

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
 Data File : BP023796.D
 Acq On : 30 Jan 2025 12:29
 Operator : RC/JU
 Sample : Q1190-19
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44X3

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-BP012925.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023796.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.293	32	35	44	rVB	69153	103302	0.19%	0.014%
2	4.028	156	160	164	rBV	54034	66048	0.12%	0.009%
3	4.246	194	197	203	rVB	46745	62223	0.11%	0.008%
4	6.169	519	524	535	rBV2	68849	152248	0.28%	0.020%
5	6.987	651	663	676	rVB2	2582874	5543483	10.06%	0.736%
6	7.110	678	684	693	rBV	886711	1359491	2.47%	0.180%
7	7.205	693	700	713	rVV	4021564	5784926	10.50%	0.768%
8	7.322	713	720	722	rVV	1115635	1757878	3.19%	0.233%
9	7.352	722	725	733	rVB	2156646	3141540	5.70%	0.417%
10	7.669	771	779	787	rBV	3710128	5864422	10.64%	0.778%
11	7.775	789	797	800	rVV	1267857	2215410	4.02%	0.294%
12	7.816	800	804	818	rVB	3227712	4991376	9.06%	0.663%
13	8.016	833	838	846	rVB	35591	59532	0.11%	0.008%
14	8.134	849	858	869	rVB	11567214	19509721	35.39%	2.590%
15	8.487	911	918	926	rBV	1392645	2032695	3.69%	0.270%
16	8.581	926	934	946	rVV	7529426	11614550	21.07%	1.542%
17	8.922	984	992	995	rBV	1005340	1693634	3.07%	0.225%
18	8.963	995	999	1008	rVB	4768859	7447507	13.51%	0.989%
19	9.334	1055	1062	1065	rBV	44971	78425	0.14%	0.010%
20	9.487	1077	1088	1092	rVB5	31996	70278	0.13%	0.009%
21	9.669	1105	1119	1131	rBV2	4550409	8579133	15.56%	1.139%
22	9.969	1162	1170	1181	rBV	4580486	6969963	12.64%	0.925%
23	10.210	1200	1211	1226	rBV2	8090452	15022067	27.25%	1.994%
24	10.551	1259	1269	1278	rBV	1837762	3196966	5.80%	0.424%
25	10.687	1284	1292	1308	rBV	1132685	1993939	3.62%	0.265%
26	10.851	1315	1320	1326	rVB2	41196	69282	0.13%	0.009%
27	11.846	1481	1489	1502	rBV	10374441	16353927	29.67%	2.171%
28	12.534	1600	1606	1612	rBV3	64133	110911	0.20%	0.015%
29	12.622	1617	1621	1626	rVB	72848	99176	0.18%	0.013%
30	12.822	1648	1655	1662	rBV	5094035	7675034	13.92%	1.019%
31	12.898	1662	1668	1682	rVB	7579764	11766921	21.35%	1.562%
32	13.228	1721	1724	1728	rVV	47403	60245	0.11%	0.008%
33	13.298	1732	1736	1744	rVV	105824	161952	0.29%	0.021%
34	13.804	1815	1822	1829	rBV	3029926	4256632	7.72%	0.565%
35	13.969	1840	1850	1856	rBV	6673330	9456687	17.16%	1.255%
36	14.087	1861	1870	1881	rVB	3001826	4884522	8.86%	0.648%

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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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Title : SVOA CALIBRATION

37	14.398	1914	1923	1928	rBV	2975020	4519590	8.20%	0.600%
38	14.469	1928	1935	1941	rVV	20097195	31356582	56.89%	4.162%
39	14.528	1941	1945	1952	rVB	418852	530226	0.96%	0.070%
40	14.634	1953	1963	1974	rBV2	5593066	9608864	17.43%	1.276%
41	14.763	1974	1985	1994	rVV	10829479	18116775	32.87%	2.405%
42	14.828	1994	1996	2004	rVB	97676	137296	0.25%	0.018%
43	14.957	2014	2018	2023	rVB	91587	119936	0.22%	0.016%
44	15.028	2023	2030	2040	rVB2	146525	263624	0.48%	0.035%
45	15.234	2056	2065	2075	rBV	9474321	11607360	21.06%	1.541%
46	15.404	2088	2094	2099	rBV	4428656	6762633	12.27%	0.898%
47	15.469	2099	2105	2110	rVV	15782743	24120438	43.76%	3.202%
48	15.516	2110	2113	2123	rVB	1399757	1763424	3.20%	0.234%
49	15.716	2142	2147	2152	rVV	224464	336688	0.61%	0.045%
50	15.775	2152	2157	2164	rVB	1735198	2206086	4.00%	0.293%
51	15.957	2184	2188	2198	rVB4	36649	82613	0.15%	0.011%
52	16.139	2211	2219	2224	rBV3	177149	306871	0.56%	0.041%
53	16.369	2249	2258	2265	rVV	9128126	12743807	23.12%	1.692%
54	16.592	2292	2296	2300	rBV3	42857	68102	0.12%	0.009%
55	16.845	2330	2339	2349	rBV	17327371	23822665	43.22%	3.162%
56	16.928	2350	2353	2358	rVV2	121748	223165	0.40%	0.030%
57	16.998	2360	2365	2376	rVV2	316790	615443	1.12%	0.082%
58	17.198	2393	2399	2403	rVV	3407089	5492195	9.96%	0.729%
59	17.245	2403	2407	2412	rVV	22545172	28080452	50.94%	3.728%
60	17.298	2412	2416	2419	rVV2	5466438	7772913	14.10%	1.032%
61	17.333	2419	2422	2429	rVB	7209237	9277748	16.83%	1.232%
62	17.616	2466	2470	2472	rBV	53428	76773	0.14%	0.010%
63	17.639	2472	2474	2479	rVB	66184	74253	0.13%	0.010%
64	17.769	2493	2496	2500	rVV4	46549	65673	0.12%	0.009%
65	17.833	2503	2507	2516	rVV3	112401	244976	0.44%	0.033%
66	17.916	2516	2521	2528	rVB	579833	738891	1.34%	0.098%
67	18.086	2546	2550	2554	rBV6	60502	88354	0.16%	0.012%
68	18.139	2554	2559	2565	rBV	466617	810214	1.47%	0.108%
69	18.204	2565	2570	2576	rVV2	24213074	34443071	62.49%	4.572%
70	18.445	2605	2611	2618	rBV5	145366	308542	0.56%	0.041%
71	18.516	2618	2623	2633	rVB4	274787	553213	1.00%	0.073%
72	18.604	2634	2638	2642	rVV3	118895	207898	0.38%	0.028%
73	18.657	2642	2647	2654	rVB	852034	1151778	2.09%	0.153%
74	18.886	2683	2686	2689	rVB2	75772	83623	0.15%	0.011%
75	19.004	2702	2706	2717	rVB3	273111	635080	1.15%	0.084%
76	19.122	2723	2726	2731	rVB4	129155	176122	0.32%	0.023%

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77	19.222	2740	2743	2747	rBV	85856	119543	0.22%	0.016%
78	19.269	2747	2751	2754	rBV	401571	498259	0.90%	0.066%
79	19.327	2755	2761	2768	rVB	19369177	24928706	45.23%	3.309%
80	19.469	2782	2785	2788	rBV2	117192	172116	0.31%	0.023%
81	19.580	2799	2804	2810	rBV	803474	1065882	1.93%	0.141%
82	19.675	2814	2820	2821	rVV	7018721	8613491	15.63%	1.143%
83	19.704	2821	2825	2836	rVB2	34420758	55120320	100.00%	7.317%
84	19.798	2838	2841	2846	rBV3	176316	268125	0.49%	0.036%
85	20.410	2941	2945	2949	rVB2	668949	873281	1.58%	0.116%
86	20.657	2981	2987	2992	rBV	18365859	22248092	40.36%	2.953%
87	21.321	3096	3100	3109	rBV	979116	1445546	2.62%	0.192%
88	21.498	3126	3130	3137	rVB	601384	856437	1.55%	0.114%
89	21.580	3138	3144	3148	rVV	21111390	34277162	62.19%	4.550%
90	21.633	3148	3153	3160	rVV2	16428067	31139394	56.49%	4.134%
91	21.698	3160	3164	3170	rVB	10283951	15382006	27.91%	2.042%
92	22.792	3347	3350	3357	rVB	435421	699127	1.27%	0.093%
93	22.886	3358	3366	3376	rBV	21407660	40679602	73.80%	5.400%
94	23.968	3542	3550	3555	rBV	5842455	13954769	25.32%	1.852%
95	24.039	3555	3562	3573	rVB	9715196	23946849	43.44%	3.179%
96	24.815	3683	3694	3698	rBV	3091480	9057783	16.43%	1.202%
97	24.880	3698	3705	3720	rVB	4328231	10707548	19.43%	1.421%
98	25.039	3725	3732	3744	rVB	2173003	6277947	11.39%	0.833%
99	28.868	4376	4383	4386	rBV	900710	2463038	4.47%	0.327%
100	28.980	4389	4402	4421	rVB	10114682	44735820	81.16%	5.938%

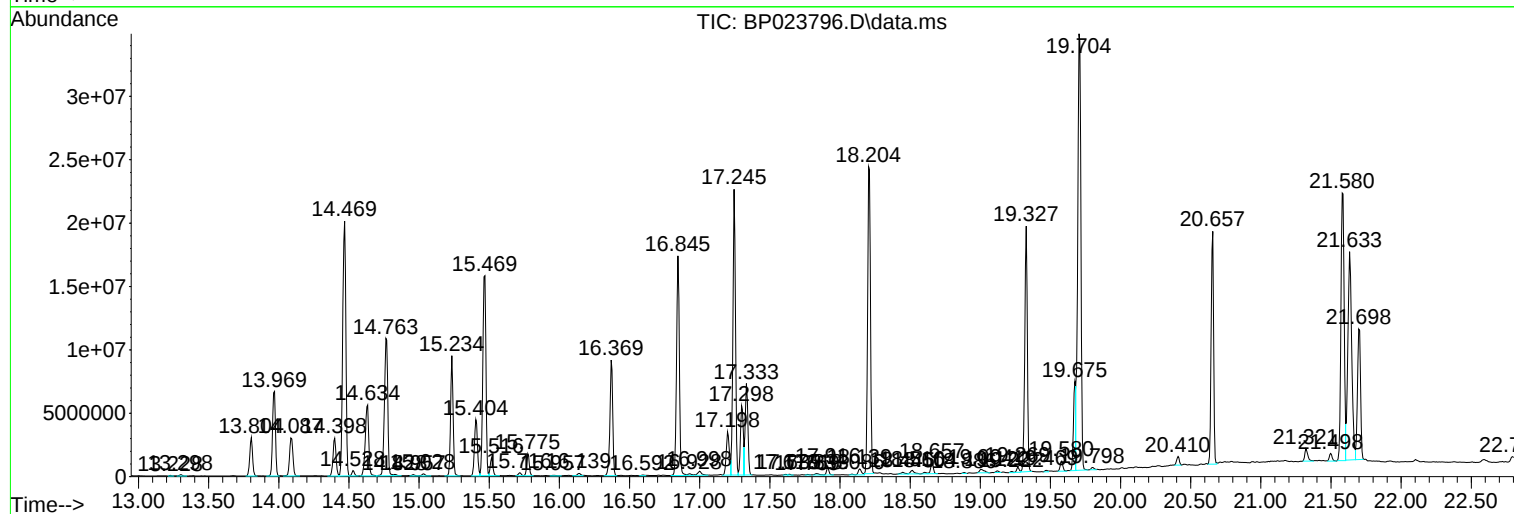
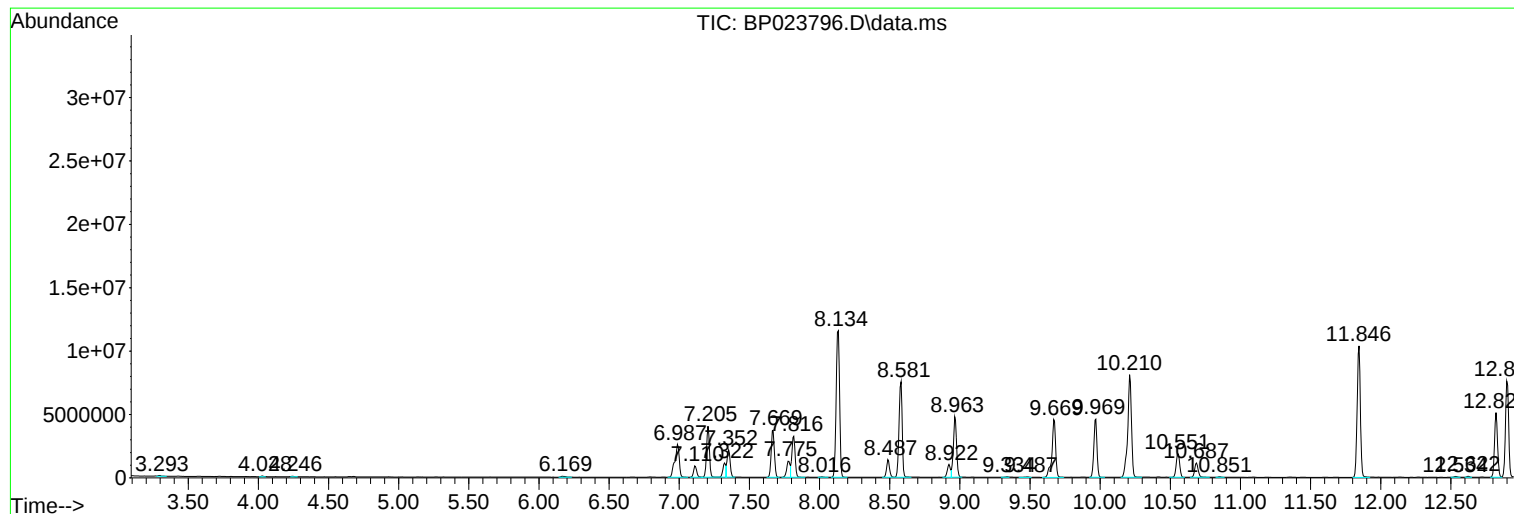
Sum of corrected areas: 753330846

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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



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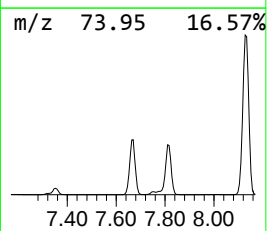
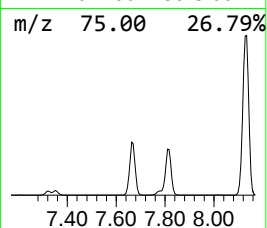
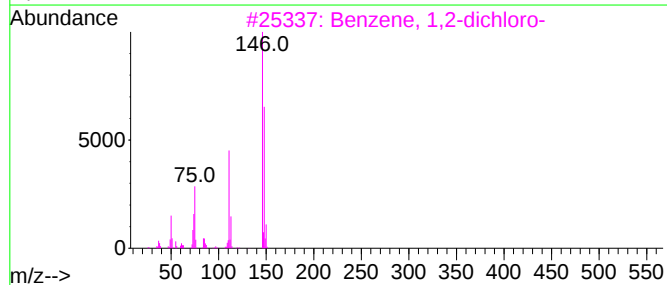
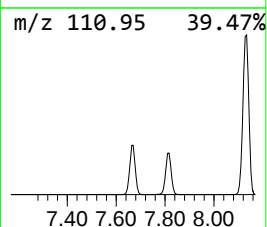
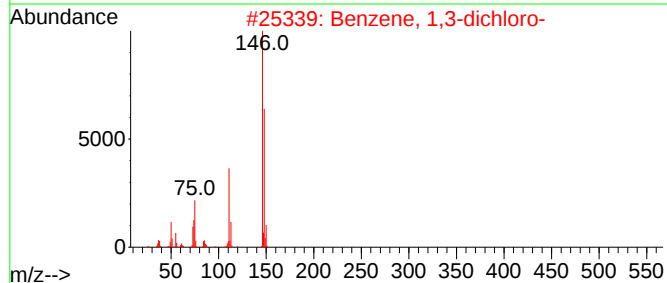
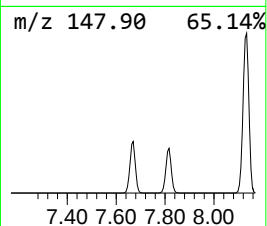
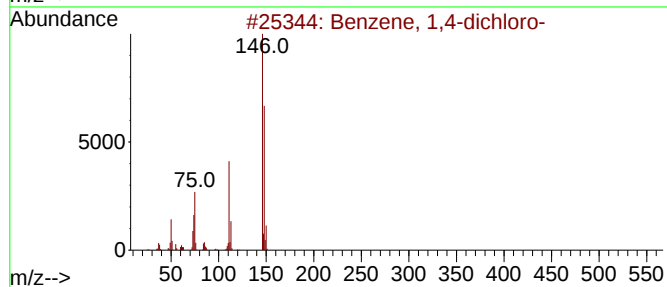
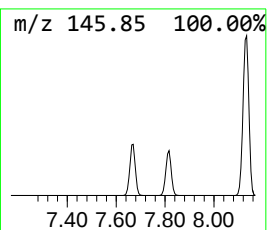
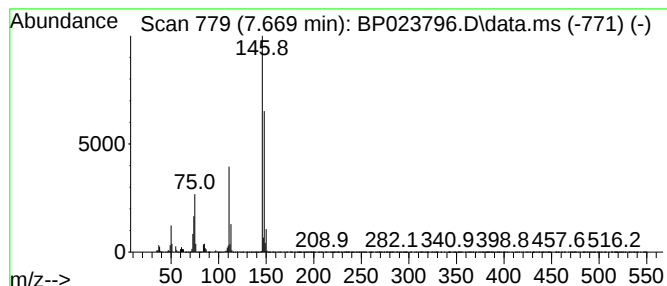
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.669	52.94 ng/ul	5864420	1,4-Dichlorobenzene-d4	7.775

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	97
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97



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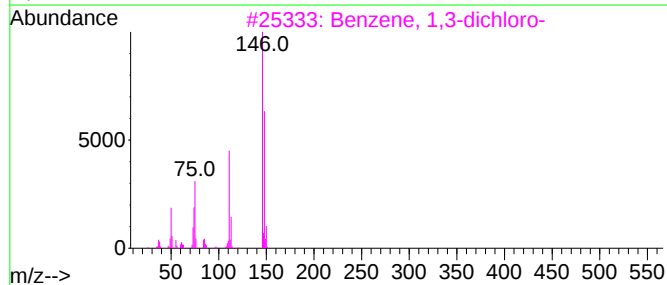
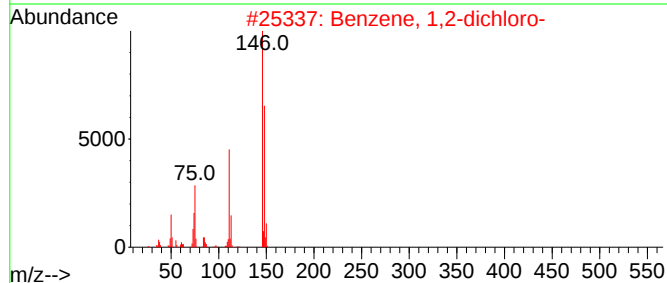
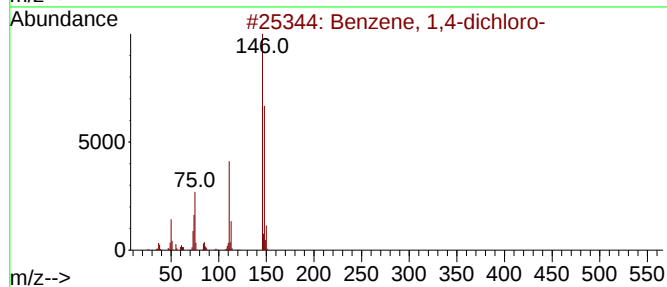
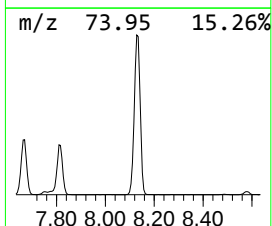
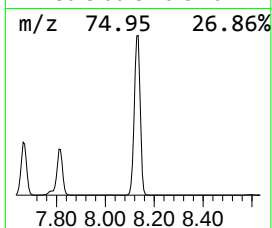
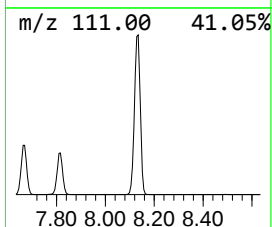
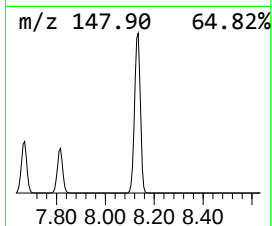
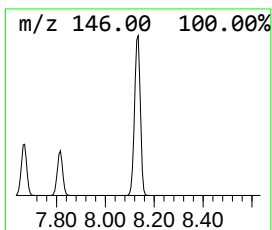
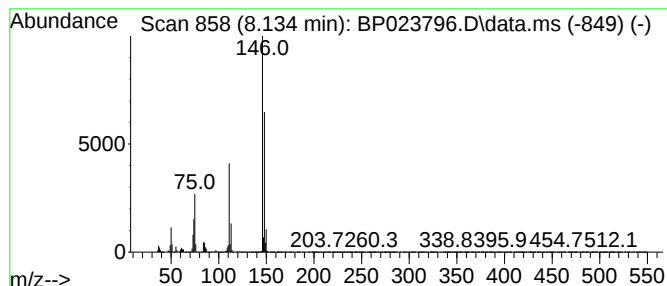
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.134	176.13 ng/ul	19509700	1,4-Dichlorobenzene-d4	7.775

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



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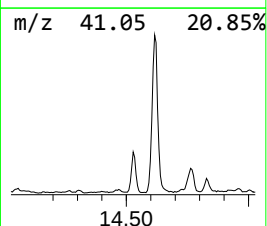
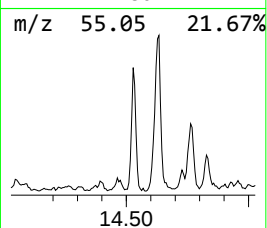
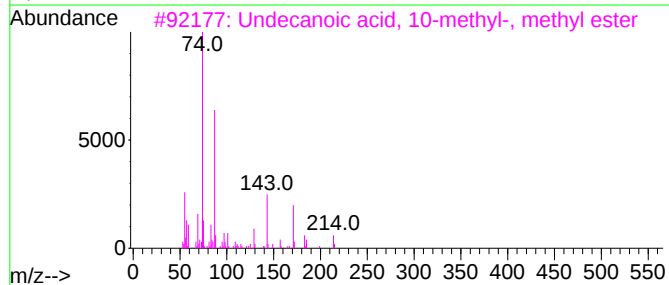
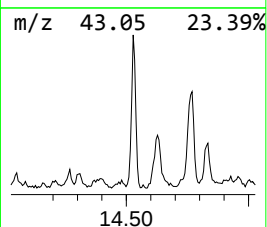
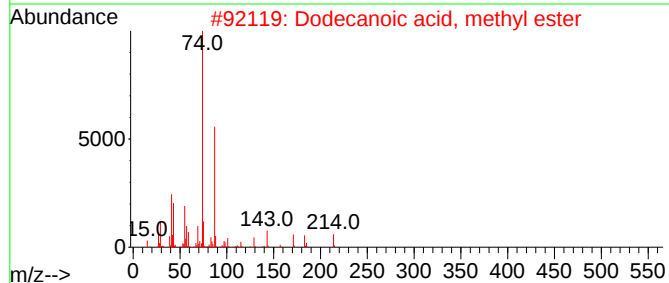
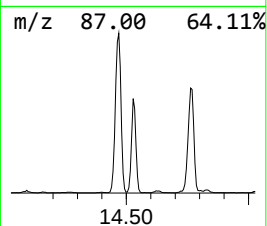
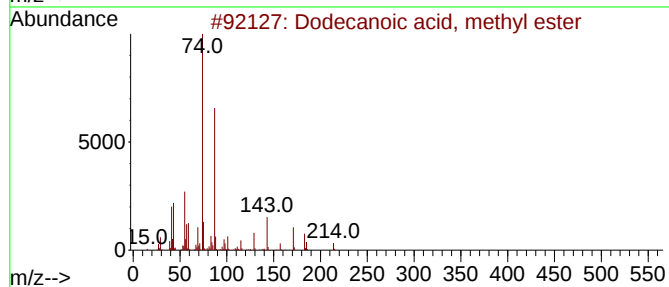
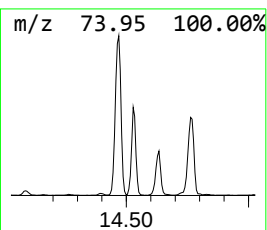
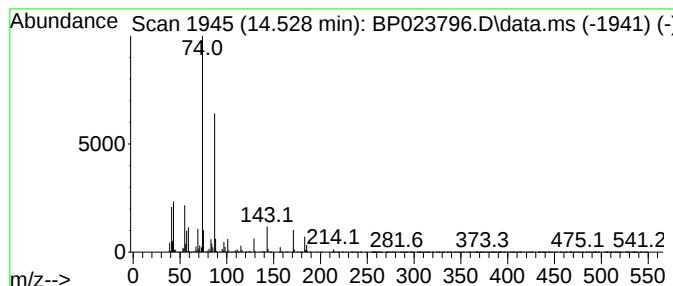
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Dodecanoic acid, methyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.528	2.35 ng/ul	530226	Acenaphthene-d10	14.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	91
2			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	91
3			Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	90
4			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	90
5			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
Operator : RC/JU
Sample : Q1190-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44X3

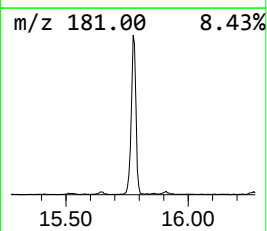
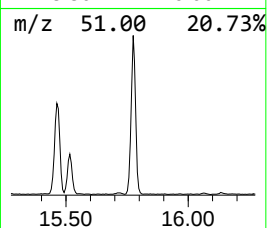
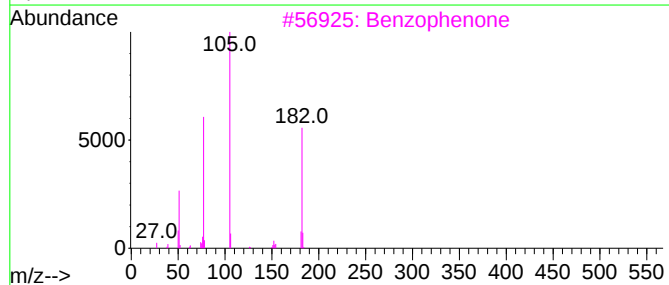
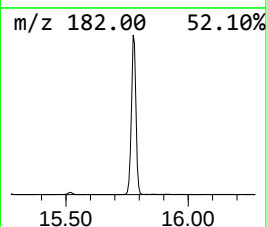
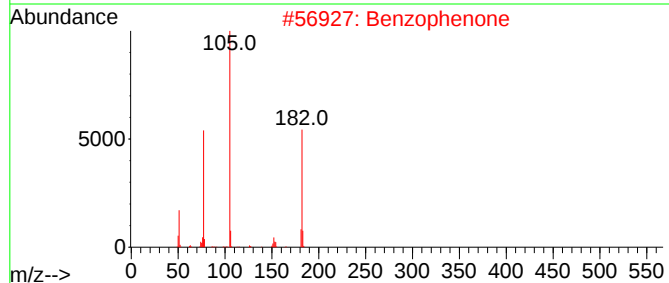
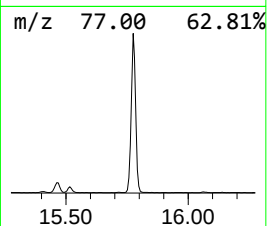
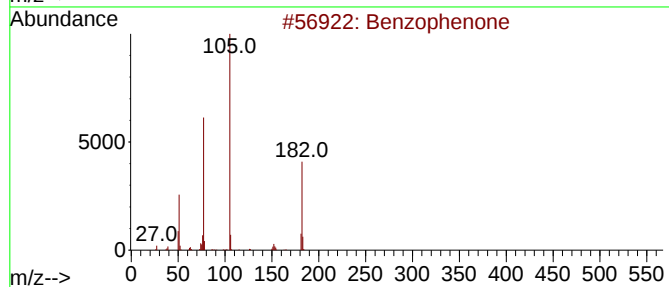
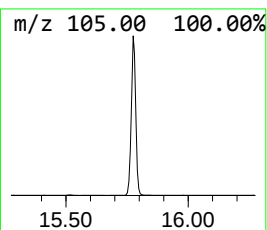
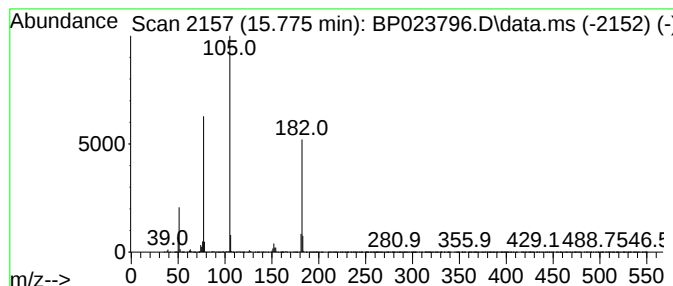
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	9.76 ng/ul	2206090	Acenaphthene-d10	14.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	95
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
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Sample : Q1190-19
Misc :
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Instrument :
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ClientSampleId :
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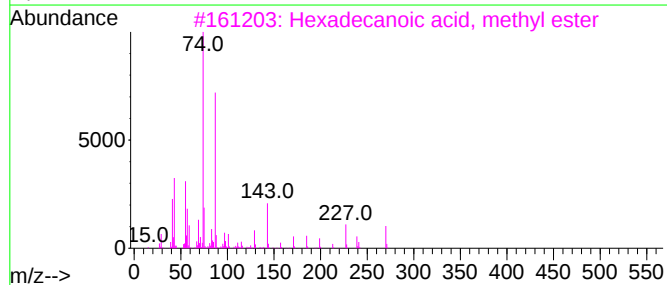
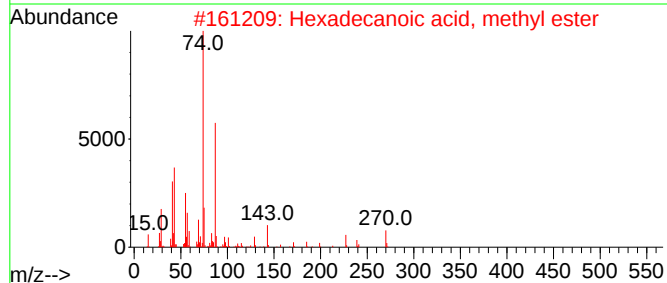
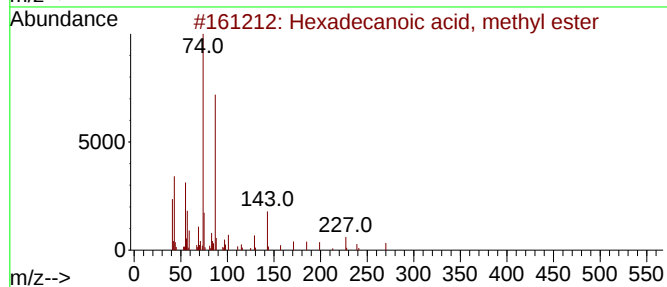
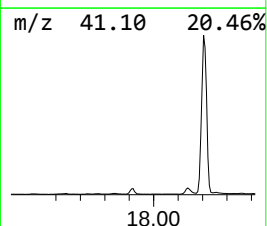
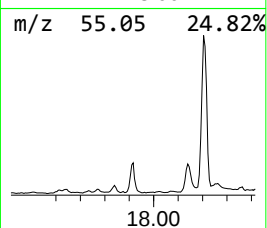
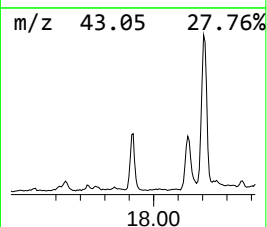
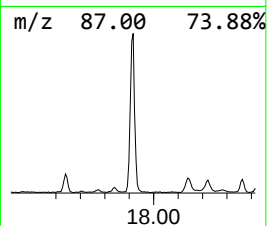
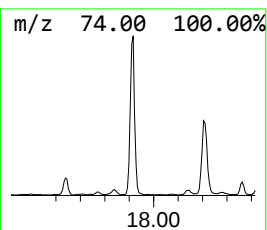
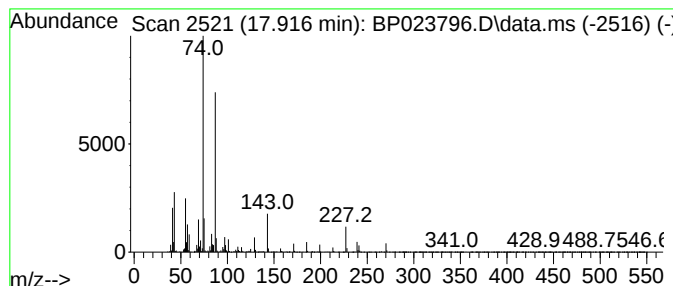
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.916	2.69 ng/ul	738891	Phenanthrene-d10	17.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	96
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	92
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
Operator : RC/JU
Sample : Q1190-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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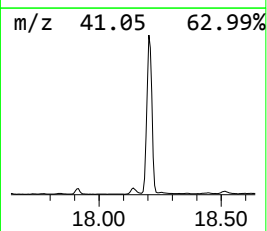
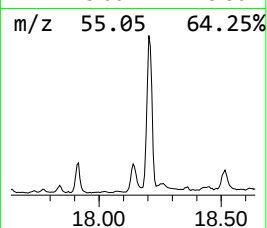
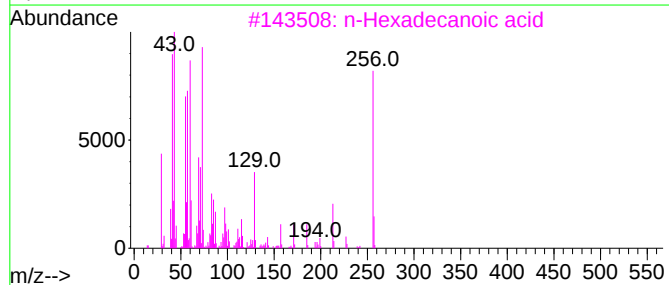
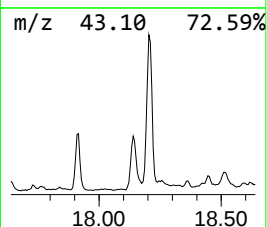
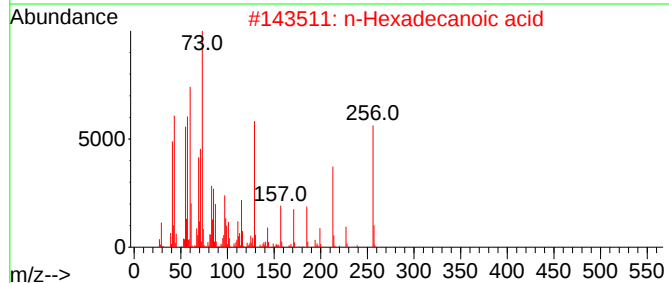
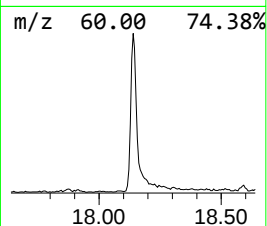
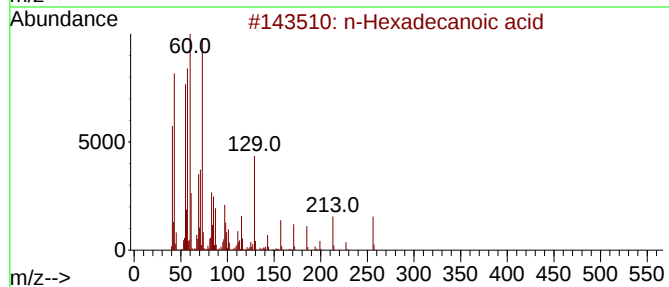
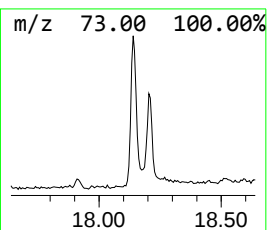
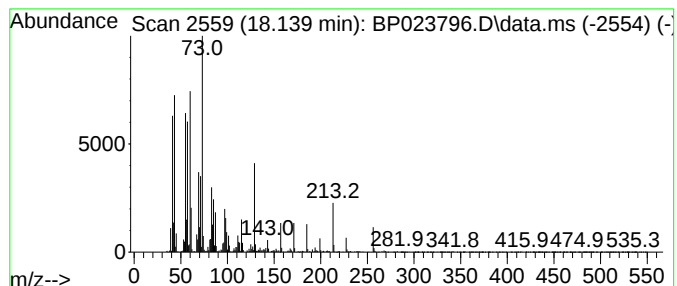
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	2.95 ng/ul	810214	Phenanthrene-d10	17.198

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
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Sample : Q1190-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

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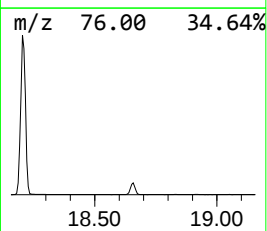
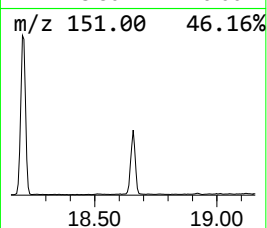
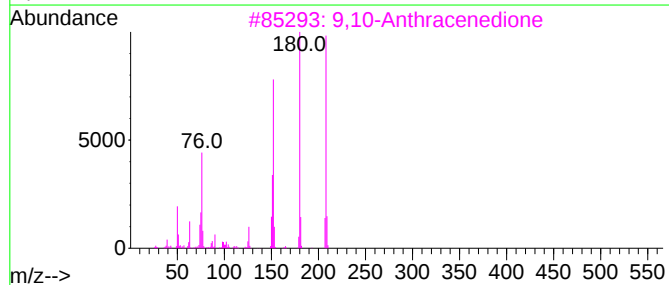
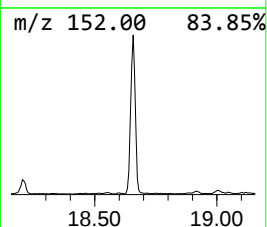
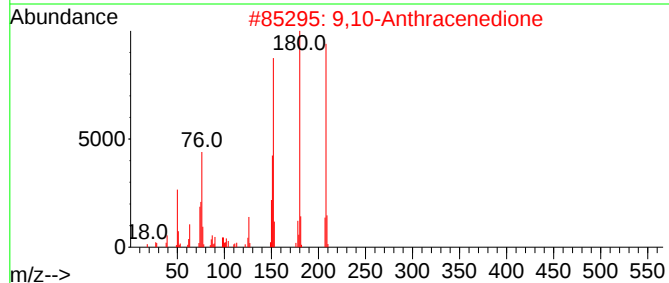
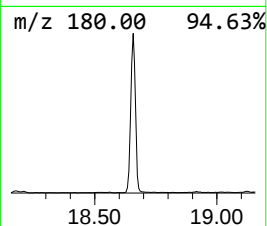
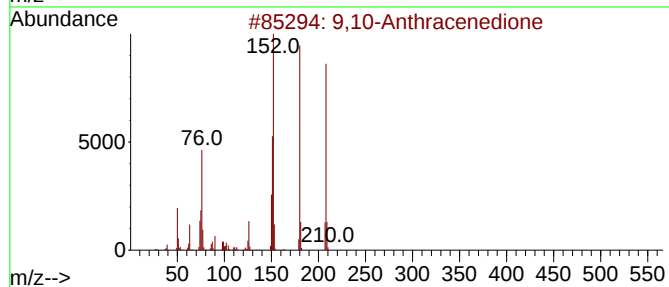
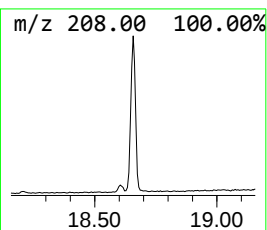
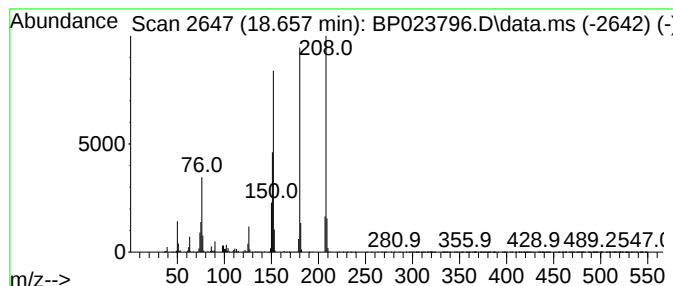
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.657	4.19 ng/ul	1151780	Phenanthrene-d10	17.198

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	95
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
Operator : RC/JU
Sample : Q1190-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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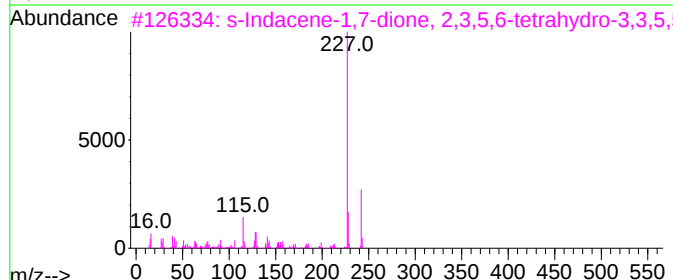
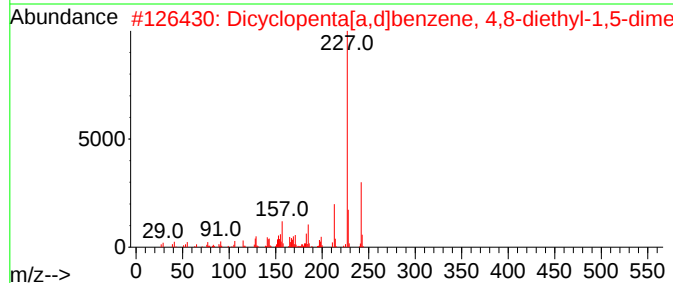
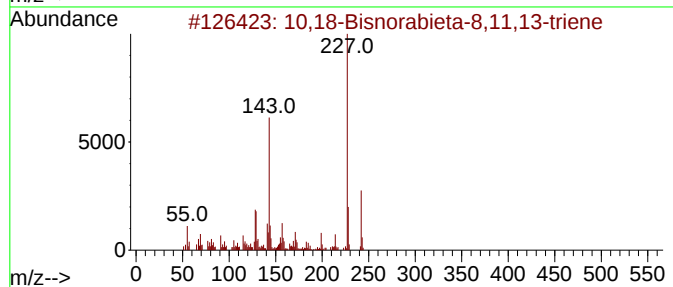
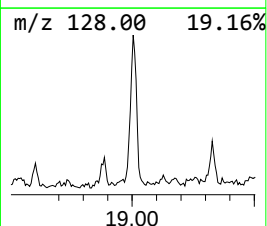
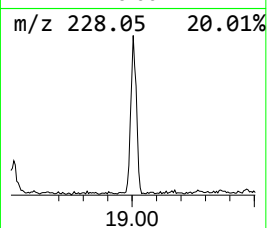
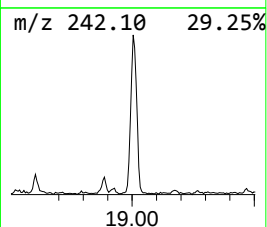
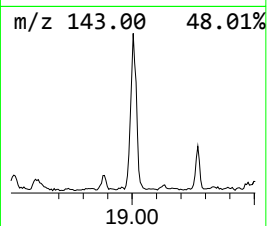
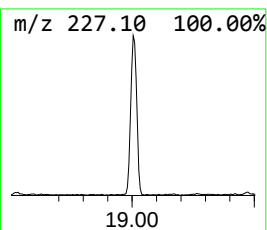
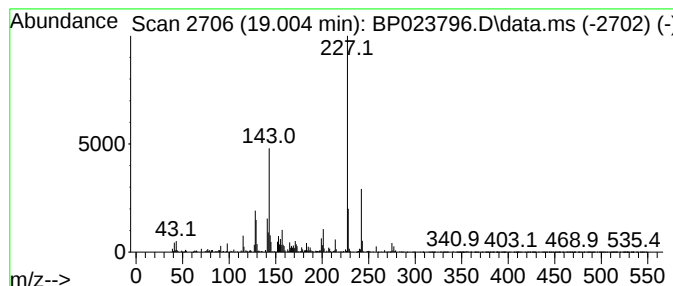
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 10,18-Bisnorabieta-8,11,13-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.004	2.31 ng/ul	635080	Phenanthrene-d10	17.198

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		10,18-Bisnorabieta-8,11,13-triene	242	C18H26	032624-67-2	93
2		Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	60
3		s-Indacene-1,7-dione, 2,3,5,6-te...	242	C16H18O2	055591-17-8	49
4		7-Diethylamino-2-oxo-2H-chromene...	242	C14H14N2O2	332411-59-3	49
5		Chlorquinaldol	227	C10H7Cl12NO	000072-80-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
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Sample : Q1190-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

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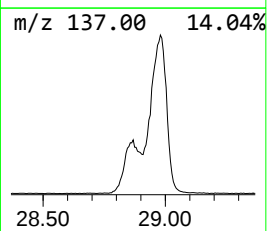
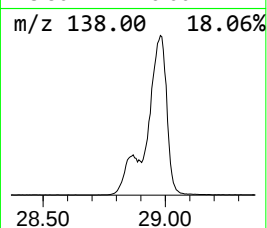
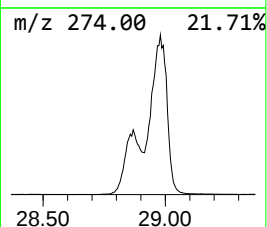
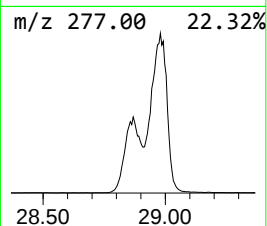
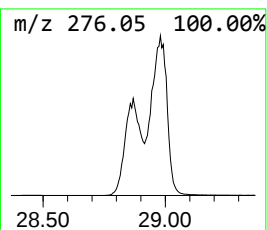
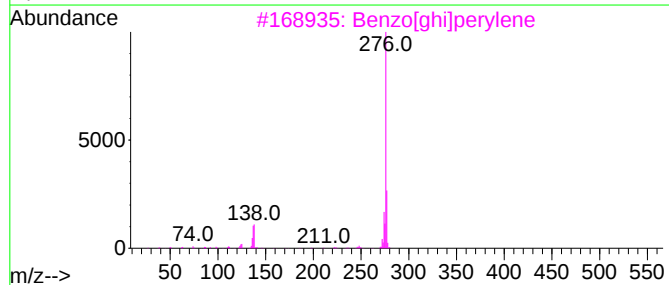
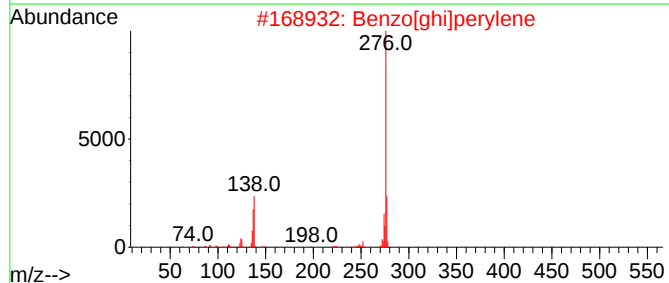
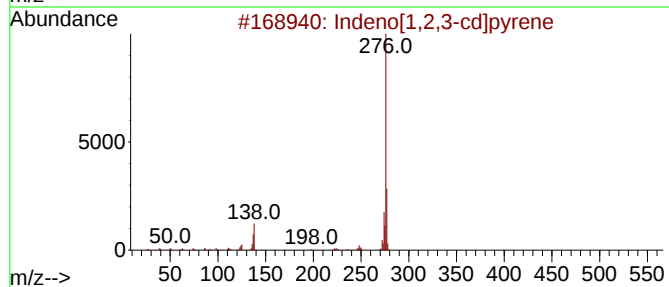
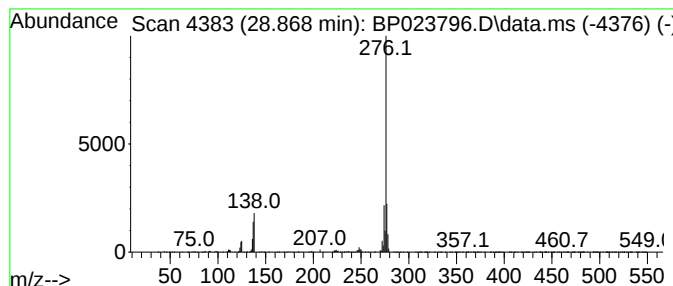
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Dibenzo[def,mno]chrysene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.868	7.85 ng/ul	2463040	Perylene-d12	25.039

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	98
2		Benzo[ghi]perylene	276	C22H12	000191-24-2	97
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	94
4		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	93
5		Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013025\
Data File : BP023796.D
Acq On : 30 Jan 2025 12:29
Operator : RC/JU
Sample : Q1190-19
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.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
Benzene, 1,4-di...	7.669	52.9	ng/ul	5864420	1	7.775	2215410	20.0
Benzene, 1,2-di...	8.134	176.1	ng/ul	19509700	1	7.775	2215410	20.0
Dodecanoic acid...	14.528	2.4	ng/ul	530226	3	14.398	4519590	20.0
Benzophenone	15.775	9.8	ng/ul	2206090	3	14.398	4519590	20.0
Hexadecanoic ac...	17.916	2.7	ng/ul	738891	4	17.198	5492200	20.0
n-Hexadecanoic ...	18.139	3.0	ng/ul	810214	4	17.198	5492200	20.0
9,10-Anthracene...	18.657	4.2	ng/ul	1151780	4	17.198	5492200	20.0
10,18-Bisnorabi...	19.004	2.3	ng/ul	635080	4	17.198	5492200	20.0
Dibenzo[def,mno...	28.868	7.8	ng/ul	2463040	6	25.039	6277950	20.0