

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022047.D
 Acq On : 26 Sep 2024 04:19
 Operator : RC/JU
 Sample : P4130-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0H10

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-BP092424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022047.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.058	9	12	20	rVB	26853	36177	0.82%	0.087%
2	3.222	36	40	49	rVB	24791	30100	0.68%	0.072%
3	3.499	82	87	93	rVB	30368	41748	0.95%	0.100%
4	3.640	107	111	121	rBV	130809	209239	4.75%	0.502%
5	3.722	121	125	128	rVV	41579	59210	1.34%	0.142%
6	3.758	128	131	136	rVV2	14903	21814	0.49%	0.052%
7	3.816	136	141	147	rVV	60388	91270	2.07%	0.219%
8	4.075	182	185	190	rVB2	8092	11615	0.26%	0.028%
9	4.199	203	206	211	rBV3	7484	14005	0.32%	0.034%
10	4.258	212	216	223	rVV4	8836	16395	0.37%	0.039%
11	4.458	245	250	255	rVB	23194	32877	0.75%	0.079%
12	4.681	283	288	291	rBV3	7449	11703	0.27%	0.028%
13	4.716	291	294	299	rVV2	8030	12181	0.28%	0.029%
14	4.863	314	319	323	rVB5	7018	11573	0.26%	0.028%
15	4.916	323	328	333	rBV2	17961	28086	0.64%	0.067%
16	4.969	333	337	341	rVB3	10285	15239	0.35%	0.037%
17	5.022	341	346	348	rBV4	9037	13715	0.31%	0.033%
18	5.069	348	354	368	rVB	2008706	2994023	67.93%	7.188%
19	5.199	369	376	380	rBV6	8820	17012	0.39%	0.041%
20	5.552	429	436	440	rVB9	5185	10467	0.24%	0.025%
21	5.616	440	447	452	rBV10	4531	10719	0.24%	0.026%
22	5.734	463	467	478	rVB10	4570	12749	0.29%	0.031%
23	5.940	497	502	507	rBV8	5223	12257	0.28%	0.029%
24	6.010	512	514	518	rVV4	5605	9993	0.23%	0.024%
25	6.058	518	522	527	rVV7	7023	12636	0.29%	0.030%
26	6.269	549	558	565	rVB10	7482	17768	0.40%	0.043%
27	6.481	589	594	598	rBV5	8267	14078	0.32%	0.034%
28	6.522	598	601	612	rVB8	5746	14253	0.32%	0.034%
29	6.887	654	663	671	rBV	135645	226053	5.13%	0.543%
30	7.046	683	690	697	rVV	291515	451799	10.25%	1.085%
31	7.116	697	702	708	rVB2	14926	24702	0.56%	0.059%
32	7.246	716	724	731	rBV	381065	600450	13.62%	1.441%
33	7.363	740	744	753	rVB6	9872	18132	0.41%	0.044%
34	7.599	776	784	790	rVB	177913	285248	6.47%	0.685%
35	7.705	795	802	804	rVV	339856	523834	11.89%	1.258%
36	7.740	804	808	827	rVB	1360770	2263499	51.36%	5.434%

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Integration Parameters: LSCINT.P

Integrator: RTE

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37	7.987	833	850	854	rBV	140544	348582	7.91%	0.837%
38	8.052	854	861	870	rVV	1735436	2748066	62.35%	6.597%
39	8.340	906	910	913	rBV4	6845	10683	0.24%	0.026%
40	8.404	913	921	932	rVB	349847	572493	12.99%	1.374%
41	8.534	938	943	952	rVB	6137	12711	0.29%	0.031%
42	8.616	952	957	961	rBV6	7316	14031	0.32%	0.034%
43	8.669	961	966	971	rVV2	21072	39972	0.91%	0.096%
44	8.734	971	977	985	rVV	69589	122881	2.79%	0.295%
45	8.840	985	995	1008	rVV	347182	647688	14.70%	1.555%
46	8.952	1010	1014	1022	rVB4	8459	18568	0.42%	0.045%
47	9.057	1022	1032	1038	rVB4	16966	39312	0.89%	0.094%
48	9.140	1038	1046	1050	rBV	9603	17070	0.39%	0.041%
49	9.269	1061	1068	1074	rBV7	7560	17997	0.41%	0.043%
50	9.481	1095	1104	1109	rBV7	10462	20842	0.47%	0.050%
51	9.557	1109	1117	1125	rBV	325003	539638	12.24%	1.296%
52	9.757	1142	1151	1160	rVB2	18027	36842	0.84%	0.088%
53	10.093	1200	1208	1216	rBV	542505	865838	19.65%	2.079%
54	10.328	1240	1248	1264	rBV	490748	886526	20.12%	2.128%
55	10.469	1264	1272	1278	rBV	420074	719223	16.32%	1.727%
56	10.522	1278	1281	1286	rVV3	6296	10256	0.23%	0.025%
57	10.598	1286	1294	1307	rVV	375595	671300	15.23%	1.612%
58	10.769	1319	1323	1329	rVB3	6634	12896	0.29%	0.031%
59	10.846	1329	1336	1342	rBV	125155	215478	4.89%	0.517%
60	11.063	1366	1373	1378	rVB7	7886	16057	0.36%	0.039%
61	11.287	1404	1411	1419	rVB7	6173	15404	0.35%	0.037%
62	11.787	1489	1496	1501	rVB2	12945	20792	0.47%	0.050%
63	11.928	1517	1520	1528	rVB10	4247	9856	0.22%	0.024%
64	12.045	1535	1540	1544	rBV3	6565	11828	0.27%	0.028%
65	12.322	1583	1587	1592	rVB6	8300	11737	0.27%	0.028%
66	12.422	1598	1604	1608	rBV5	12362	20248	0.46%	0.049%
67	12.493	1612	1616	1620	rVB3	12044	19760	0.45%	0.047%
68	13.098	1714	1719	1725	rBV	35728	59809	1.36%	0.144%
69	13.298	1749	1753	1759	rVB8	7635	14513	0.33%	0.035%
70	13.716	1814	1824	1830	rBV	997533	1447026	32.83%	3.474%
71	13.922	1856	1859	1866	rVB9	7647	13209	0.30%	0.032%
72	13.998	1866	1872	1880	rBV	1037642	1455359	33.02%	3.494%
73	14.310	1917	1925	1931	rBV	721458	1106204	25.10%	2.656%
74	14.363	1931	1934	1937	rVV4	13527	21346	0.48%	0.051%
75	14.398	1937	1940	1944	rVB6	6055	10147	0.23%	0.024%
76	14.522	1955	1961	1973	rBV	133919	251351	5.70%	0.603%

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Integrator: RTE

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Sampling : 1

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Stop Thrs : 0

Peak Location: TOP

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77	14.610	1973	1976	1983	rVB8	9966	18204	0.41%	0.044%
78	14.939	2026	2032	2035	rBV8	6141	13188	0.30%	0.032%
79	15.039	2044	2049	2052	rBV7	5545	9720	0.22%	0.023%
80	15.139	2061	2066	2072	rBV3	15153	27148	0.62%	0.065%
81	15.316	2089	2096	2104	rBV	1456186	2070105	46.97%	4.970%
82	15.439	2111	2117	2126	rBV	539977	776044	17.61%	1.863%
83	15.569	2133	2139	2147	rVB3	37534	68857	1.56%	0.165%
84	15.692	2153	2160	2168	rBV2	66236	117819	2.67%	0.283%
85	15.910	2195	2197	2203	rVB5	7893	12286	0.28%	0.029%
86	16.757	2337	2341	2350	rVB5	6629	13933	0.32%	0.033%
87	17.110	2394	2401	2408	rBV	969678	1527183	34.65%	3.666%
88	17.204	2411	2417	2430	rVB2	1788888	2763094	62.69%	6.633%
89	17.928	2535	2540	2550	rVB4	17595	45288	1.03%	0.109%
90	18.045	2555	2560	2564	rBV	95242	135770	3.08%	0.326%
91	18.122	2570	2573	2580	rVB	37111	44835	1.02%	0.108%
92	18.922	2706	2709	2712	rBV5	8304	11405	0.26%	0.027%
93	19.133	2741	2745	2749	rVB2	29378	39365	0.89%	0.095%
94	19.204	2753	2757	2762	rBV2	27661	49770	1.13%	0.119%
95	19.380	2783	2787	2793	rBV4	41281	55919	1.27%	0.134%
96	19.533	2811	2813	2816	rBV3	9764	10442	0.24%	0.025%
97	19.586	2816	2822	2829	rBV2	2670460	3892318	88.32%	9.344%
98	21.539	3147	3154	3165	rVB2	1174593	2063594	46.82%	4.954%
99	24.586	3662	3672	3686	rVB	1505572	4407227	100.00%	10.580%
100	24.827	3703	3713	3724	rVB	881404	2211124	50.17%	5.308%

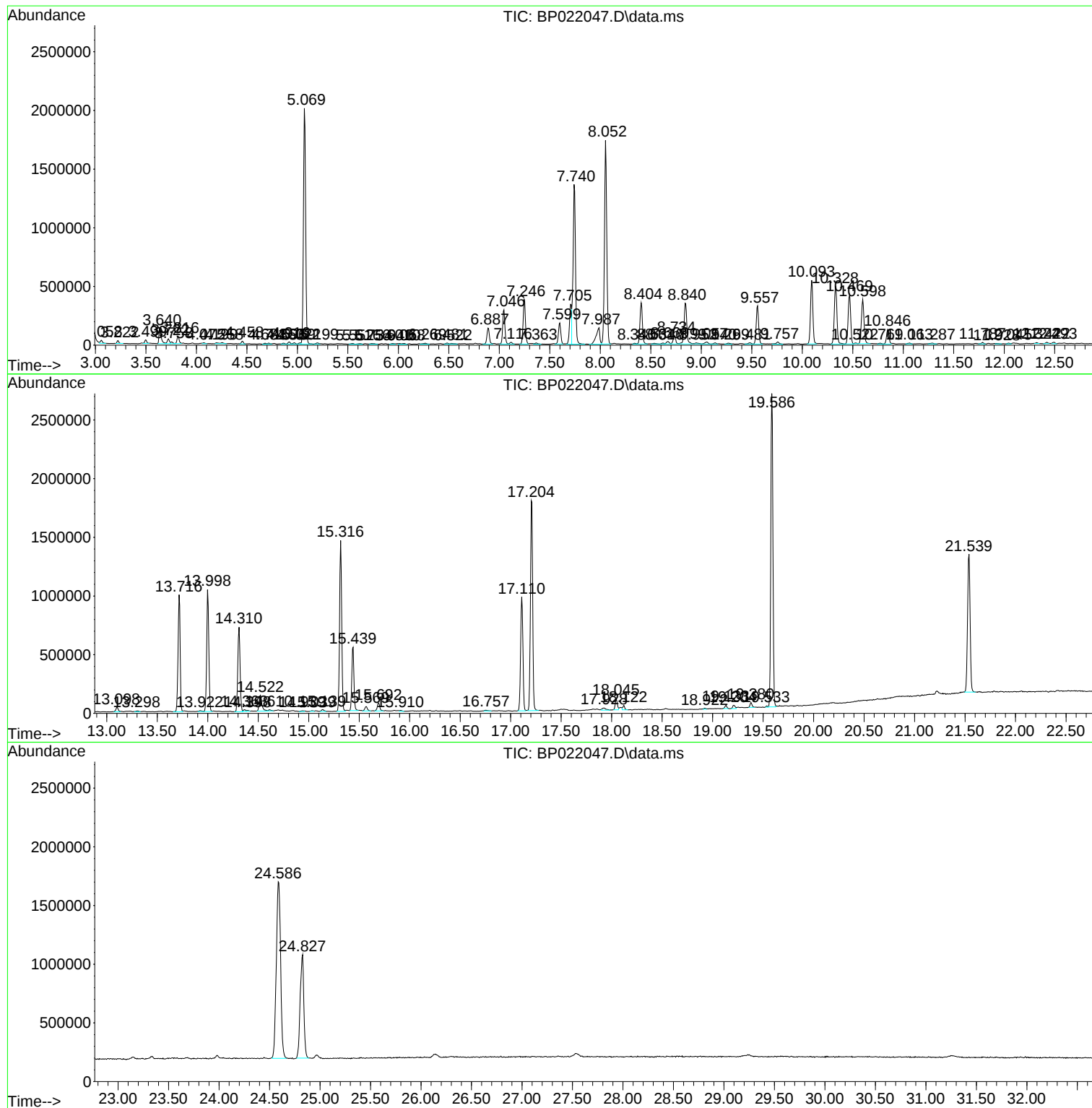
Sum of corrected areas: 41654796

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Instrument :
BNA_P
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H0H10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_P
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H0H10

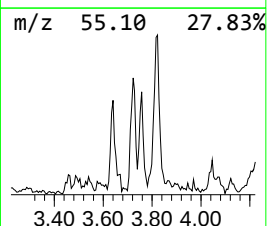
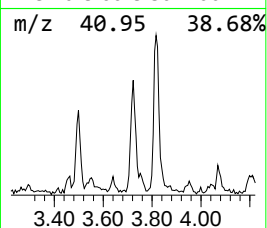
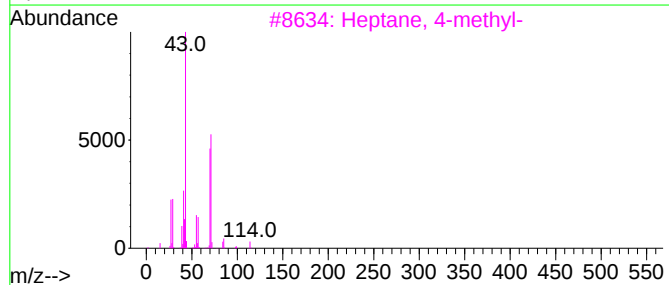
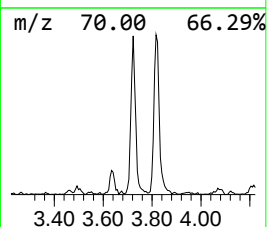
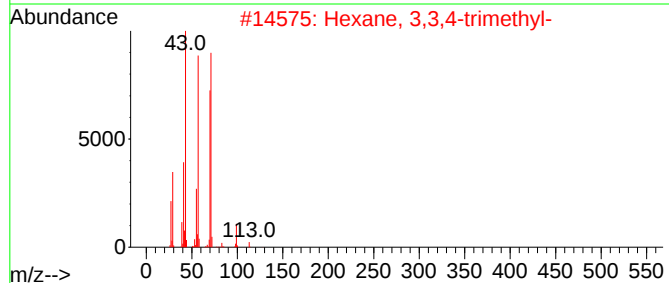
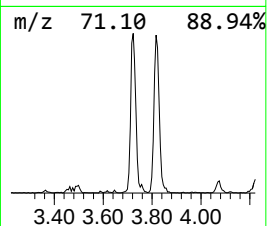
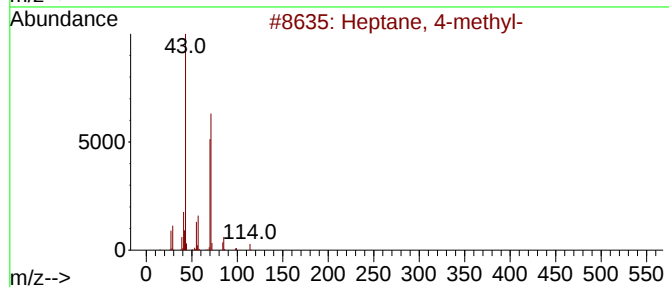
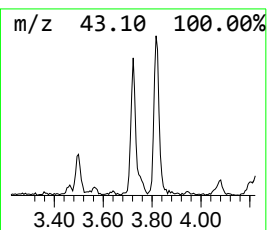
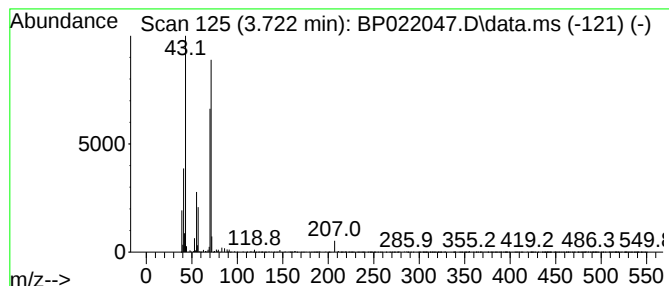
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.722	2.26 ng/ul	59210	1,4-Dichlorobenzene-d4	7.705

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-methyl-	114	C8H18	000589-53-7	78
2		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	78
4		Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	64



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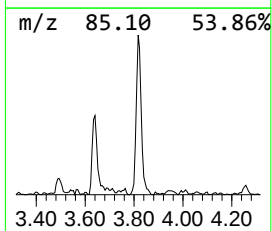
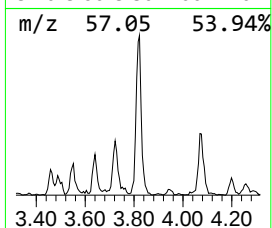
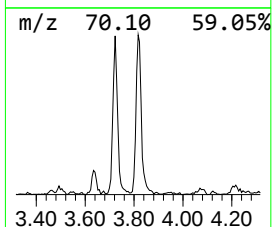
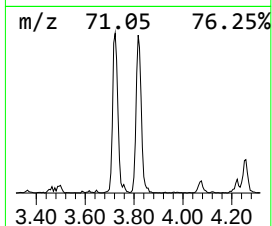
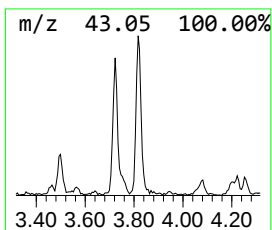
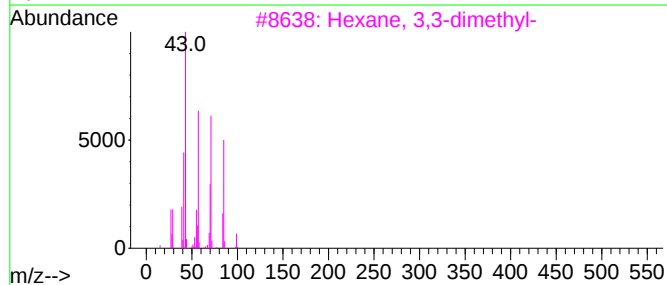
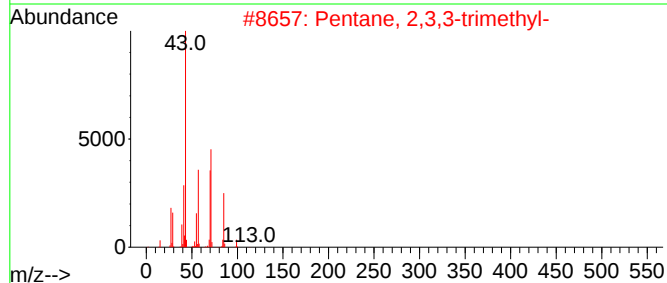
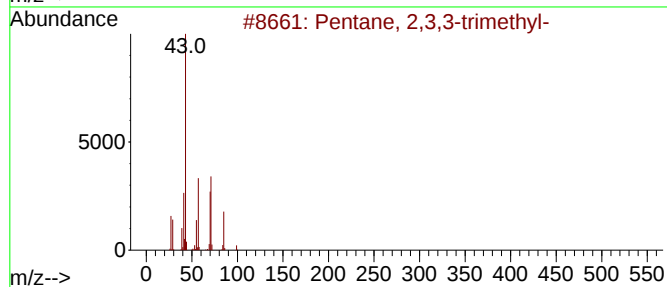
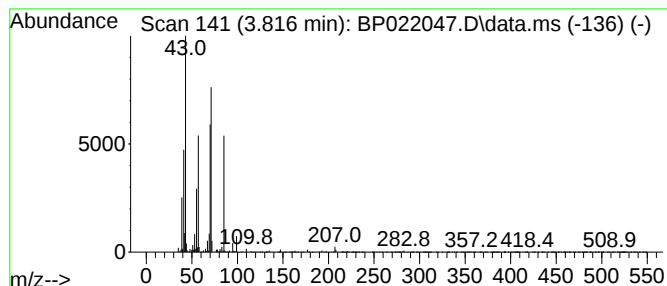
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.816	3.48 ng/ul	91270	1,4-Dichlorobenzene-d4	7.705

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	78
2		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	78
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5		Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	59



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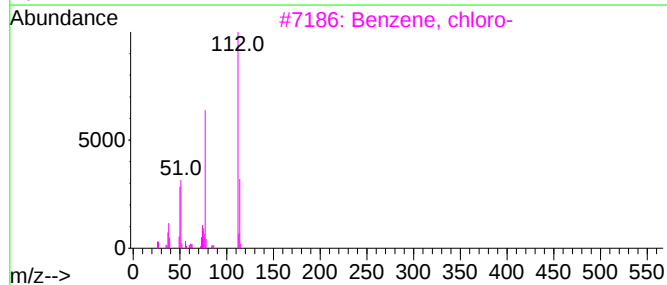
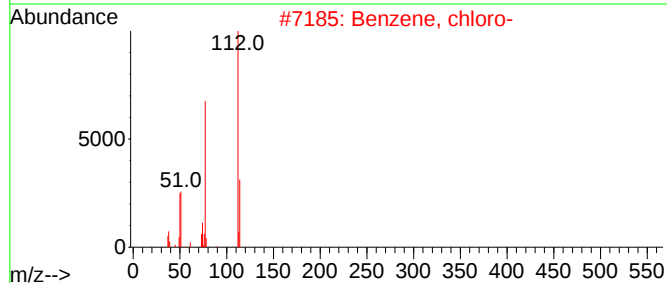
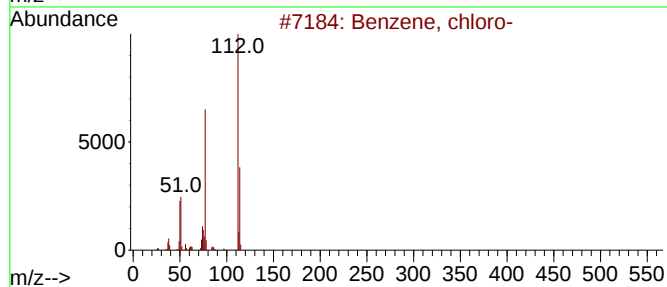
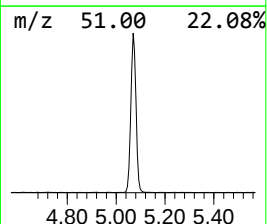
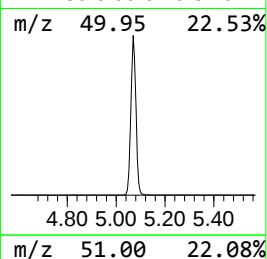
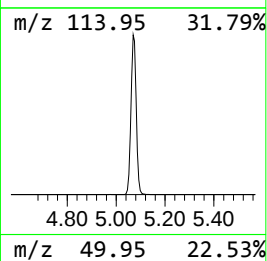
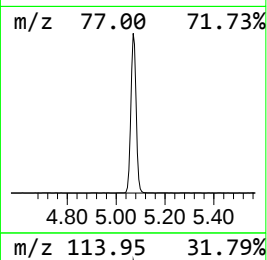
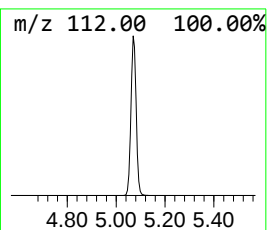
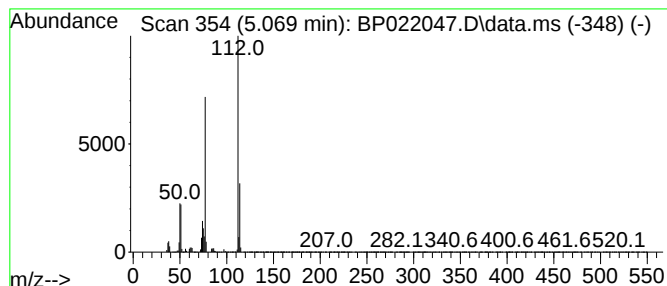
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.069	114.31 ng/ul	2994020	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	97
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	95
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
Operator : RC/JU
Sample : P4130-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0H10

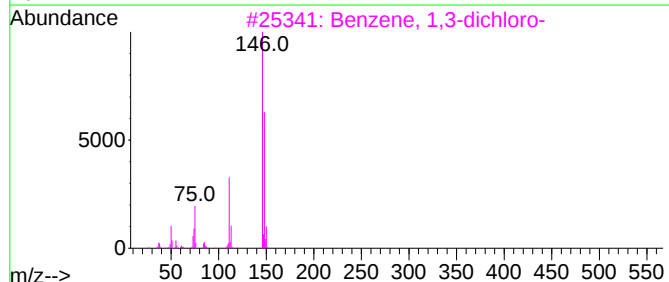
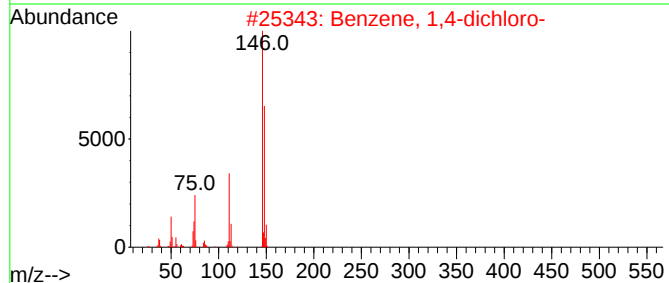
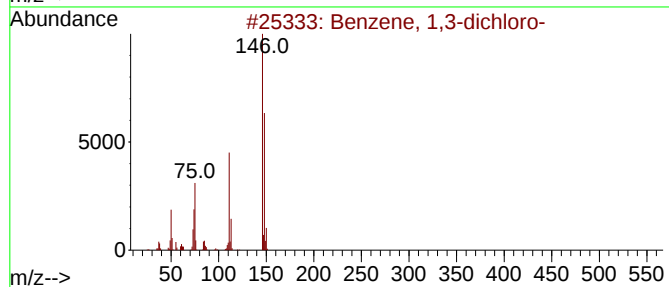
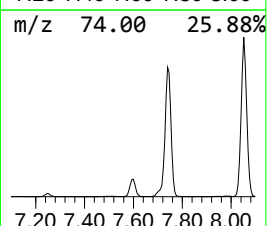
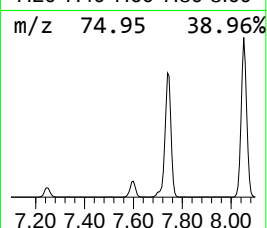
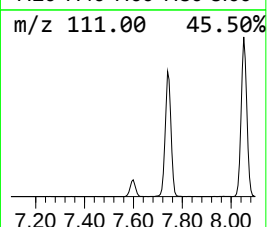
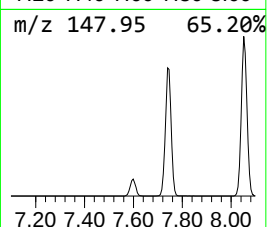
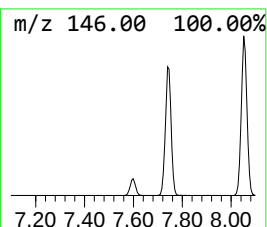
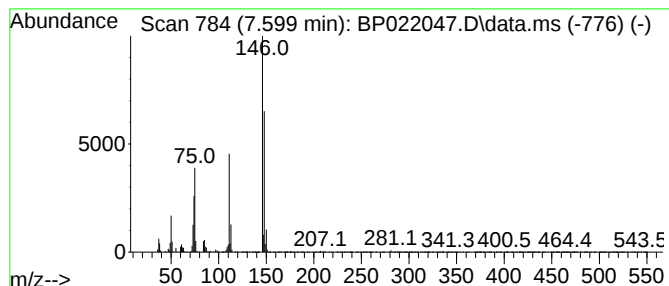
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.599	10.89 ng/ul	285248	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
Operator : RC/JU
Sample : P4130-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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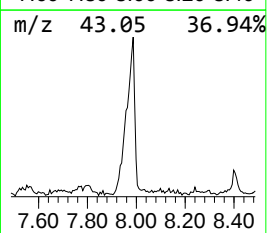
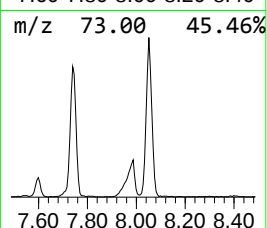
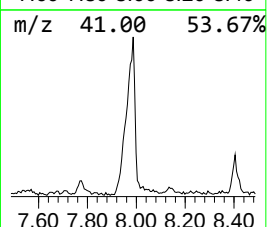
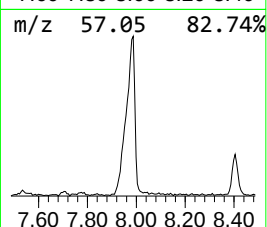
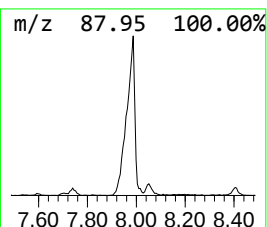
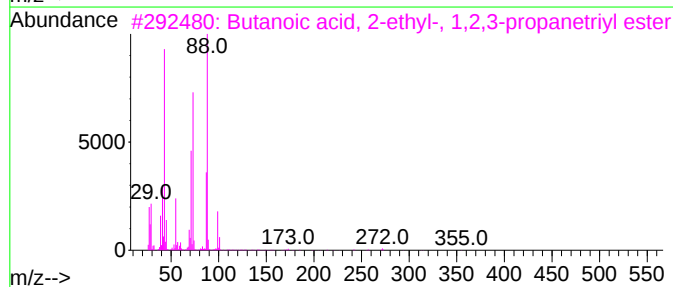
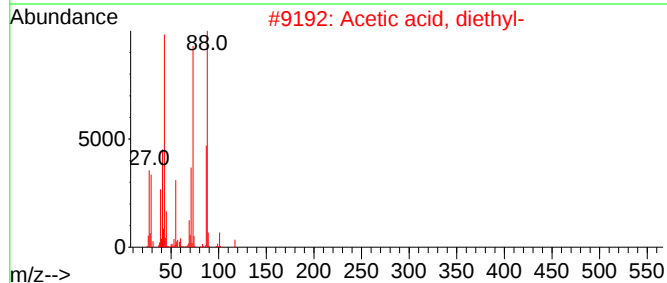
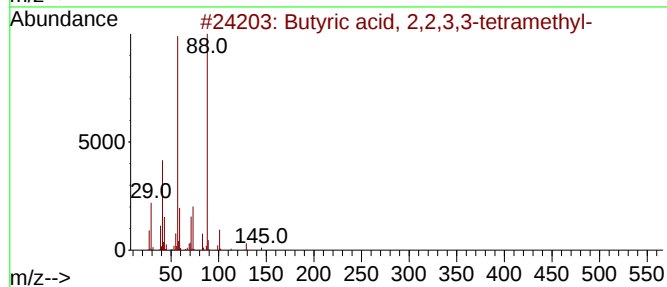
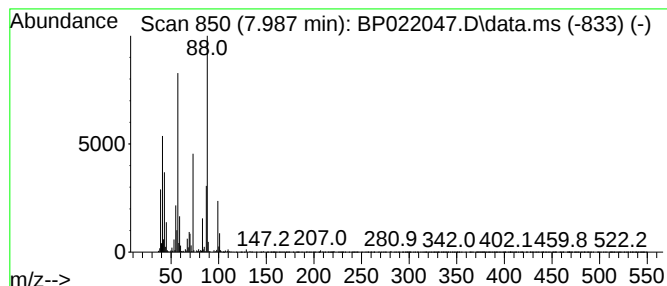
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Butyric acid, 2,2,3,3-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.987	13.31 ng/ul	348582	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butyric acid, 2,2,3,3-tetramethyl-	144	C8H16O2	030407-41-1	64
2			Acetic acid, diethyl-	116	C6H12O2	000088-09-5	53
3			Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	43
4			Propanoic acid, 2-[dimethylamino...	281	C8H14BF6NO2	1000161-96-0	38
5			Heptanoic acid, 2-methyl-, methy...	158	C9H18O2	051209-78-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
Operator : RC/JU
Sample : P4130-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
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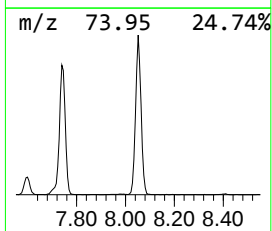
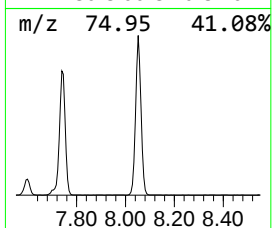
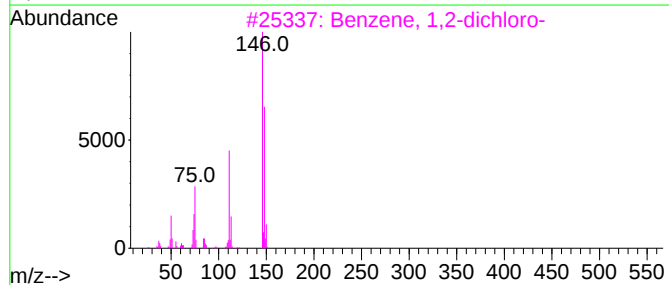
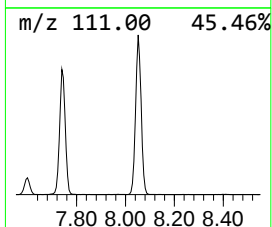
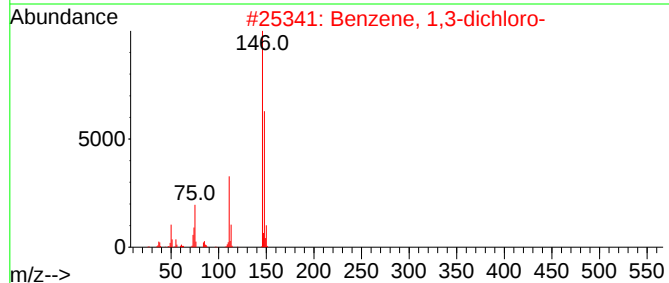
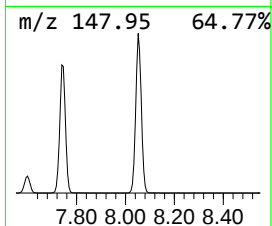
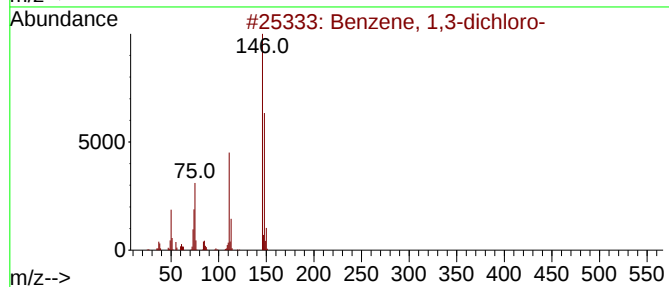
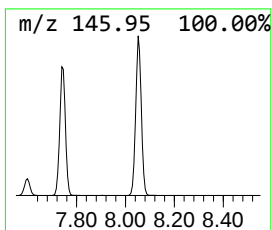
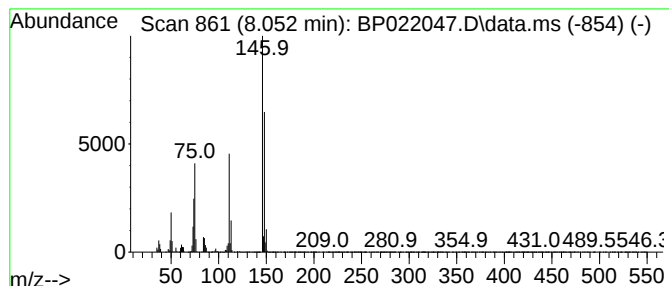
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.052	104.92 ng/ul	2748070	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
Operator : RC/JU
Sample : P4130-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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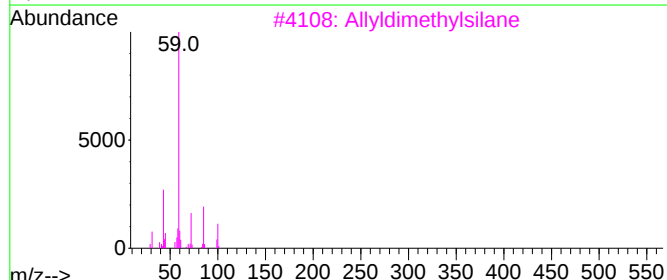
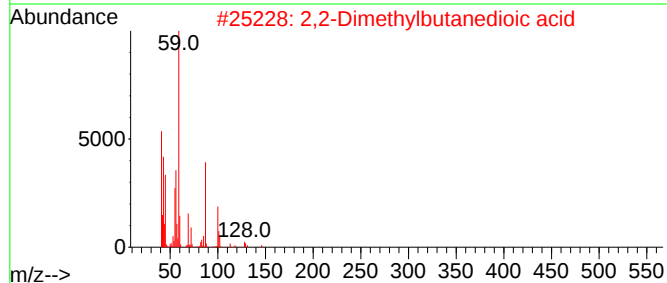
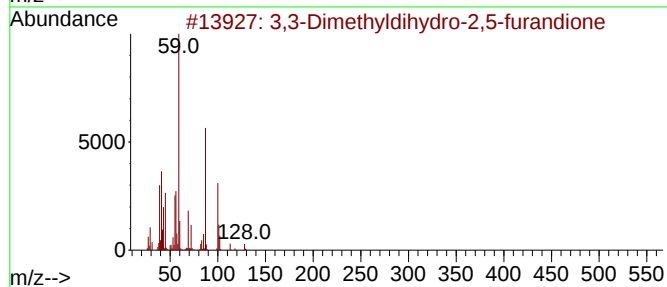
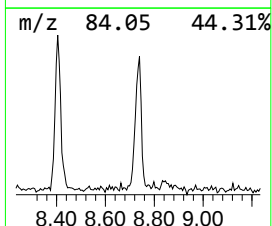
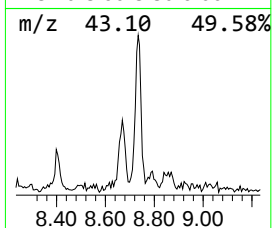
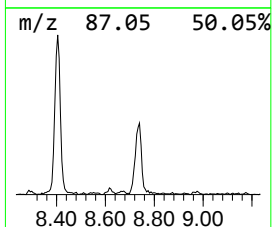
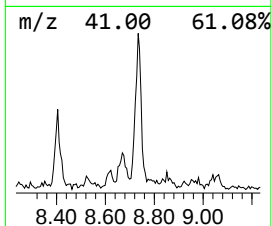
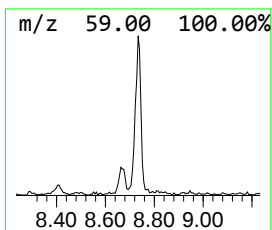
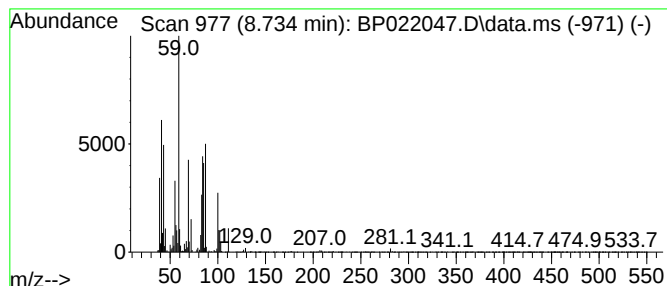
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.734	4.69 ng/ul	122881	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,3-Dimethyldihydro-2,5-furandione	128	C6H8O3	017347-61-4	38
2			2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	37
3			Allyldimethylsilane	100	C5H12Si	003937-30-2	32
4			3-Pentadecanol	228	C15H32O	053346-71-7	27
5			N-(1-Carbomethoxyethyloxycarbonyl...)	325	C9H9F6NO5	1000283-30-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
Operator : RC/JU
Sample : P4130-06
Misc :
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Instrument :
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ClientSampleId :
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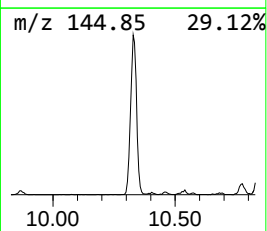
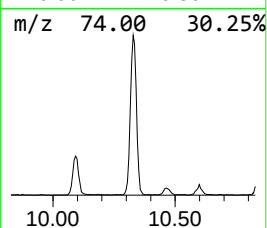
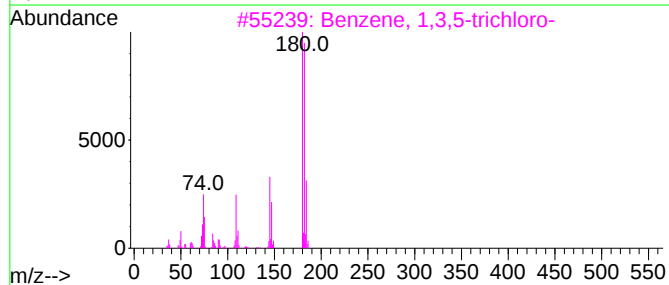
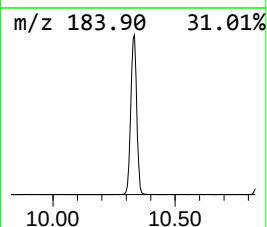
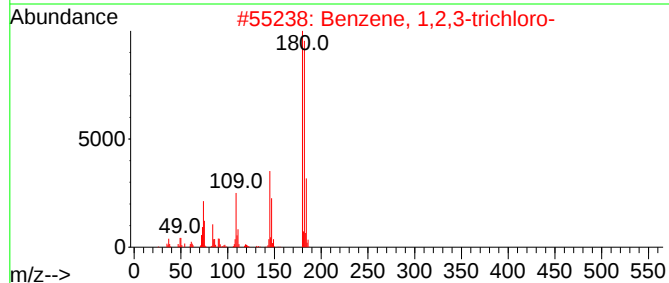
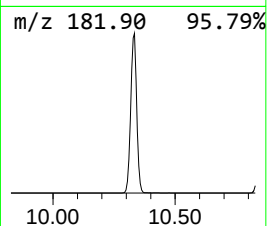
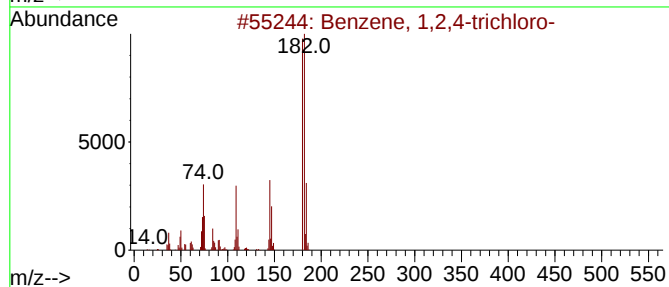
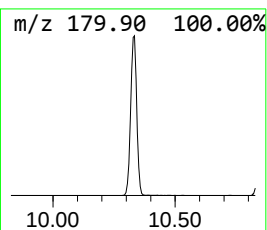
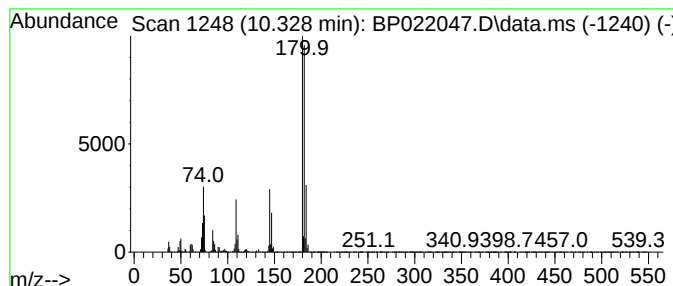
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,4-trichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.328	24.65 ng/ul	886526	Naphthalene-d8	10.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
4			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
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Sample : P4130-06
Misc :
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Instrument :
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ClientSampleId :
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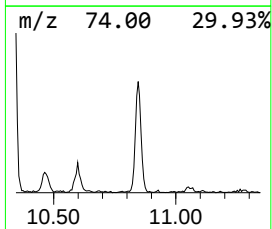
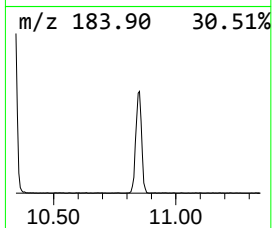
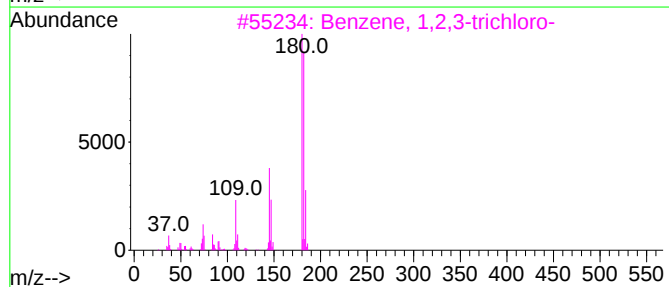
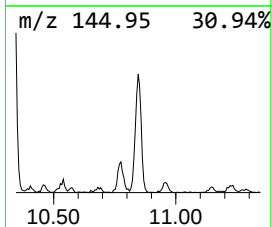
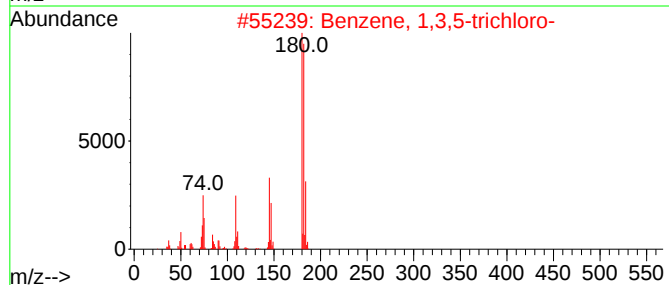
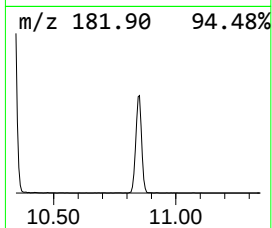
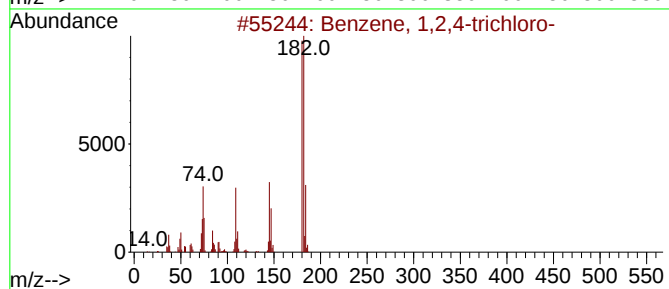
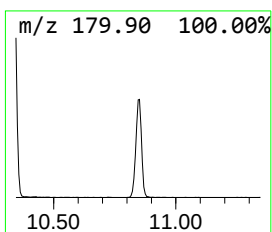
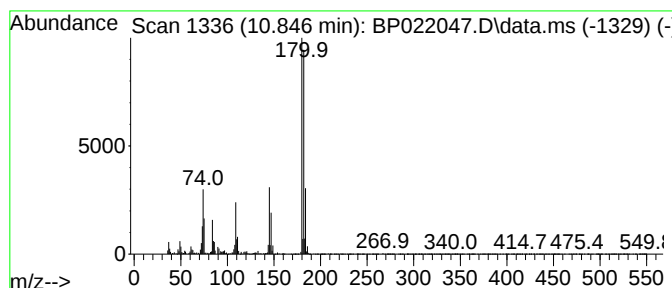
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,3,5-trichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.845	5.99 ng/ul	215478	Naphthalene-d8	10.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
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ClientSampleId :
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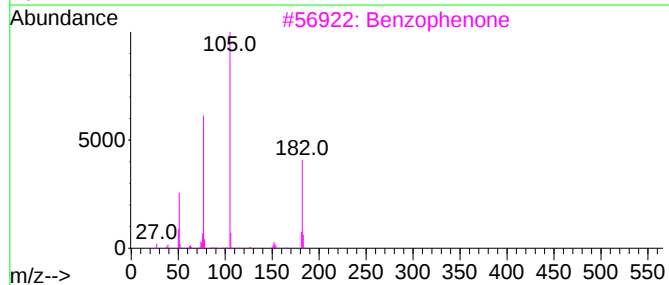
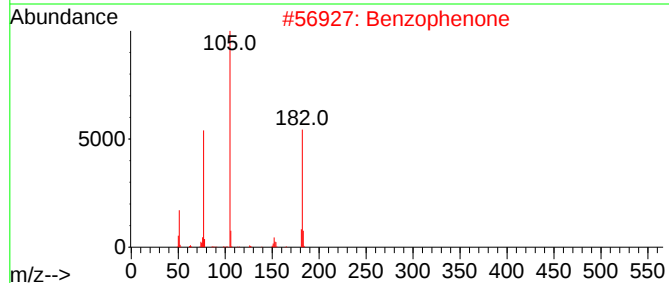
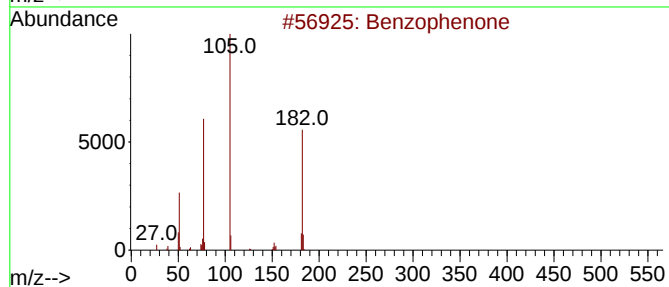
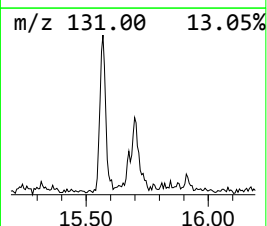
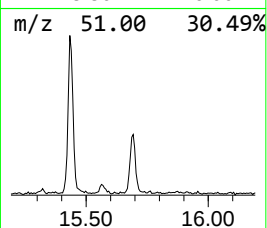
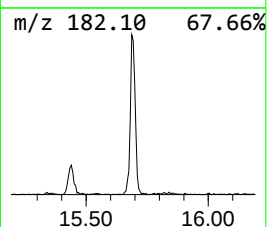
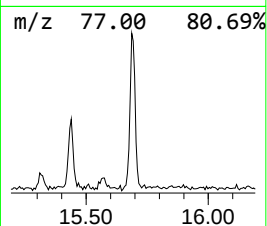
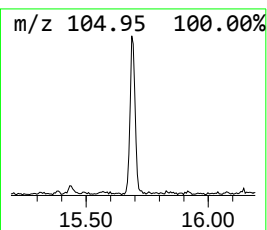
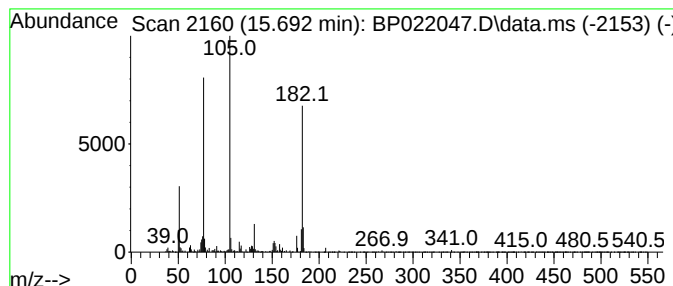
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.692	2.13 ng/ul	117819	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	93
2			Benzophenone	182	C13H10O	000119-61-9	91
3			Benzophenone	182	C13H10O	000119-61-9	87
4			Benzophenone	182	C13H10O	000119-61-9	86
5			Benzophenone	182	C13H10O	000119-61-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
Data File : BP022047.D
Acq On : 26 Sep 2024 04:19
Operator : RC/JU
Sample : P4130-06
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0H10

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.722	2.3	ng/ul	59210	1	7.705	523834	20.0
(DEL) Alkane: S...	3.816	3.5	ng/ul	91270	1	7.705	523834	20.0
Benzene, chloro-	5.069	114.3	ng/ul	2994020	1	7.705	523834	20.0
Benzene, 1,3-di...	7.599	10.9	ng/ul	285248	1	7.705	523834	20.0
Butyric acid, 2...	7.987	13.3	ng/ul	348582	1	7.705	523834	20.0
Benzene, 1,2-di...	8.052	104.9	ng/ul	2748070	1	7.705	523834	20.0
unknown-01	8.734	4.7	ng/ul	122881	1	7.705	523834	20.0
Benzene, 1,2,4-...	10.328	24.6	ng/ul	886526	2	10.469	719223	20.0
Benzene, 1,3,5-...	10.845	6.0	ng/ul	215478	2	10.469	719223	20.0
Benzophenone	15.692	2.1	ng/ul	117819	3	14.310	1106200	20.0