

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
 Data File : BP013339.D
 Acq On : 27 Dec 2022 15:26
 Operator : CG/JU
 Sample : N6133-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXGJ1

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013339.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.323	19	23	30	rVB	103055	136464	1.04%	0.091%
2	3.670	79	82	88	rVB	27460	37248	0.28%	0.025%
3	3.752	92	96	107	rBV	734050	1090193	8.31%	0.726%
4	3.981	130	135	139	rVB	36151	57467	0.44%	0.038%
5	4.075	147	151	156	rVB	60676	76323	0.58%	0.051%
6	5.022	306	312	319	rVB	256965	361868	2.76%	0.241%
7	5.193	336	341	347	rVV	121567	173185	1.32%	0.115%
8	5.681	420	424	433	rBV	83572	159667	1.22%	0.106%
9	6.534	563	569	572	rBV4	23998	39531	0.30%	0.026%
10	7.093	653	664	673	rBV	487907	788483	6.01%	0.525%
11	7.205	679	683	686	rBV	26705	31687	0.24%	0.021%
12	7.258	686	692	699	rBV	1741283	2608109	19.89%	1.737%
13	7.458	716	726	736	rBV	1662836	2619873	19.98%	1.745%
14	7.769	768	779	786	rVB2	55607	209077	1.59%	0.139%
15	7.928	798	806	812	rBV	1892158	3000646	22.89%	1.999%
16	7.993	812	817	833	rVB	997405	1740214	13.27%	1.159%
17	8.181	844	849	854	rVB5	20507	40414	0.31%	0.027%
18	8.640	921	927	937	rBV	1100263	1786500	13.63%	1.190%
19	9.022	986	992	996	rBV7	21748	43959	0.34%	0.029%
20	9.099	998	1005	1014	rVV	2087750	3293007	25.12%	2.194%
21	9.181	1014	1019	1020	rVV3	27650	37446	0.29%	0.025%
22	9.246	1020	1030	1042	rVB	714204	1481937	11.30%	0.987%
23	9.499	1067	1073	1077	rVV	78819	128693	0.98%	0.086%
24	9.540	1077	1080	1088	rVB8	18821	41849	0.32%	0.028%
25	9.822	1121	1128	1141	rBV	1674369	2894991	22.08%	1.928%
26	9.916	1142	1144	1150	rVV7	17954	35198	0.27%	0.023%
27	9.981	1150	1155	1164	rVB3	55367	102942	0.79%	0.069%
28	10.052	1164	1167	1176	rVB10	19197	37472	0.29%	0.025%
29	10.157	1179	1185	1188	rBV7	20491	35811	0.27%	0.024%
30	10.363	1213	1220	1237	rVB	2613132	4373594	33.36%	2.913%
31	10.522	1241	1247	1257	rBV2	100387	195596	1.49%	0.130%
32	10.640	1259	1267	1278	rVV4	53486	157512	1.20%	0.105%
33	10.746	1278	1285	1291	rVV	2725708	4466408	34.06%	2.975%
34	10.799	1291	1294	1299	rVB3	53799	76247	0.58%	0.051%
35	10.881	1301	1308	1319	rBV	772897	1357795	10.36%	0.904%
36	11.210	1359	1364	1370	rVB5	24221	38961	0.30%	0.026%

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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-BP120722.M
 Title : SVOA CALIBRATION

37	11.275	1370	1375	1382	rBV	48370	90789	0.69%	0.060%
38	11.399	1390	1396	1405	rVB	232901	414193	3.16%	0.276%
39	11.563	1417	1424	1433	rVB	262569	445831	3.40%	0.297%
40	11.993	1491	1497	1503	rBV5	34744	55900	0.43%	0.037%
41	12.151	1519	1524	1529	rVB2	36519	65931	0.50%	0.044%
42	12.234	1529	1538	1546	rBV	489761	828798	6.32%	0.552%
43	12.328	1548	1554	1559	rBV2	45298	81390	0.62%	0.054%
44	12.516	1580	1586	1596	rBV6	33848	82978	0.63%	0.055%
45	12.734	1610	1623	1628	rBV	156606	340986	2.60%	0.227%
46	12.910	1646	1653	1654	rBV7	21130	37532	0.29%	0.025%
47	13.181	1694	1699	1712	rVB5	72140	165629	1.26%	0.110%
48	13.328	1717	1724	1729	rBV5	56791	135596	1.03%	0.090%
49	13.387	1729	1734	1742	rVV2	71671	129231	0.99%	0.086%
50	13.463	1742	1747	1756	rVV	164990	273527	2.09%	0.182%
51	13.681	1778	1784	1789	rBV8	19193	45990	0.35%	0.031%
52	13.840	1807	1811	1816	rBV	102959	144861	1.10%	0.096%
53	13.910	1816	1823	1829	rVV	566815	1254361	9.57%	0.836%
54	13.981	1829	1835	1841	rVV	5247003	7379122	56.28%	4.915%
55	14.028	1841	1843	1850	rVV4	89055	141826	1.08%	0.094%
56	14.098	1850	1855	1861	rVB7	42219	78751	0.60%	0.052%
57	14.216	1867	1875	1878	rBV5	48535	98386	0.75%	0.066%
58	14.269	1878	1884	1893	rVB	5459264	8058718	61.46%	5.368%
59	14.398	1903	1906	1909	rVB2	31878	36357	0.28%	0.024%
60	14.434	1909	1912	1914	rBV	43468	52241	0.40%	0.035%
61	14.575	1930	1936	1945	rVB	4219197	6164980	47.02%	4.107%
62	14.781	1965	1971	1982	rBV	418826	708746	5.41%	0.472%
63	14.928	1992	1996	2002	rBV6	45718	79199	0.60%	0.053%
64	14.981	2002	2005	2015	rVV8	55613	139395	1.06%	0.093%
65	15.063	2016	2019	2023	rVB6	25407	37764	0.29%	0.025%
66	15.151	2029	2034	2042	rBV	71729	132404	1.01%	0.088%
67	15.316	2056	2062	2066	rBV	363171	491566	3.75%	0.327%
68	15.357	2066	2069	2071	rVV2	65838	93491	0.71%	0.062%
69	15.392	2071	2075	2082	rVB	170682	255067	1.95%	0.170%
70	15.569	2099	2105	2112	rBV	7637142	10607624	80.90%	7.066%
71	15.692	2120	2126	2140	rBV	2653584	3677075	28.04%	2.449%
72	15.810	2143	2146	2149	rBV3	18274	30740	0.23%	0.020%
73	15.934	2162	2167	2173	rBV2	177179	311236	2.37%	0.207%
74	16.081	2186	2192	2196	rBV	52175	86298	0.66%	0.057%
75	16.216	2211	2215	2219	rBV5	67260	95081	0.73%	0.063%
76	16.304	2227	2230	2238	rVB4	41878	71379	0.54%	0.048%

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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-BP120722.M
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77	16.398	2242	2246	2255	rVB	117820	171538	1.31%	0.114%
78	16.481	2255	2260	2262	rBV3	120156	195669	1.49%	0.130%
79	16.522	2262	2267	2274	rVB	1044307	1379919	10.52%	0.919%
80	16.722	2297	2301	2307	rBV8	36095	58650	0.45%	0.039%
81	16.998	2345	2348	2354	rVB2	40391	51016	0.39%	0.034%
82	17.081	2358	2362	2365	rBV4	38855	56046	0.43%	0.037%
83	17.122	2365	2369	2373	rVB4	48310	65896	0.50%	0.044%
84	17.181	2375	2379	2382	rBV	106400	162411	1.24%	0.108%
85	17.334	2399	2405	2411	rBV	5280632	7122573	54.32%	4.745%
86	17.434	2416	2422	2430	rBV2	8103303	11295867	86.15%	7.525%
87	17.522	2431	2437	2451	rBV	3873208	5568950	42.47%	3.710%
88	17.992	2513	2517	2520	rBV4	56985	66039	0.50%	0.044%
89	18.039	2520	2525	2531	rVV2	562091	737744	5.63%	0.491%
90	18.210	2548	2554	2563	rBV	2024700	3093392	23.59%	2.061%
91	18.722	2639	2641	2645	rBV5	25598	31247	0.24%	0.021%
92	19.239	2726	2729	2735	rVB2	119470	166609	1.27%	0.111%
93	19.310	2737	2741	2745	rBV	117461	197893	1.51%	0.132%
94	19.439	2758	2763	2776	rBV	741046	1377220	10.50%	0.917%
95	19.669	2796	2802	2807	rBV	10138423	13111488	100.00%	8.734%
96	19.886	2835	2839	2846	rVB	156173	216930	1.65%	0.145%
97	21.116	3044	3048	3055	rBV	893020	1428671	10.90%	0.952%
98	21.422	3095	3100	3105	rBV	5368764	6988922	53.30%	4.656%
99	23.733	3485	3493	3508	rVB2	5983250	12281579	93.67%	8.181%
100	23.892	3513	3520	3530	rVB2	3497712	7116765	54.28%	4.741%

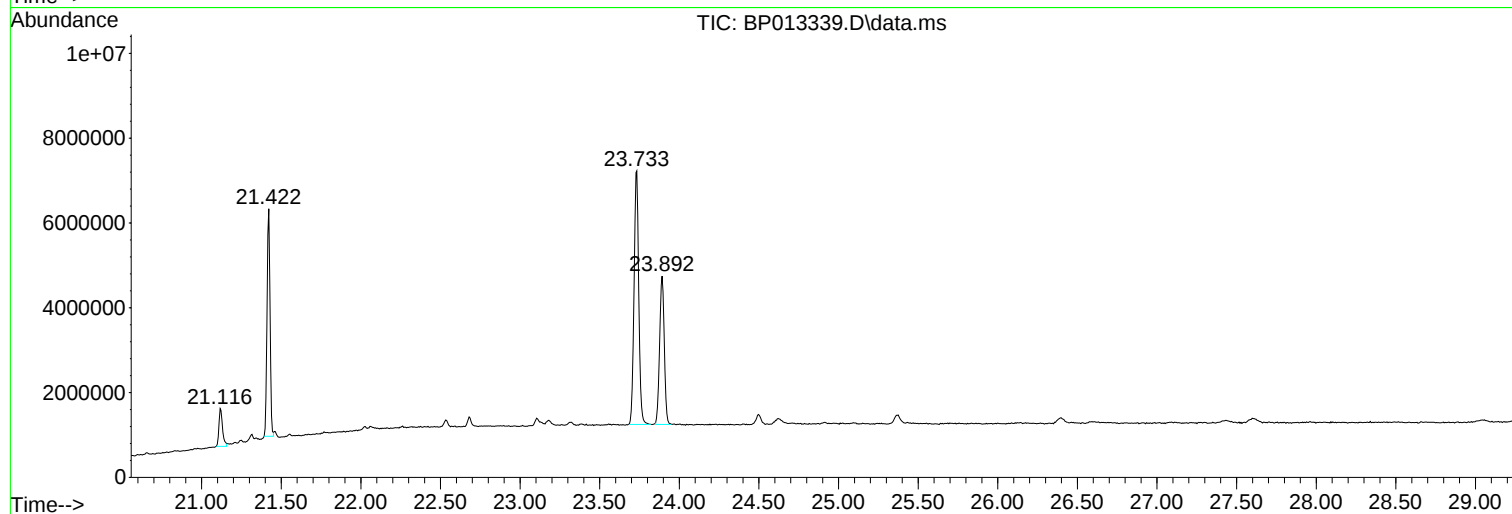
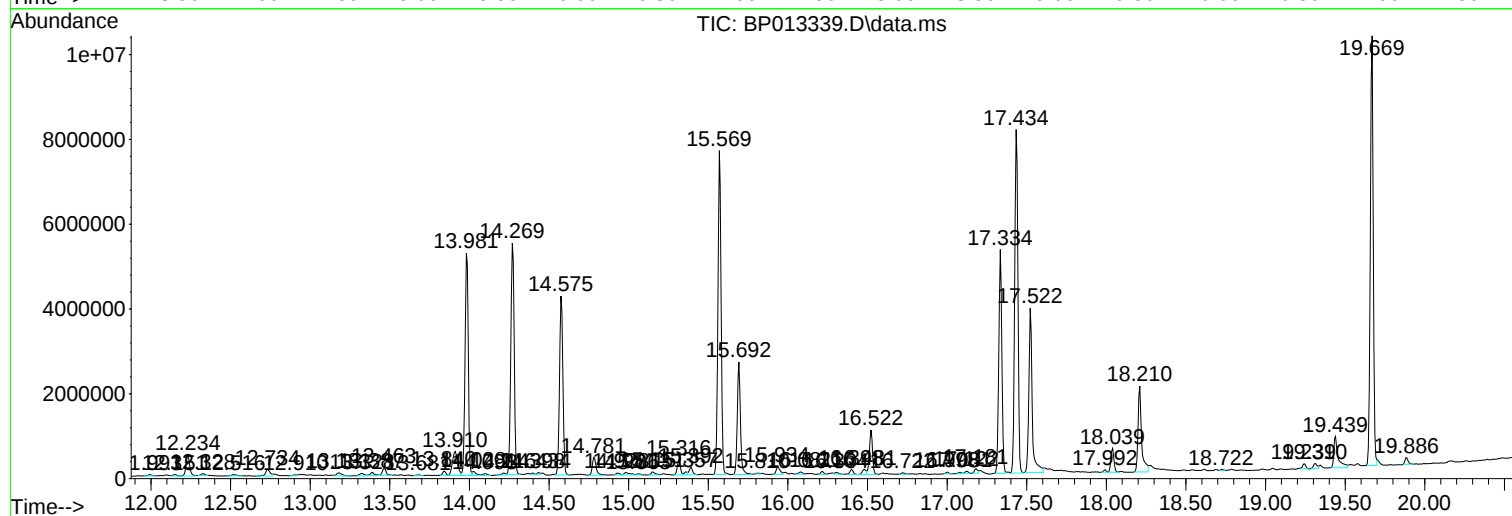
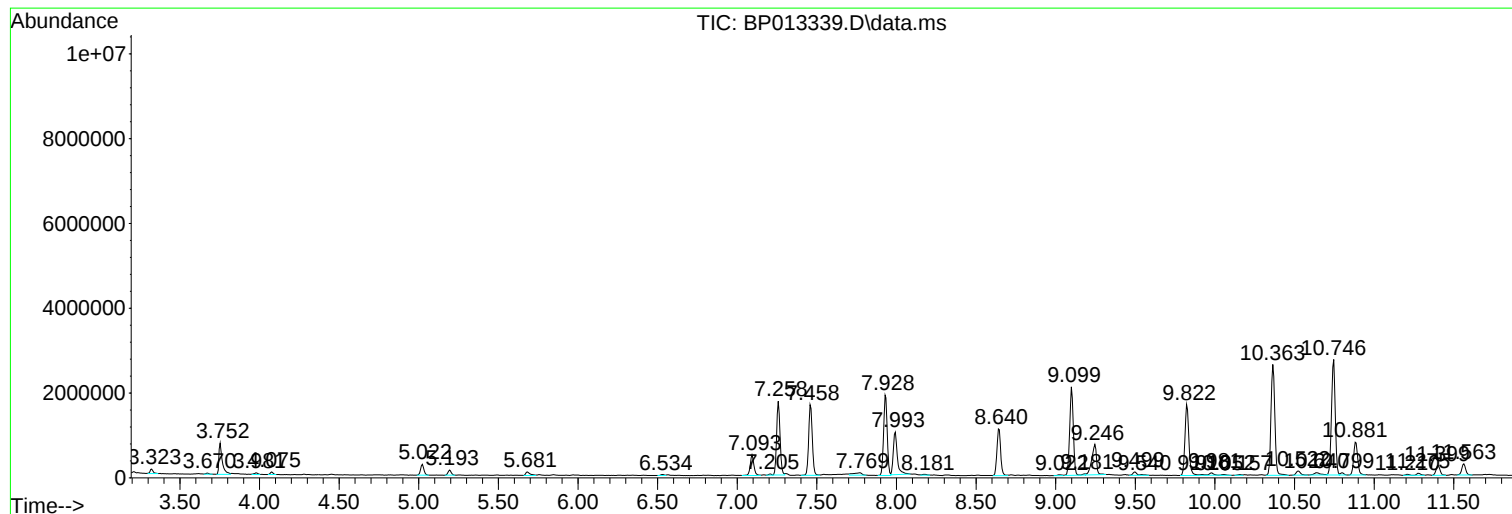
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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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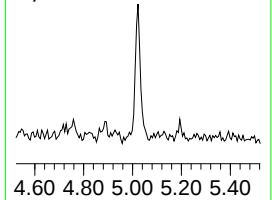
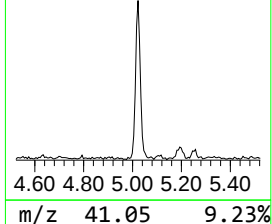
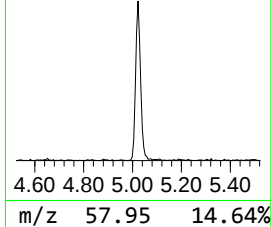
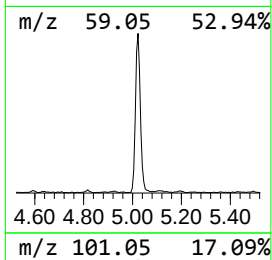
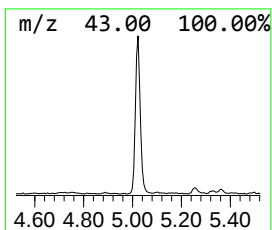
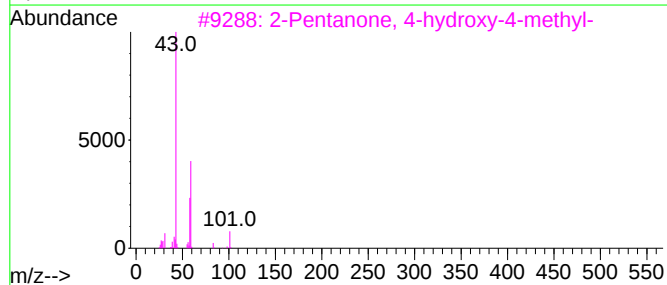
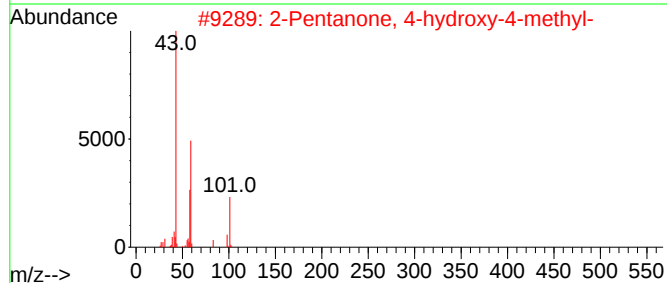
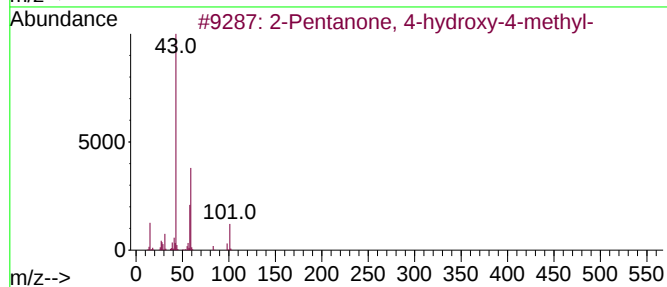
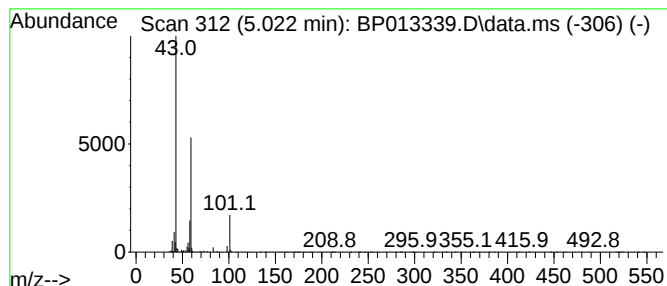
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.022	2.41 ng/ul	361868	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	37		
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25		



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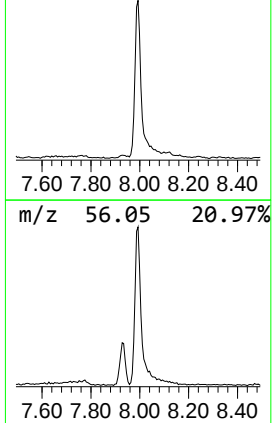
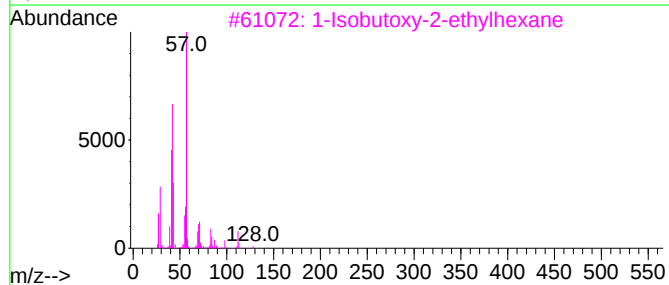
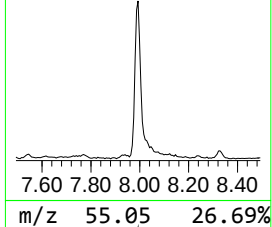
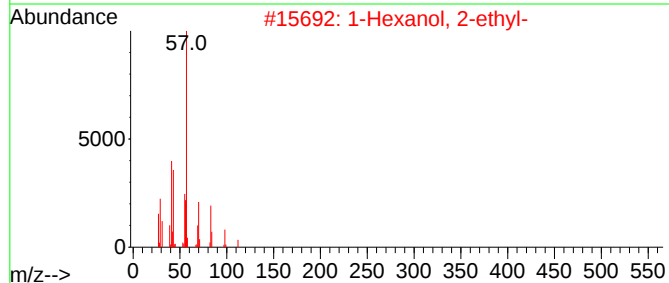
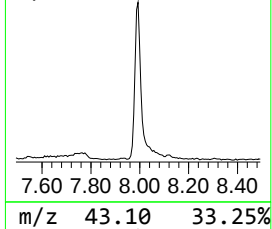
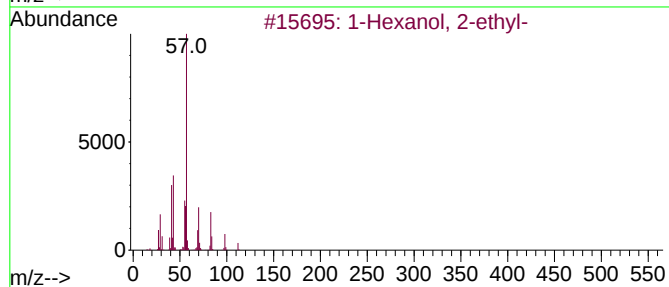
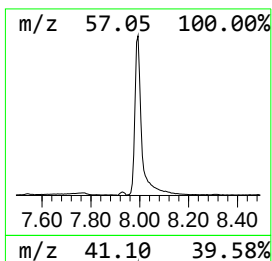
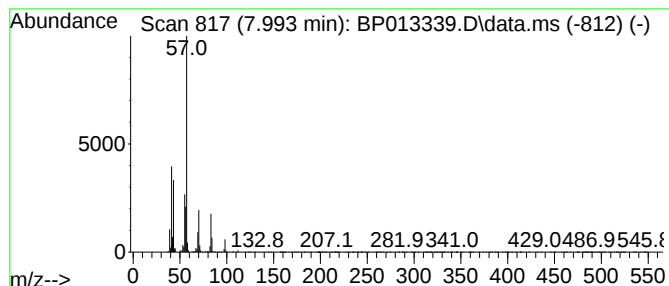
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	11.60 ng/ul	1740210	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	74
3	1-Isobutoxy-2-ethylhexane			186	C12H26O	1000139-90-3	56
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	45
5	Pentadecafluorooctanoic acid, 2-...			526	C16H17F15O2	1000406-80-9	45



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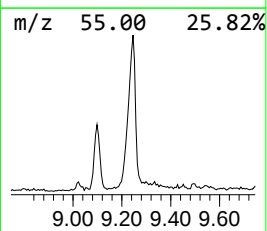
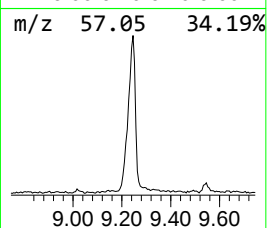
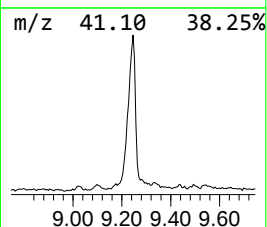
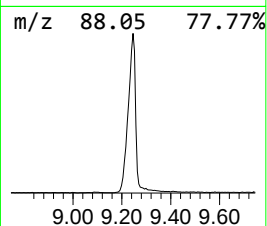
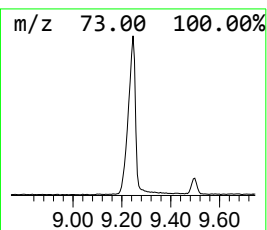
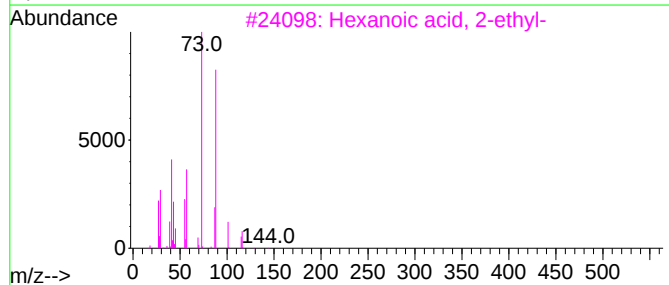
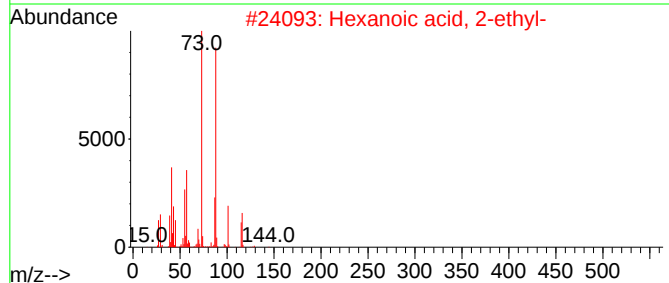
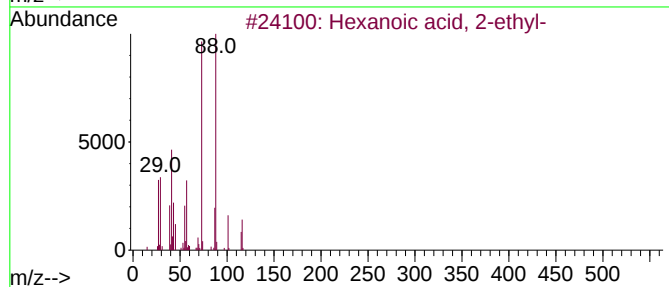
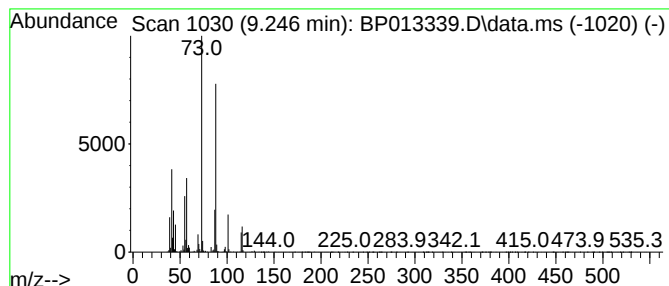
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Hexanoic acid, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.246	9.88 ng/ul	1481940	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
Operator : CG/JU
Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXGJ1

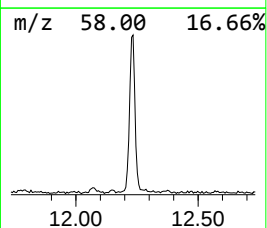
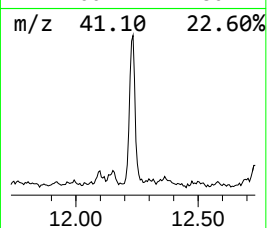
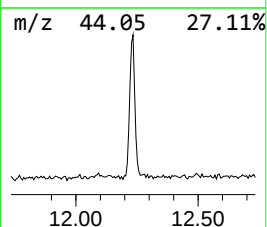
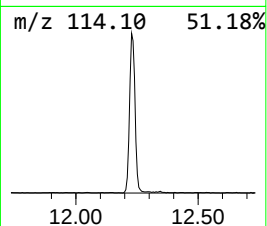
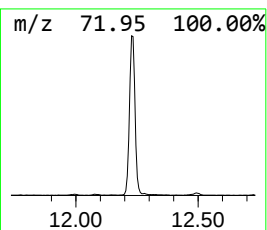
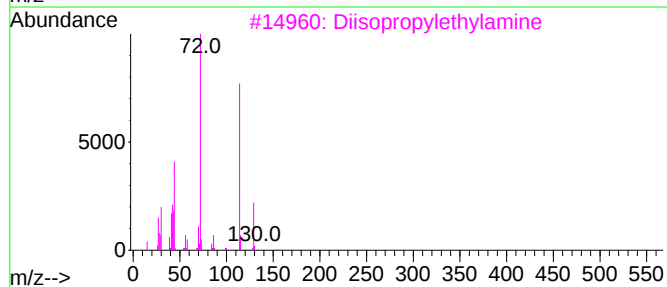
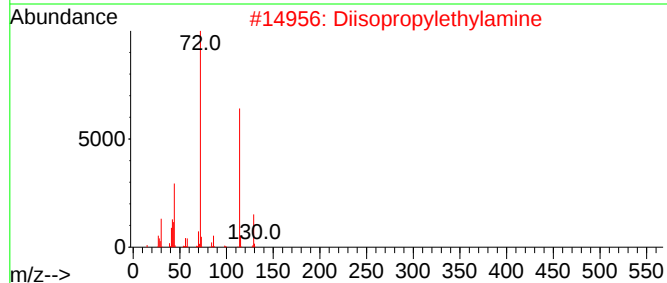
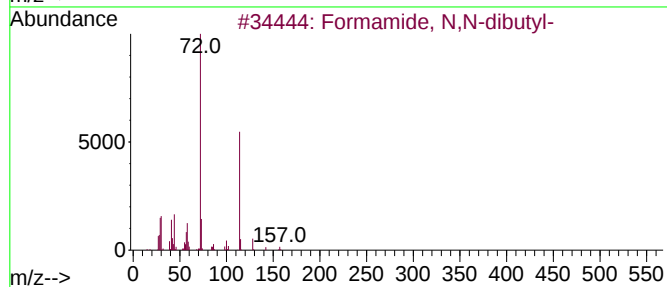
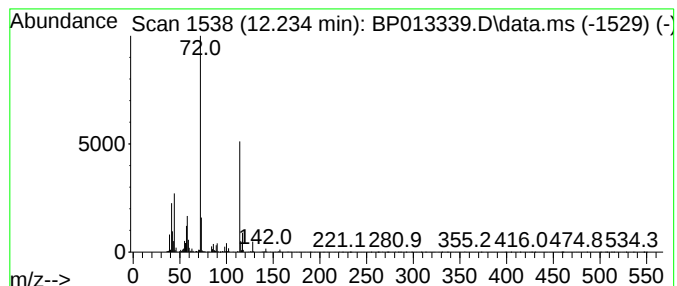
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Formamide, N,N-dibutyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.234	3.71 ng/ul	828798	Naphthalene-d8	10.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Formamide, N,N-dibutyl-	157	C9H19NO	000761-65-9	87
2			Diisopropylethylamine	129	C8H19N	007087-68-5	64
3			Diisopropylethylamine	129	C8H19N	007087-68-5	64
4			1-Methylsilacyclopentane	100	C5H12Si	1000433-92-4	50
5			Diisopropylethylamine	129	C8H19N	007087-68-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
Operator : CG/JU
Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

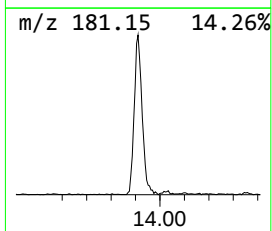
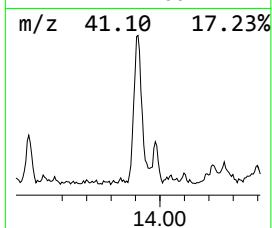
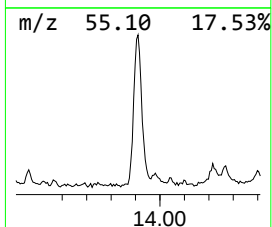
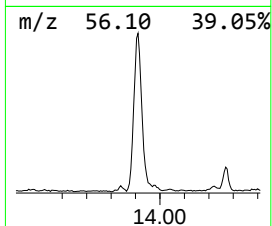
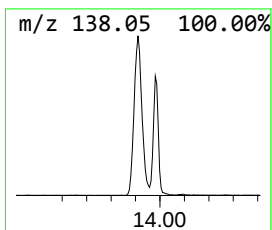
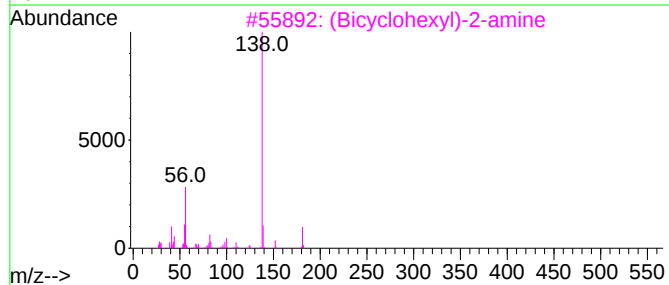
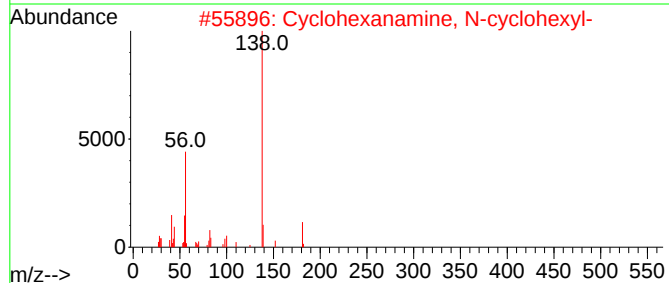
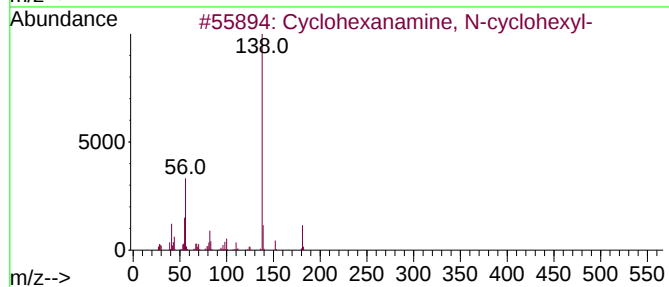
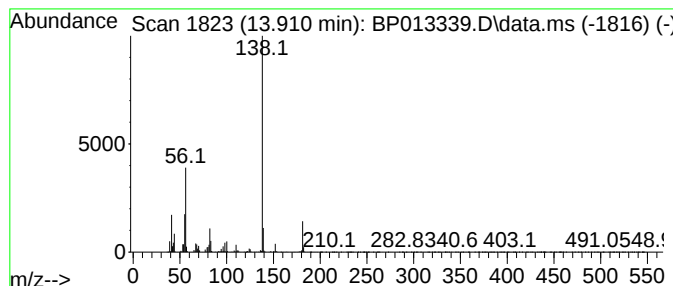
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclohexanamine, N-cyclohexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.910	4.07 ng/ul	1254360	Acenaphthene-d10	14.575	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	96
2	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	96
3	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	96
4	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
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Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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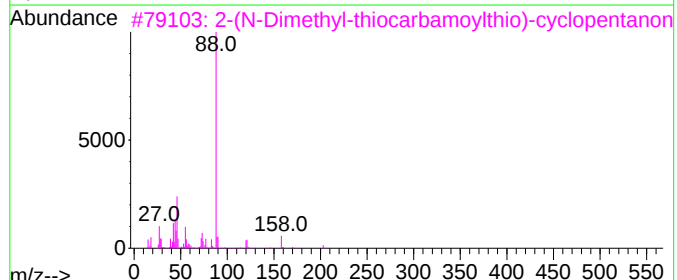
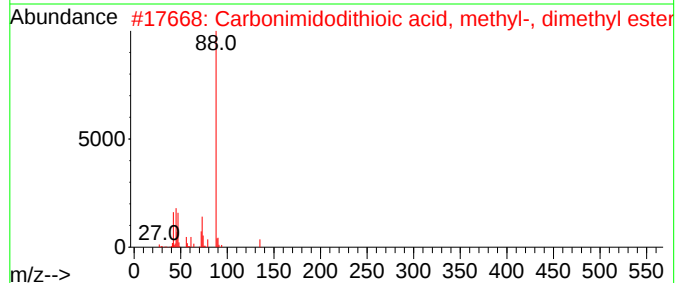
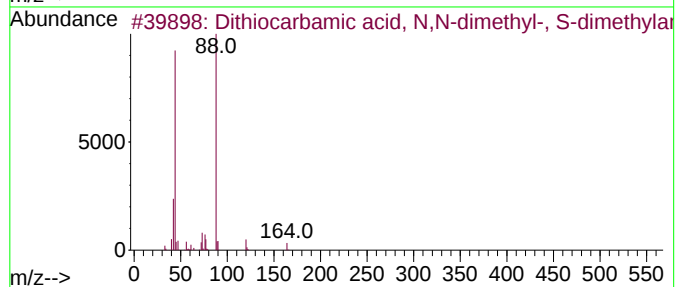
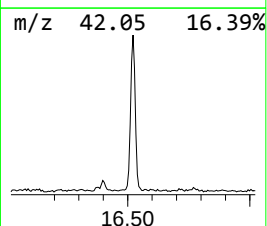
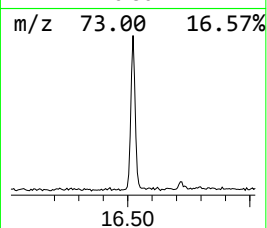
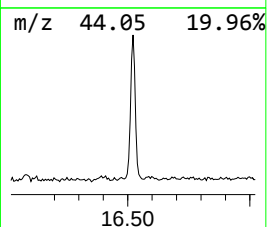
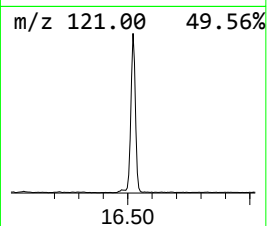
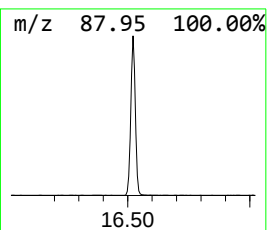
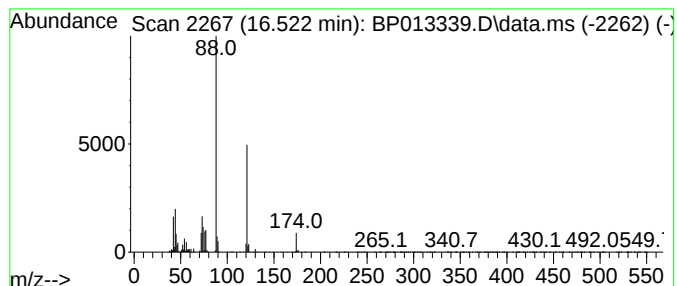
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.522	3.87 ng/ul	1379920	Phenanthrene-d10	17.334

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dithiocarbamic acid, N,N-dimethy...	164	C5H12N2S2	002801-22-1	42
2			Carbonimidodithioic acid, methyl...	135	C4H9NS2	018805-25-9	38
3			2-(N-Dimethyl-thiocarbamoylthio)...	203	C8H13NOS2	023839-30-7	38
4			Allyl dimethyldithiocarbamate	161	C6H11NS2	020821-66-3	38
5			Ethyl 1,3-thiazolidine-4-carboxy...	161	C6H11N02S	051977-21-0	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
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Acq On : 27 Dec 2022 15:26
Operator : CG/JU
Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

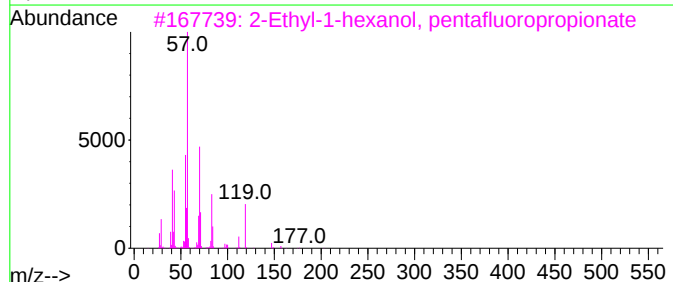
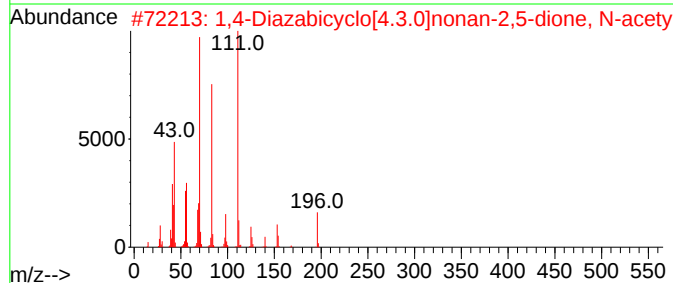
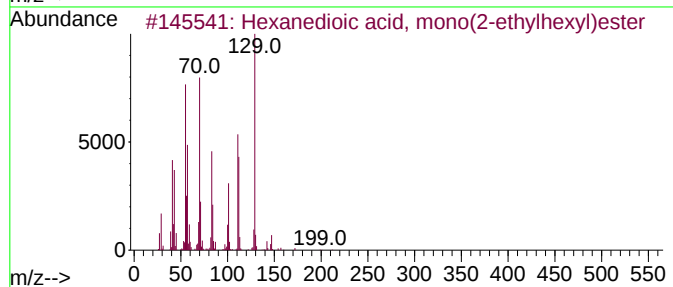
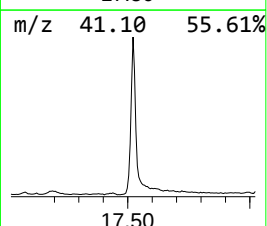
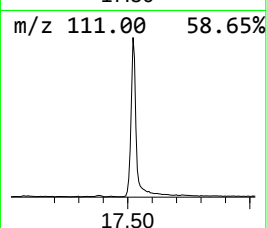
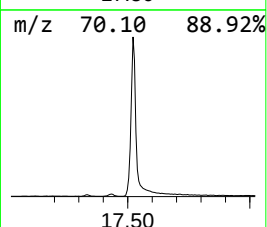
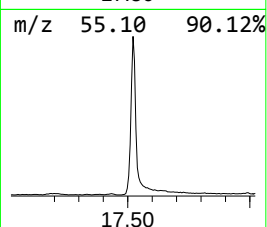
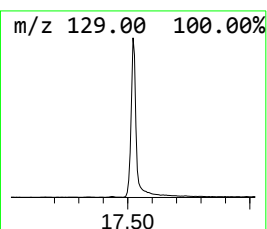
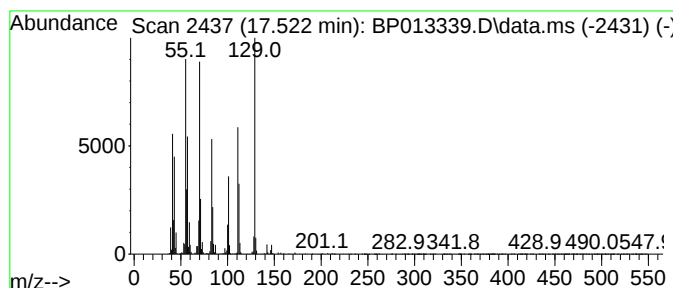
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexanedioic acid, mono(2-et... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.522	15.64 ng/ul	5568950	Phenanthrene-d10		17.334	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexanedioic acid, mono(2-ethylhe...		258	C14H26O4	004337-65-9	83
2	1,4-Diazabicyclo[4.3.0]nonan-2,5...		196	C9H12N2O3	1000512-90-4	38
3	2-Ethyl-1-hexanol, pentafluoropr...		276	C11H17F5O2	959012-82-9	38
4	iso-Amyl tiglate		170	C10H18O2	066917-62-2	35
5	2-Ethyl-1-hexanol, pentafluoropr...		276	C11H17F5O2	959012-82-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
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Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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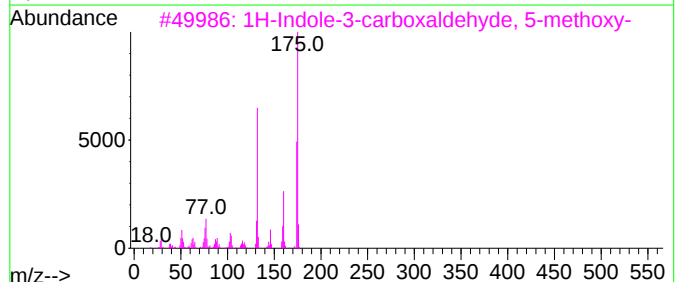
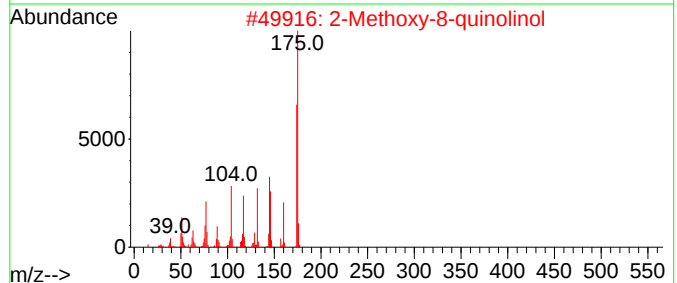
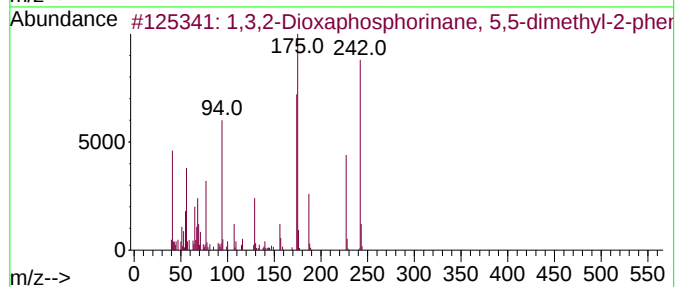
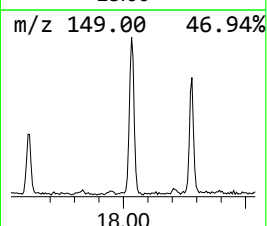
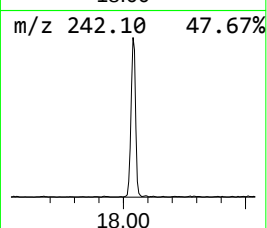
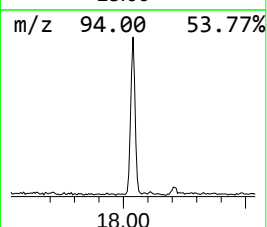
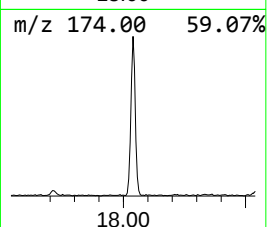
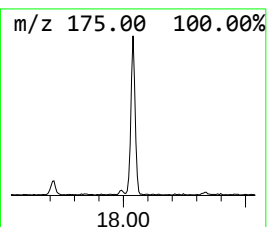
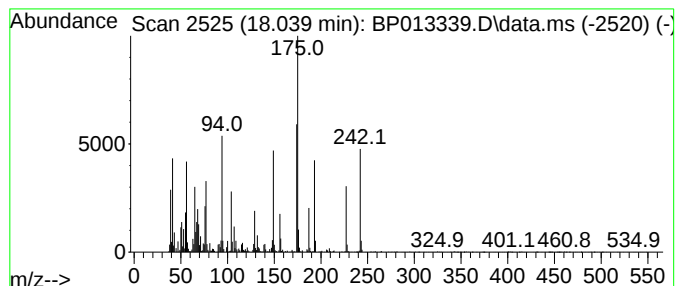
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,3,2-Dioxaphosphorinane, 5... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	2.07 ng/ul	737744	Phenanthrene-d10	17.334

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,2-Dioxaphosphorinane, 5,5-di...	242	C11H15O4P	000884-89-9	99
2			2-Methoxy-8-quinolinol	175	C10H9NO2	074668-72-7	22
3			1H-Indole-3-carboxaldehyde, 5-me...	175	C10H9NO2	010601-19-1	22
4			Phosphoric acid, dibutyl phenyl ...	286	C14H23O4P	002528-36-1	16
5			2-Methoxy-8-quinolinol, acetate	217	C12H11NO3	1000852-74-3	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
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Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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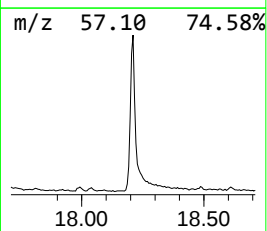
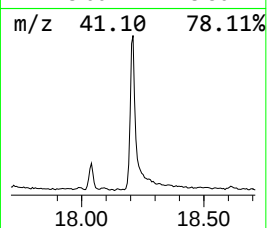
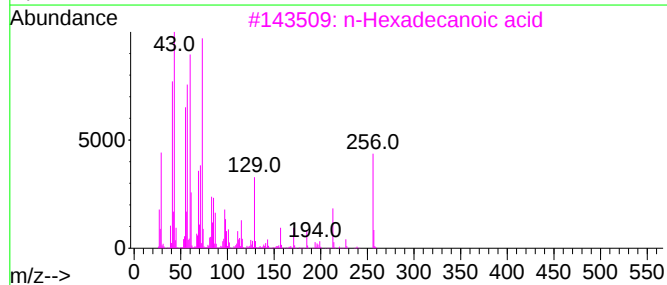
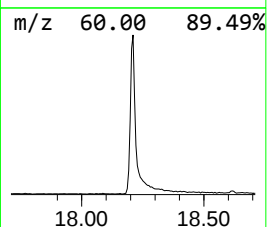
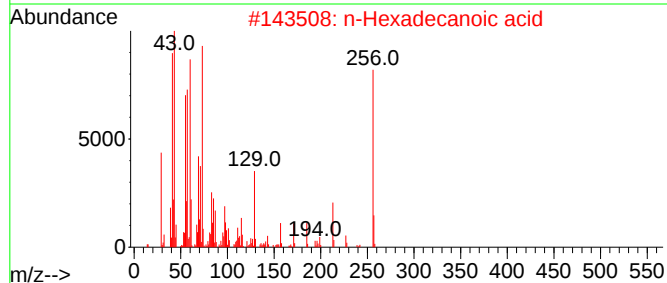
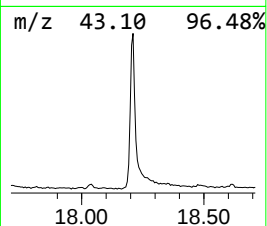
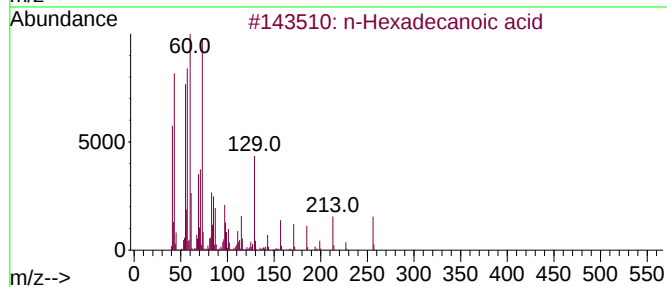
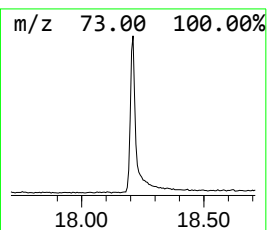
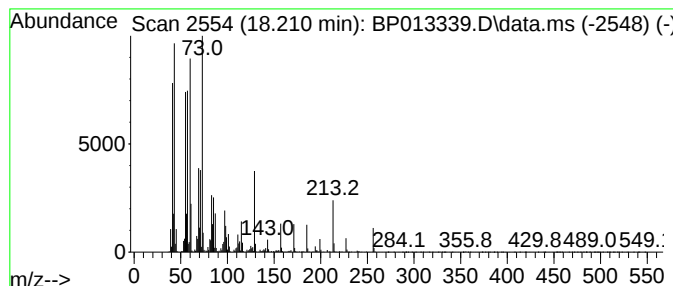
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	8.69 ng/ul	3093390	Phenanthrene-d10	17.334

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
Operator : CG/JU
Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EXGJ1

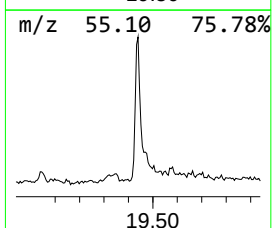
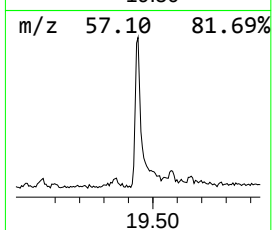
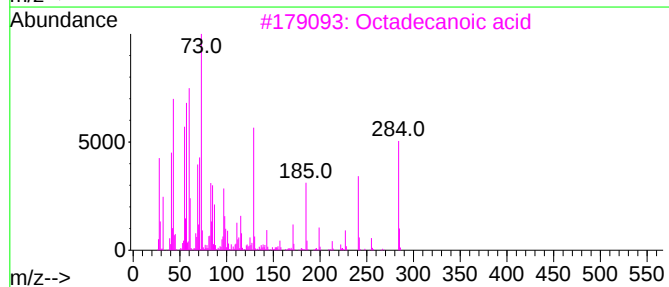
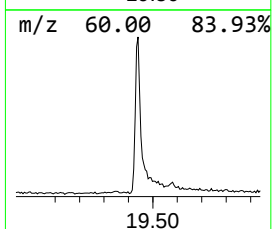
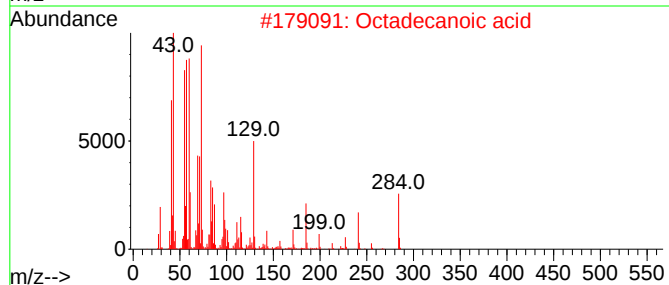
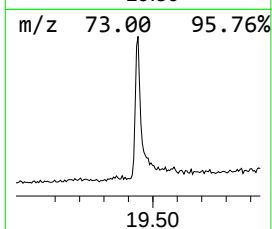
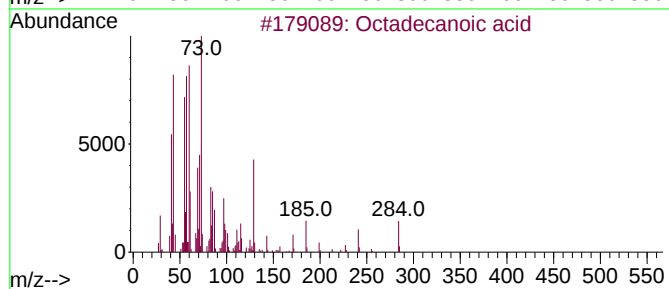
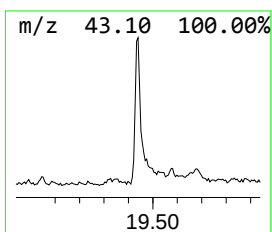
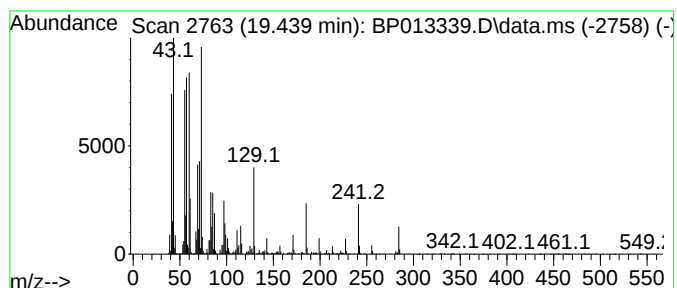
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.439	3.94 ng/ul	1377220	Chrysene-d12	21.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
Operator : CG/JU
Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

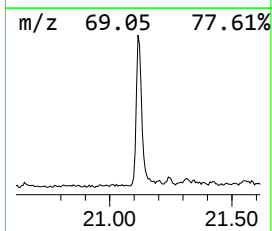
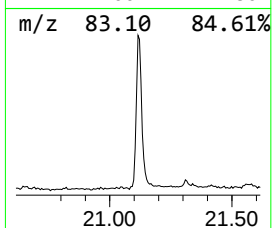
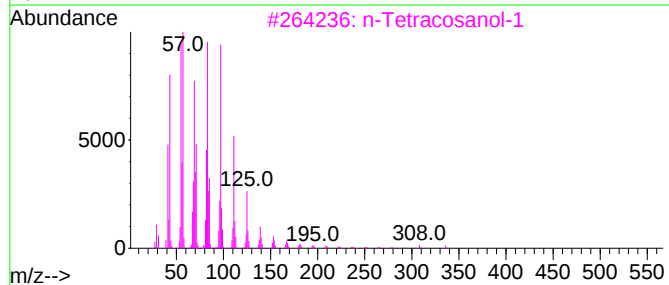
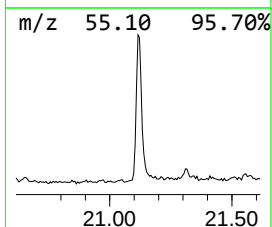
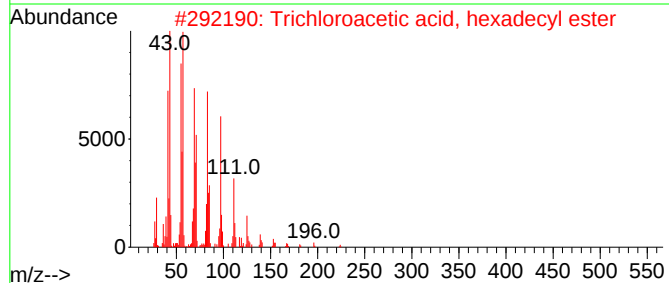
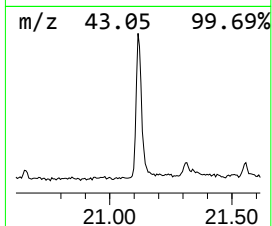
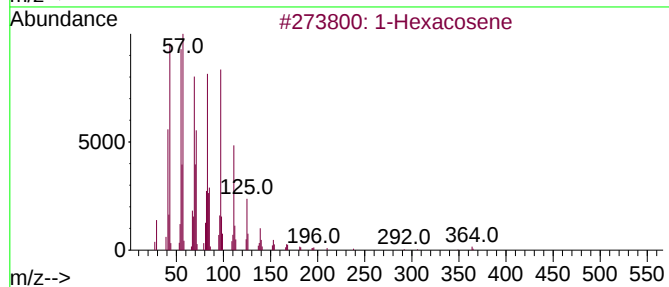
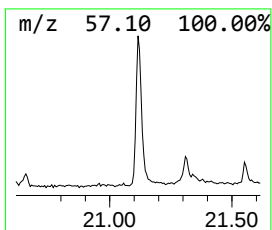
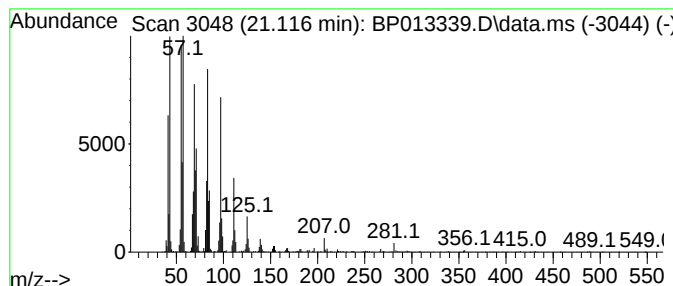
Instrument :
BNA_P
ClientSampleId :
EXGJ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.116	4.09 ng/ul	1428670	Chrysene-d12	21.422	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	94
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122722\
Data File : BP013339.D
Acq On : 27 Dec 2022 15:26
Operator : CG/JU
Sample : N6133-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXGJ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
2-Pentanone, 4-...	5.022	2.4	ng/ul	361868	1	7.934	3000650	20.0
1-Hexanol, 2-et...	7.993	11.6	ng/ul	1740210	1	7.934	3000650	20.0
Hexanoic acid, ...	9.246	9.9	ng/ul	1481940	1	7.934	3000650	20.0
Formamide, N,N-...	12.234	3.7	ng/ul	828798	2	10.746	4466410	20.0
Cyclohexanamine...	13.910	4.1	ng/ul	1254360	3	14.575	6164980	20.0
unknown-01	16.522	3.9	ng/ul	1379920	4	17.334	7122570	20.0
Hexanedioic aci...	17.522	15.6	ng/ul	5568950	4	17.334	7122570	20.0
1,3,2-Dioxaphos...	18.039	2.1	ng/ul	737744	4	17.334	7122570	20.0
n-Hexadecanoic ...	18.210	8.7	ng/ul	3093390	4	17.334	7122570	20.0
Octadecanoic acid	19.439	3.9	ng/ul	1377220	5	21.422	6988920	20.0
(DEL) Alkane: S...	21.116	4.1	ng/ul	1428670	5	21.422	6988920	20.0