

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
 Data File : BP019398.D
 Acq On : 19 Jan 2024 10:29
 Operator : MA/JU
 Sample : P1160-09
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 XA318

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-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019398.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	18	22	28	rVB	28377	39586	0.18%	0.010%
2	6.140	514	519	522	rBV	25147	38316	0.17%	0.010%
3	6.993	651	664	679	rVB2	3550304	6539868	29.57%	1.686%
4	7.117	679	685	692	rVV	427947	656316	2.97%	0.169%
5	7.211	695	701	711	rVB	961694	1351914	6.11%	0.349%
6	7.340	718	723	733	rVB	2378660	3850917	17.41%	0.993%
7	7.658	769	777	785	rBV	1408392	2207583	9.98%	0.569%
8	7.769	788	796	797	rBV	516801	651198	2.94%	0.168%
9	7.811	797	803	812	rVB	5373142	8598829	38.88%	2.217%
10	8.122	847	856	866	rVV	3753335	5979181	27.03%	1.542%
11	8.511	914	922	928	rBV	735542	1122805	5.08%	0.289%
12	8.599	928	937	949	rVV	2735784	4644924	21.00%	1.198%
13	8.840	969	978	987	rVB	3618153	5786347	26.16%	1.492%
14	8.940	987	995	998	rBV	558229	920805	4.16%	0.237%
15	8.981	998	1002	1011	rVB	1137020	1778143	8.04%	0.458%
16	9.340	1055	1063	1070	rVB4	26918	60048	0.27%	0.015%
17	9.658	1110	1117	1118	rBV	500933	664218	3.00%	0.171%
18	9.693	1118	1123	1136	rVB	1298755	2287064	10.34%	0.590%
19	9.999	1166	1175	1185	rVB	1825851	2900150	13.11%	0.748%
20	10.228	1203	1214	1228	rBV2	2017810	4459190	20.16%	1.150%
21	10.428	1242	1248	1255	rVB7	13705	31103	0.14%	0.008%
22	10.569	1255	1272	1275	rBV	649660	1201204	5.43%	0.310%
23	10.622	1275	1281	1290	rVV	6787941	11307404	51.12%	2.915%
24	10.722	1290	1298	1311	rVB	427293	803528	3.63%	0.207%
25	10.869	1318	1323	1331	rVB	17981	28189	0.13%	0.007%
26	11.881	1487	1495	1509	rBV	3841648	6079677	27.49%	1.568%
27	12.663	1624	1628	1634	rVB	31411	43982	0.20%	0.011%
28	12.852	1653	1660	1666	rVV	3122024	4749398	21.47%	1.225%
29	12.928	1666	1673	1686	rVV	6321002	9737653	44.02%	2.511%
30	13.293	1728	1735	1745	rVV2	4038737	7392156	33.42%	1.906%
31	13.357	1745	1746	1756	rVV2	41617	55518	0.25%	0.014%
32	13.469	1760	1765	1775	rVB	54957	84478	0.38%	0.022%
33	13.846	1820	1829	1836	rBV	1322471	2001667	9.05%	0.516%
34	14.028	1851	1860	1865	rBV	6269665	10282811	46.49%	2.651%
35	14.151	1868	1881	1886	rVV	9093612	15824102	71.54%	4.080%
36	14.310	1901	1908	1913	rVB3	20581	36572	0.17%	0.009%

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Integrator: RTE

Smoothing : OFF

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Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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Stop Thrs : 0

Peak Location: TOP

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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37	14.422	1917	1927	1932	rBV	966095	1495584	6.76%	0.386%
38	14.487	1932	1938	1944	rVV	7583128	11401301	51.55%	2.940%
39	14.540	1944	1947	1952	rVB	101237	135049	0.61%	0.035%
40	14.693	1963	1973	1977	rBV2	3357279	5854631	26.47%	1.510%
41	14.734	1977	1980	1987	rVB	146905	236353	1.07%	0.061%
42	14.816	1987	1994	1999	rBV	3827335	5604609	25.34%	1.445%
43	15.057	2030	2035	2041	rVB	98383	147597	0.67%	0.038%
44	15.263	2063	2070	2075	rBV	10366425	14671544	66.33%	3.783%
45	15.304	2075	2077	2084	rVB4	27809	38387	0.17%	0.010%
46	15.416	2089	2096	2100	rBV	1941270	2764686	12.50%	0.713%
47	15.475	2100	2106	2112	rVB	12277386	16800041	75.95%	4.332%
48	15.557	2114	2120	2129	rBV	628435	845250	3.82%	0.218%
49	15.657	2133	2137	2143	rVB8	13222	28191	0.13%	0.007%
50	15.722	2143	2148	2153	rVB2	50592	67895	0.31%	0.018%
51	15.969	2185	2190	2193	rBV4	17639	28115	0.13%	0.007%
52	16.016	2193	2198	2202	rVV	59445	92073	0.42%	0.024%
53	16.057	2202	2205	2214	rVB5	19476	42643	0.19%	0.011%
54	16.151	2214	2221	2227	rBV4	70027	142592	0.64%	0.037%
55	16.334	2246	2252	2256	rBV	87937	103514	0.47%	0.027%
56	16.487	2268	2278	2283	rBV	15242750	22118888	100.00%	5.703%
57	16.834	2329	2337	2346	rBV	9021432	12871365	58.19%	3.319%
58	16.898	2346	2348	2351	rVV	53786	71511	0.32%	0.018%
59	16.928	2351	2353	2356	rVV2	35975	45815	0.21%	0.012%
60	16.975	2356	2361	2373	rVB2	135100	250054	1.13%	0.064%
61	17.175	2387	2395	2399	rVV	1260463	1938854	8.77%	0.500%
62	17.222	2399	2403	2408	rVV	4644711	6182282	27.95%	1.594%
63	17.281	2408	2413	2415	rVV	1967559	3025608	13.68%	0.780%
64	17.316	2415	2419	2430	rVB	6319624	8519282	38.52%	2.197%
65	17.463	2437	2444	2449	rVB3	30478	64147	0.29%	0.017%
66	17.592	2458	2466	2472	rVB3	50252	93294	0.42%	0.024%
67	17.722	2484	2488	2492	rBV6	18515	28869	0.13%	0.007%
68	17.787	2495	2499	2505	rVV2	128593	183221	0.83%	0.047%
69	17.851	2505	2510	2516	rVB	303483	380080	1.72%	0.098%
70	18.075	2543	2548	2554	rBV	642895	872021	3.94%	0.225%
71	18.151	2554	2561	2565	rVV	11721471	14769915	66.78%	3.808%
72	18.275	2580	2582	2590	rVB6	22570	30213	0.14%	0.008%
73	18.345	2590	2594	2598	rBV4	33998	57739	0.26%	0.015%
74	18.422	2603	2607	2613	rVV	110527	178639	0.81%	0.046%
75	18.475	2613	2616	2620	rVV4	28422	45536	0.21%	0.012%
76	18.516	2620	2623	2630	rVV	69028	97807	0.44%	0.025%

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 Title : SVOA CALIBRATION

77	18.587	2630	2635	2640	rVV	560780	685038	3.10%	0.177%
78	18.957	2694	2698	2700	rBV3	85701	100145	0.45%	0.026%
79	18.986	2700	2703	2707	rVB	330536	350058	1.58%	0.090%
80	19.116	2718	2725	2731	rBV3	134374	232284	1.05%	0.060%
81	19.204	2732	2740	2745	rVV	14994732	21098164	95.39%	5.440%
82	19.310	2754	2758	2766	rBV	346002	445630	2.01%	0.115%
83	19.434	2775	2779	2784	rBV	143882	181173	0.82%	0.047%
84	19.528	2789	2795	2796	rVV	2731017	3560391	16.10%	0.918%
85	19.557	2796	2800	2807	rVB	7563738	9771180	44.18%	2.519%
86	19.622	2808	2811	2815	rBV3	81050	98660	0.45%	0.025%
87	19.810	2841	2843	2848	rVB6	32766	35704	0.16%	0.009%
88	20.022	2876	2879	2881	rBV2	50519	66159	0.30%	0.017%
89	20.186	2904	2907	2912	rBV	192068	208873	0.94%	0.054%
90	20.986	3040	3043	3054	rVB	468425	654684	2.96%	0.169%
91	21.122	3063	3066	3073	rBV	325413	376477	1.70%	0.097%
92	21.198	3073	3079	3083	rBV	15726946	17186677	77.70%	4.431%
93	21.275	3086	3092	3097	rVV	11723006	16682278	75.42%	4.301%
94	21.328	3097	3101	3106	rVB	15604740	19722204	89.16%	5.085%
95	22.028	3216	3220	3224	rVB2	249462	325028	1.47%	0.084%
96	22.898	3365	3368	3373	rBV	155808	235070	1.06%	0.061%
97	22.980	3373	3382	3390	rVB	10992562	20350245	92.00%	5.247%
98	23.486	3461	3468	3481	rVB2	1593448	3267553	14.77%	0.842%
99	23.639	3488	3494	3507	rVB	932702	1875082	8.48%	0.483%
100	26.116	3902	3915	3932	rVB	4606037	13815096	62.46%	3.562%

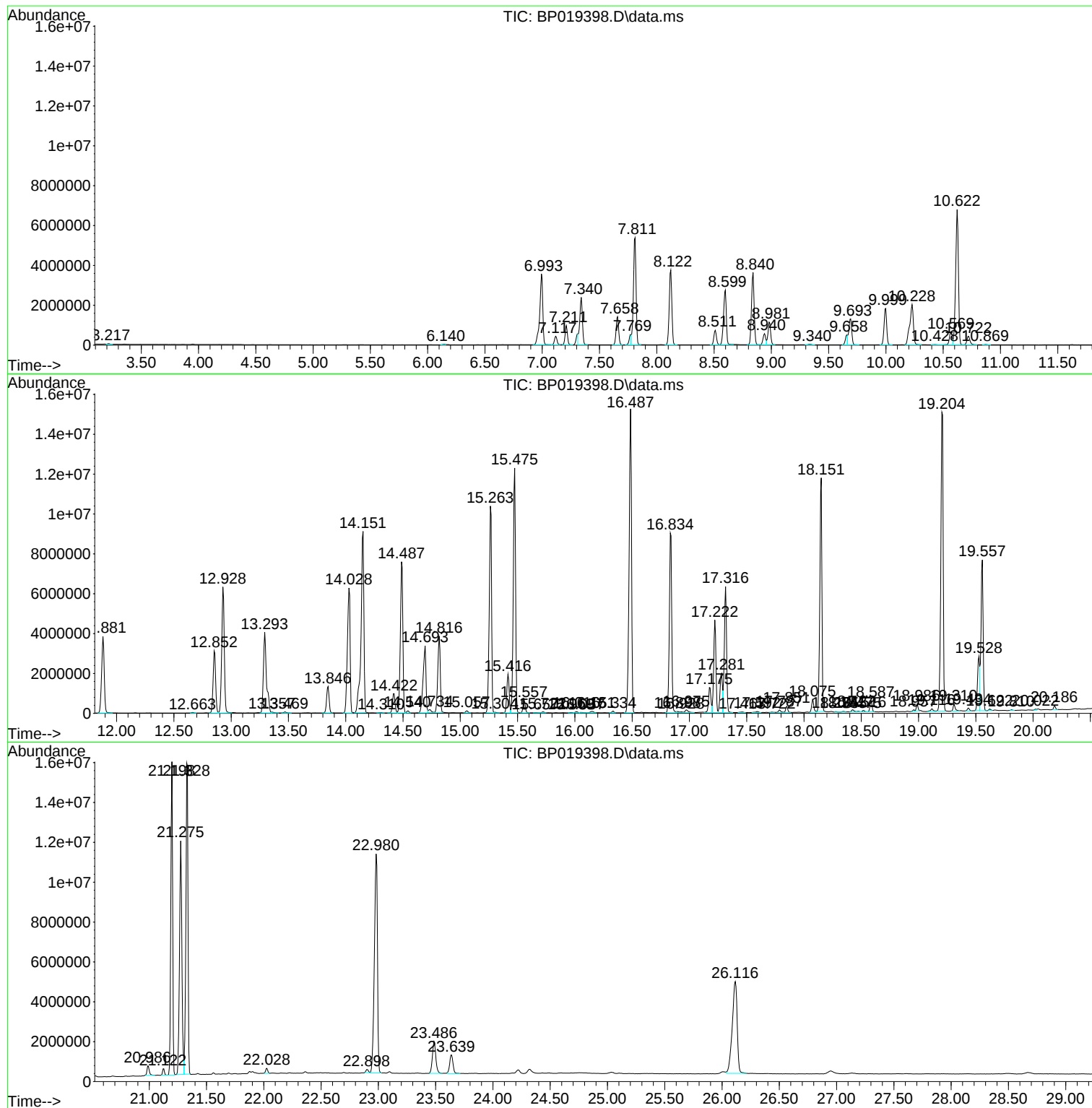
Sum of corrected areas: 387845892

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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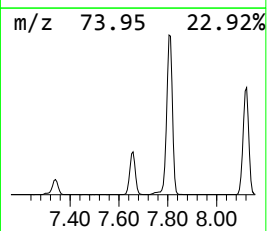
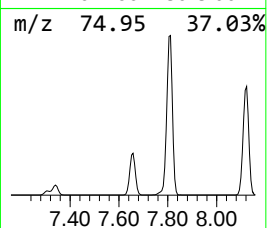
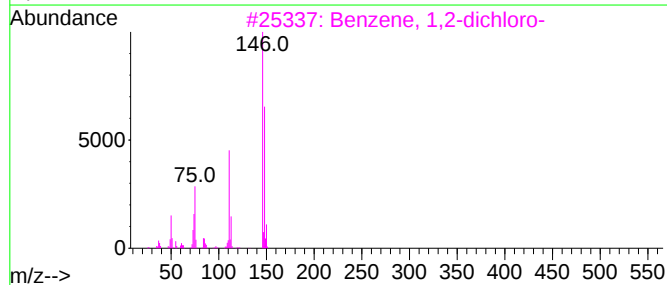
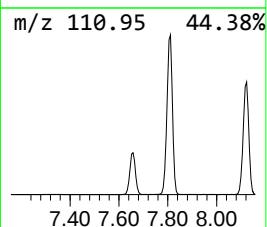
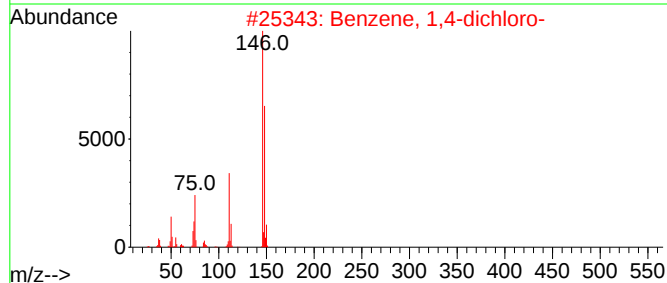
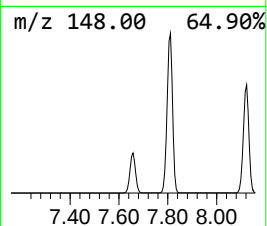
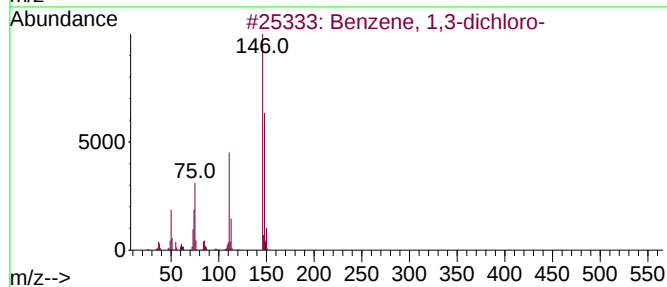
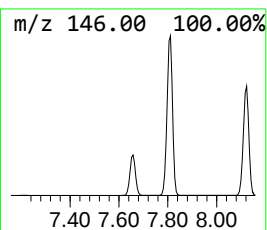
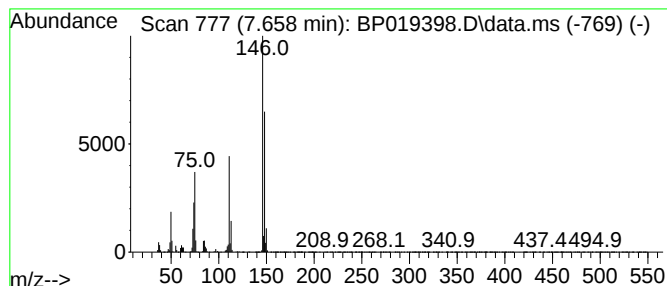
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,3-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.658	67.80 ng/ul	2207580	1,4-Dichlorobenzene-d4	7.769

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	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, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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ClientSampleId :
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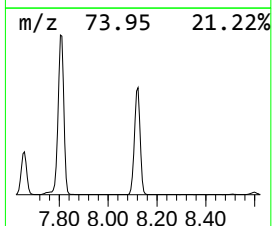
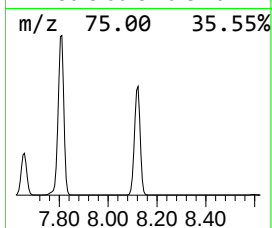
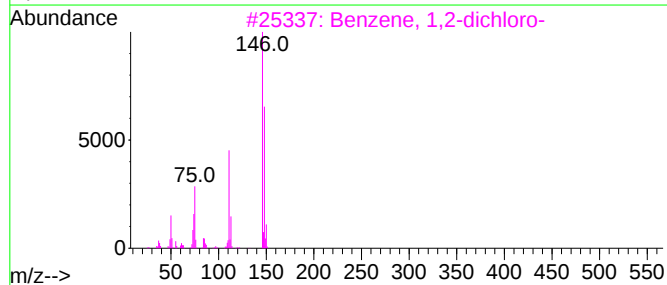
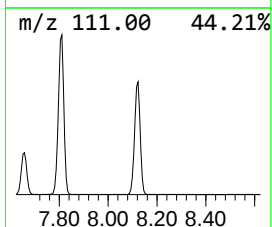
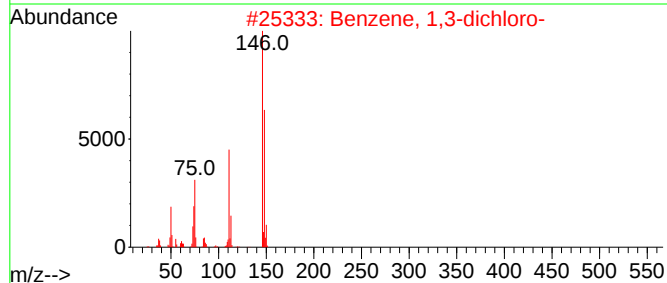
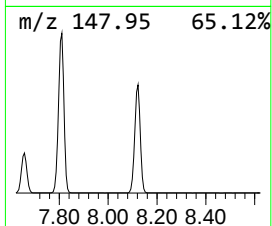
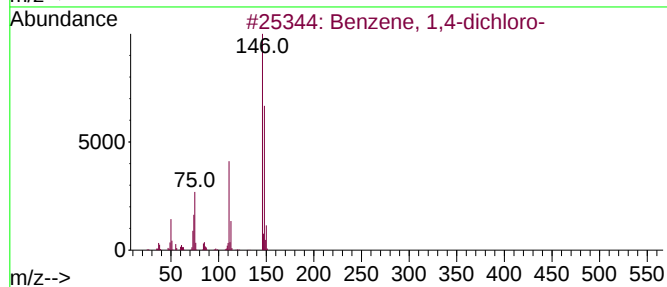
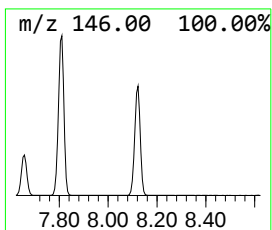
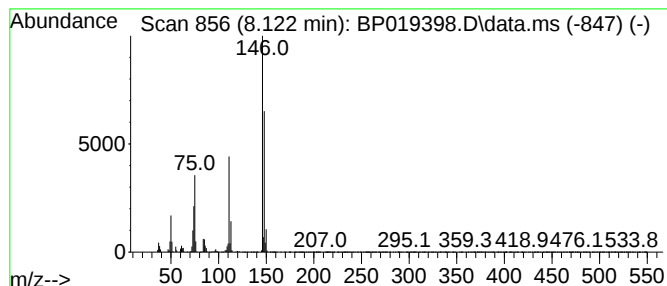
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	183.64 ng/ul	5979180	1,4-Dichlorobenzene-d4	7.769

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



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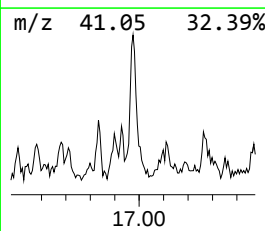
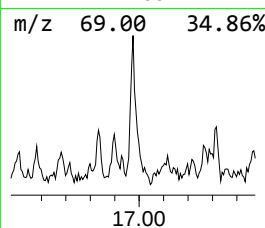
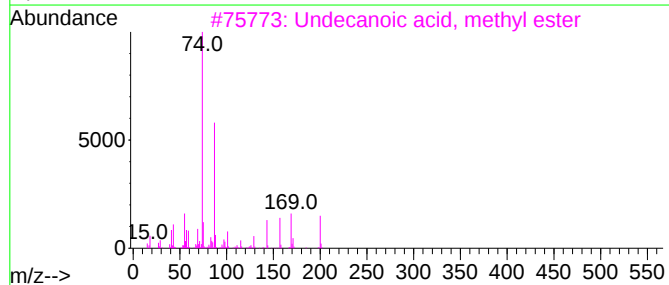
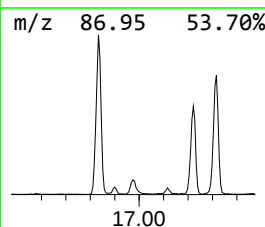
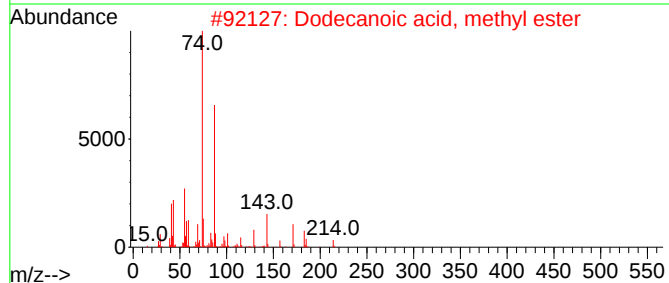
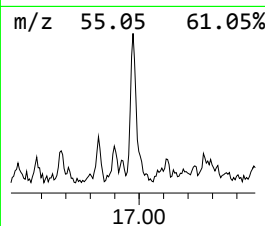
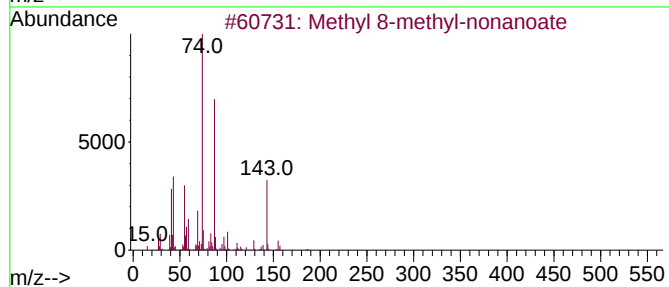
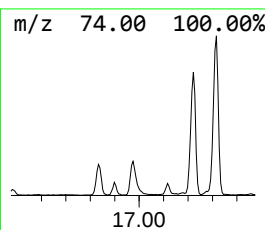
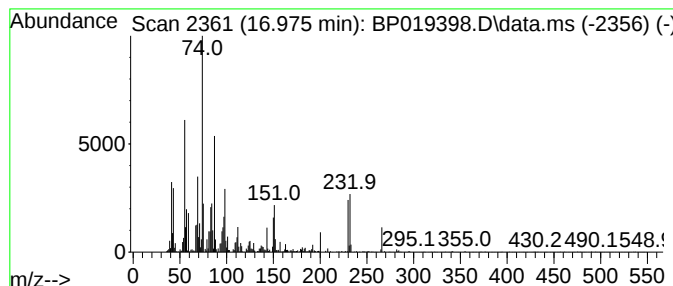
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.975	2.58 ng/ul	250054	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 8-methyl-nonanoate	186	C11H22O2	1000452-00-9	43
2			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	38
3			Undecanoic acid, methyl ester	200	C12H24O2	001731-86-8	38
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	38
5			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
Operator : MA/JU
Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

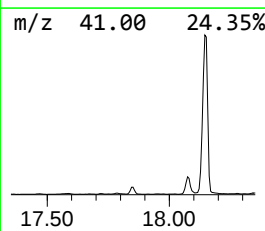
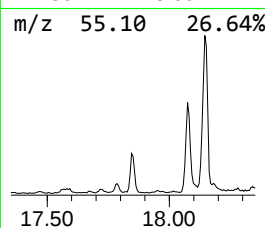
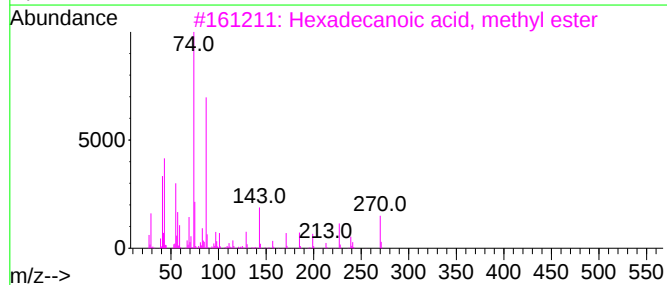
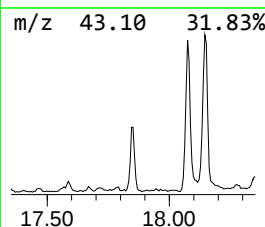
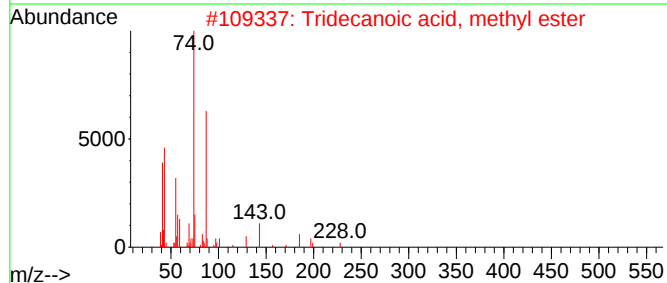
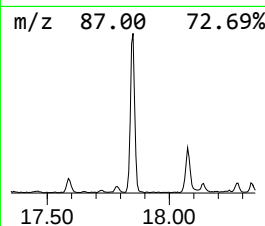
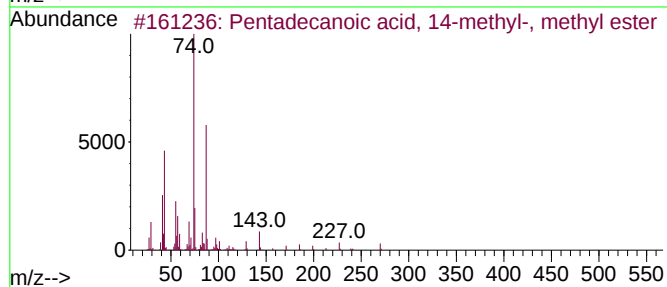
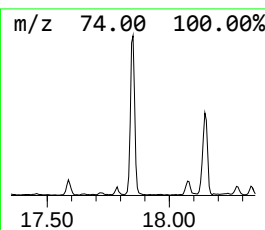
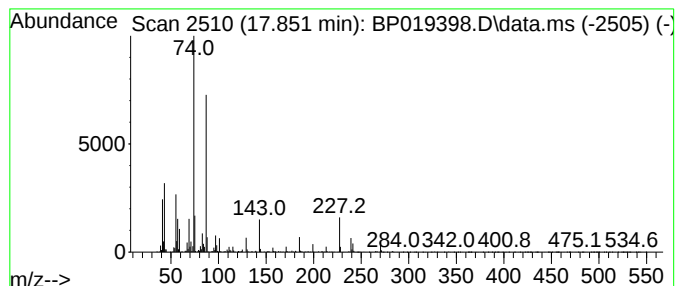
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Pentadecanoic acid, 14-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.851	3.92 ng/ul	380080	Phenanthrene-d10	17.175	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	97
2	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
3	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
4	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
Operator : MA/JU
Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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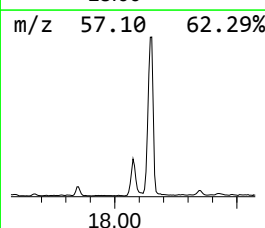
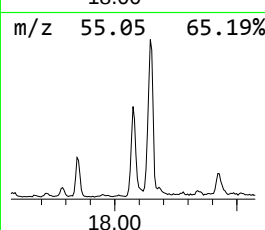
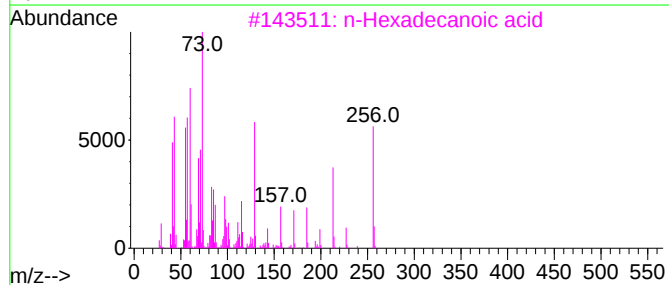
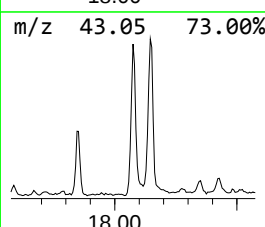
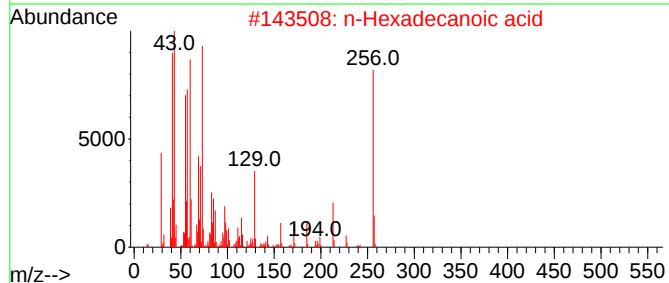
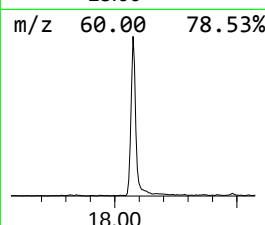
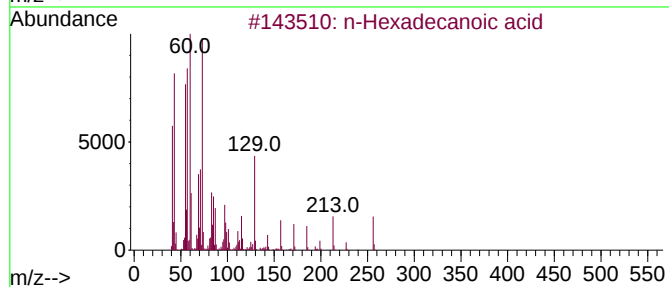
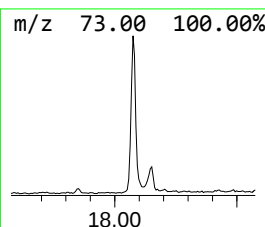
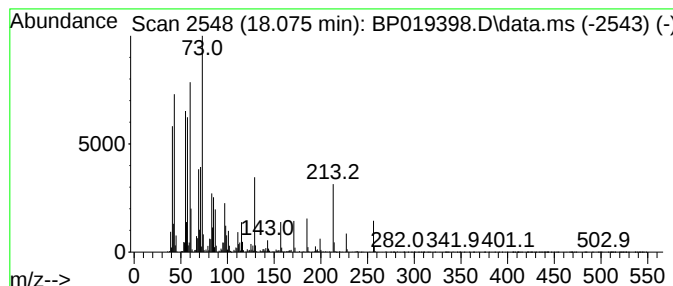
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.075	9.00 ng/ul	872021	Phenanthrene-d10	17.175

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
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Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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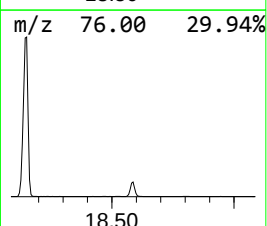
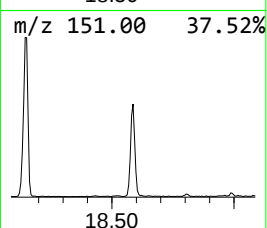
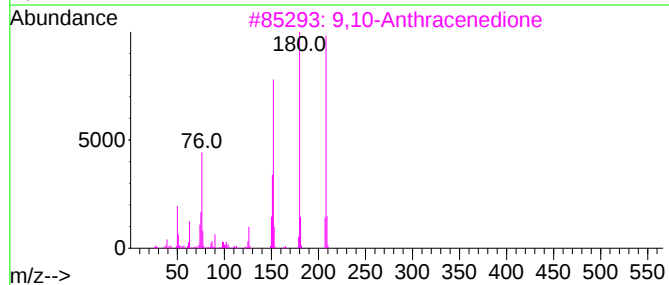
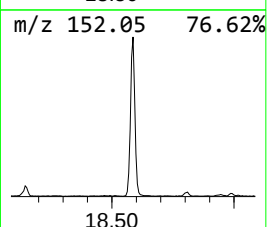
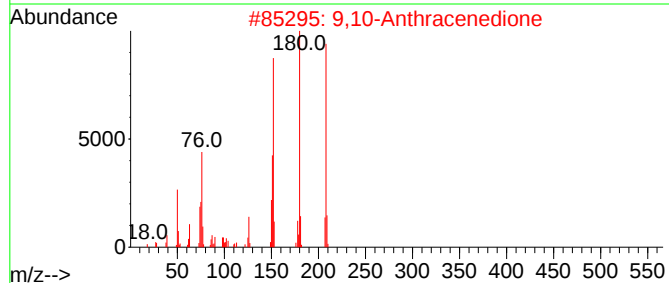
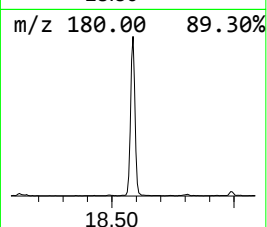
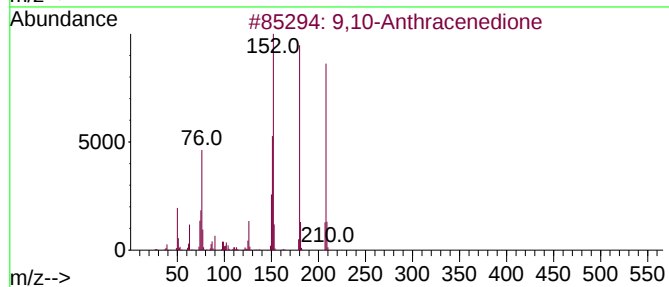
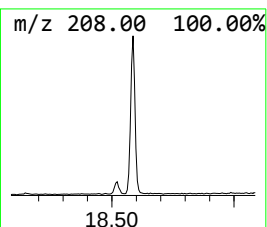
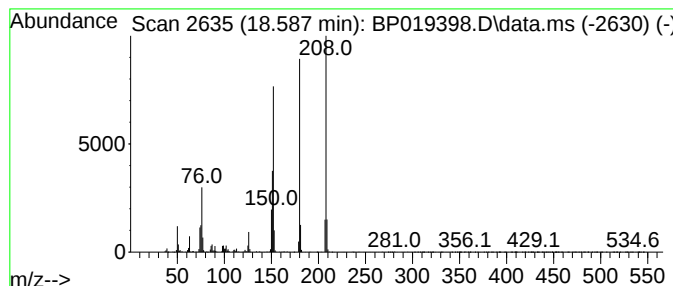
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.587	7.07 ng/ul	685038	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	91
5	1H-Phenalen-1-one			180	C13H8O	000548-39-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
Operator : MA/JU
Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

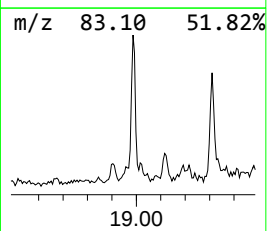
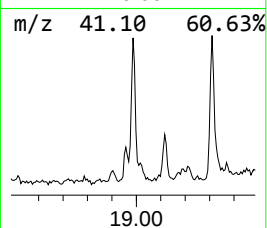
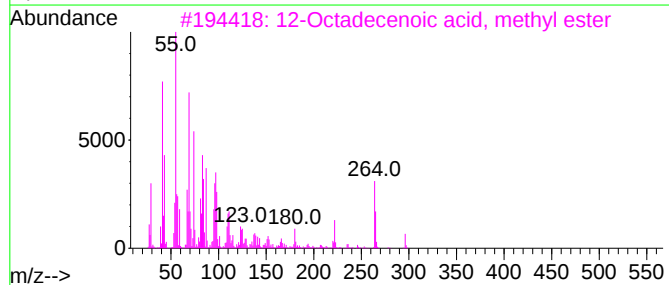
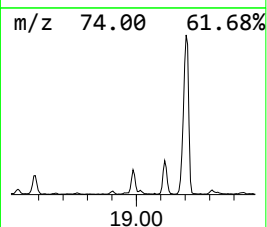
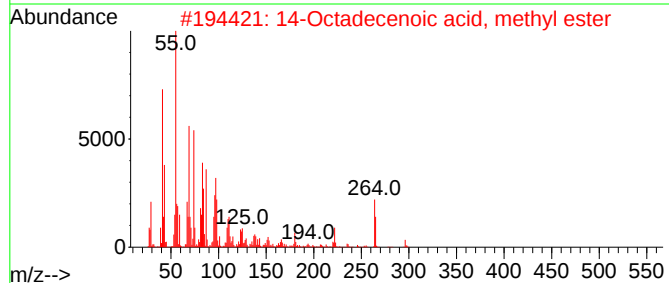
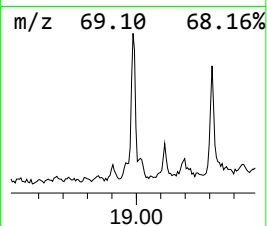
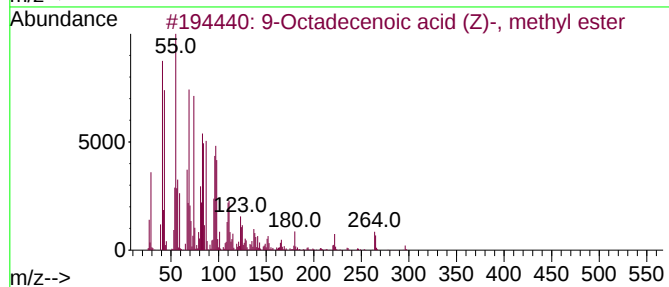
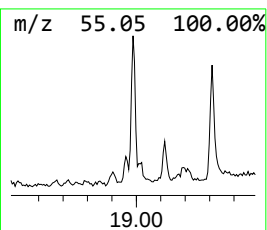
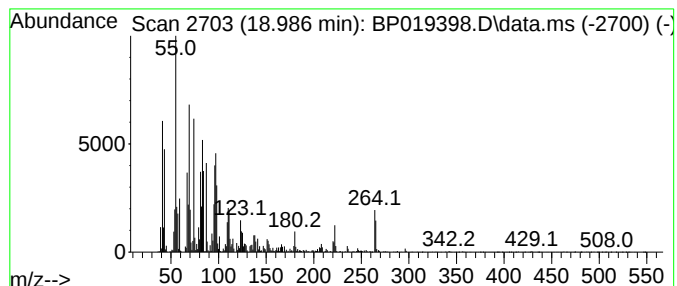
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 9-Octadecenoic acid (Z)-, m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.986	3.61 ng/ul	350058	Phenanthrene-d10	17.175	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
3	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
4	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	98
5	6-Octadecenoic acid, methyl este...	296	C19H36O2	002777-58-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
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Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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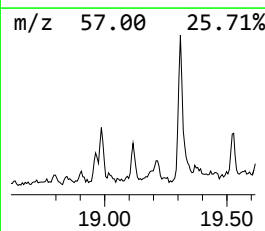
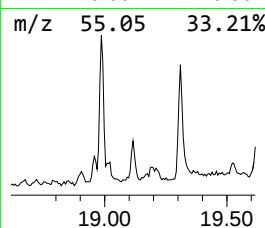
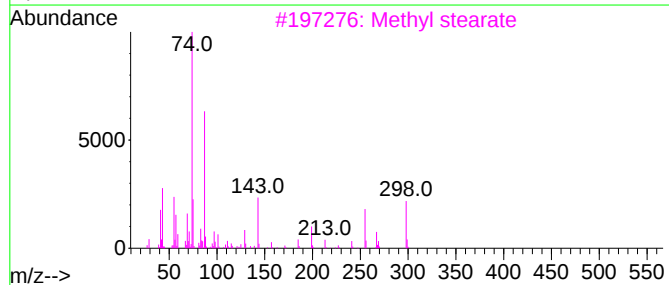
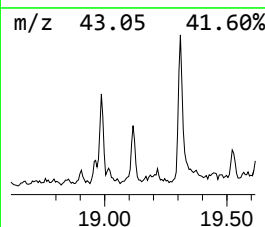
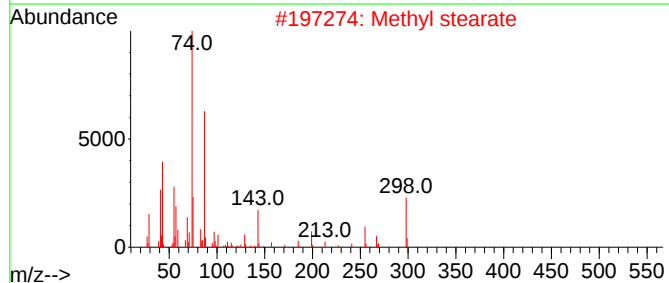
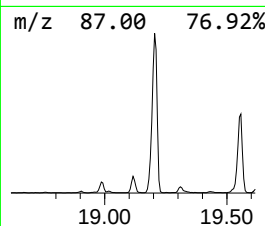
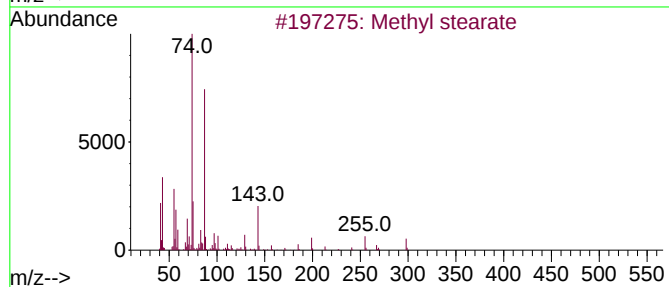
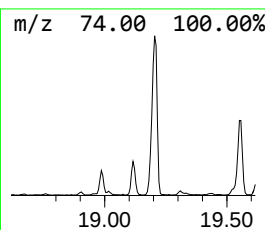
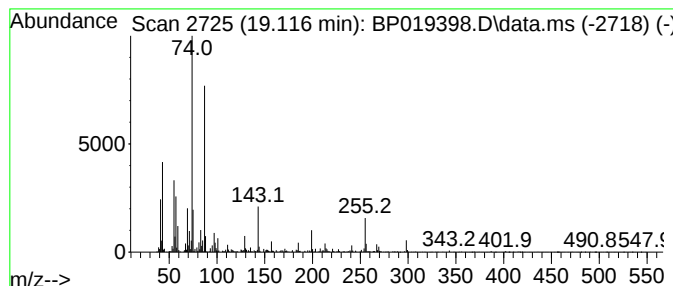
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Methyl stearate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.116	2.40 ng/ul	232284	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	96
2			Methyl stearate	298	C19H38O2	000112-61-8	95
3			Methyl stearate	298	C19H38O2	000112-61-8	94
4			Methyl stearate	298	C19H38O2	000112-61-8	93
5			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
Operator : MA/JU
Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

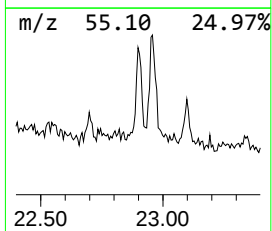
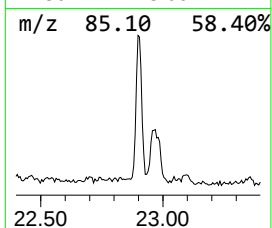
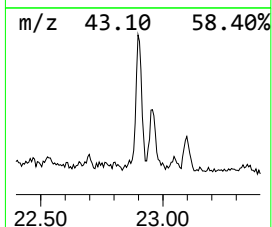
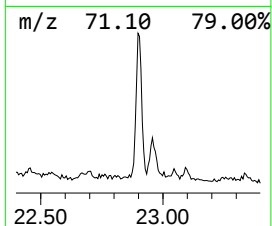
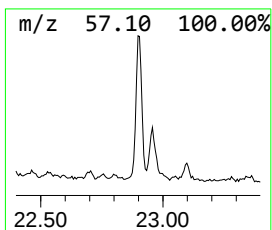
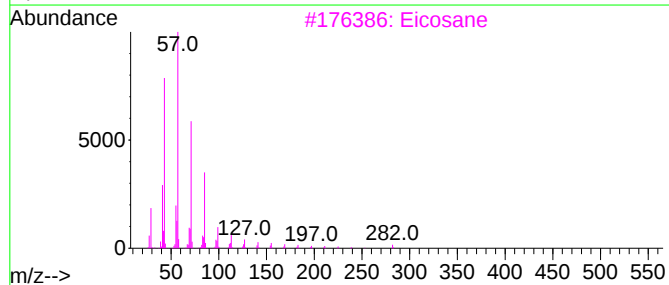
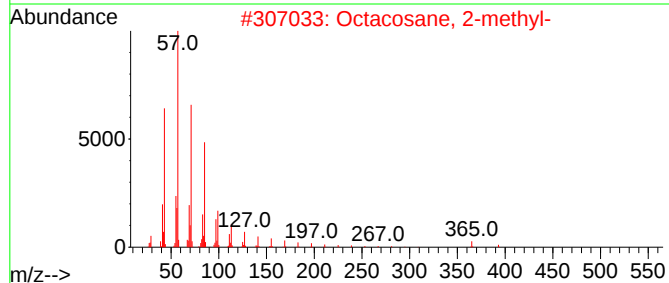
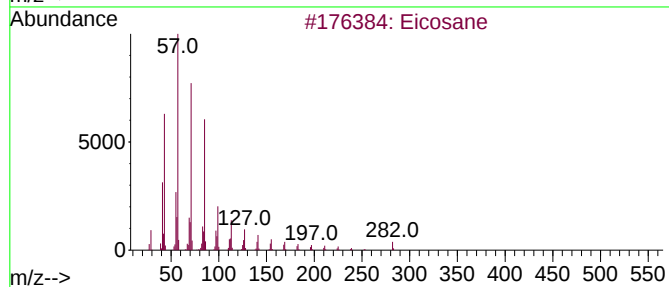
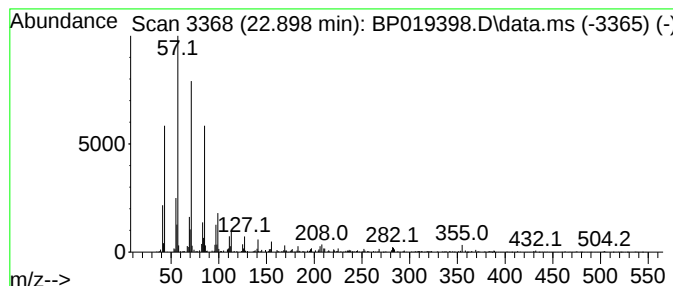
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.898	2.51 ng/ul	235070	Perylene-d12	23.639	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
3	Eicosane	282	C20H42	000112-95-8	90
4	Tetracosane	338	C24H50	000646-31-1	90
5	Nonadecane	268	C19H40	000629-92-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011924\
Data File : BP019398.D
Acq On : 19 Jan 2024 10:29
Operator : MA/JU
Sample : P1160-09
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.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
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Benzene, 1,4-di...	8.122	183.6	ng/ul	5979180	1	7.769	651198	20.0
unknown-01	16.975	2.6	ng/ul	250054	4	17.175	1938850	20.0
Pentadecanoic a...	17.851	3.9	ng/ul	380080	4	17.175	1938850	20.0
n-Hexadecanoic ...	18.075	9.0	ng/ul	872021	4	17.175	1938850	20.0
9,10-Anthracene...	18.587	7.1	ng/ul	685038	4	17.175	1938850	20.0
9-Octadecenoic ...	18.986	3.6	ng/ul	350058	4	17.175	1938850	20.0
Methyl stearate	19.116	2.4	ng/ul	232284	4	17.175	1938850	20.0
(DEL) Alkane: S...	22.898	2.5	ng/ul	235070	6	23.639	1875080	20.0