	1	8.151	1537820	20.0
3-Methyl-hexano...	8.063	3.4	ng/ul	257723	1	8.151	1537820	20.0
1-Hexanol, 2-et...	8.216	13.0	ng/ul	1000070	1	8.151	1537820	20.0
Hexanoic acid, ...	9.463	3.8	ng/ul	290290	1	8.151	1537820	20.0
unknown-01	9.581	4.3	ng/ul	484534	2	10.986	2231620	20.0
Benzoic acid	10.245	3.0	ng/ul	338810	2	10.986	2231620	20.0
unknown-02	10.492	3.5	ng/ul	392573	2	10.986	2231620	20.0
(DEL) Alkane: C...	10.816	4.9	ng/ul	548078	2	10.986	2231620	20.0
2-Fluoropyridine	11.704	2.0	ng/ul	227907	2	10.986	2231620	20.0
unknown-03	11.775	2.7	ng/ul	300456	2	10.986	2231620	20.0
Decanoic acid, ...	12.445	4.0	ng/ul	449084	2	10.986	2231620	20.0
Cyclopropanecar...	13.281	2.0	ng/ul	245104	3	14.792	2427760	20.0
3-Methylundecan...	13.633	3.6	ng/ul	438177	3	14.792	2427760	20.0
Phenol, 2,6-bis...	14.133	11.3	ng/ul	1373370	3	14.792	2427760	20.0
3-Methyldodecan...	14.680	2.7	ng/ul	331411	3	14.792	2427760	20.0
(DEL) Alkane: S...	16.010	2.4	ng/ul	287568	3	14.792	2427760	20.0
Tricyclo[3.3.3....	16.874	3.0	ng/ul	396443	4	17.551	2654910	20.0
unknown-04	17.016	2.7	ng/ul	352546	4	17.551	2654910	20.0
Tridecane, 1-iodo-	18.033	3.2	ng/ul	428283	4	17.551	2654910	20.0
(DEL) Alkane: S...	18.292	4.0	ng/ul	533974	4	17.551	2654910	20.0
n-Hexadecanoic ...	18.404	9.4	ng/ul	1246260	4	17.551	2654910	20.0
(DEL) Alkane: S...	18.692	5.1	ng/ul	675534	4	17.551	2654910	20.0
unknown-05	18.927	4.6	ng/ul	605077	4	17.551	2654910	20.0
1,5-Dihydroxy-6...	19.245	9.7	ng/ul	1284980	4	17.551	2654910	20.0
(DEL) Alkane: S...	19.292	6.2	ng/ul	822274	4	17.551	2654910	20.0