

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
 Data File : BP016659.D
 Acq On : 19 Aug 2023 02:37
 Operator : MA/JU
 Sample : 04049-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH8T3

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-BP081423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016659.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.346	6	10	15	rBV	190899	283608	2.44%	0.252%
2	3.387	15	17	31	rVB	66689	126841	1.09%	0.113%
3	3.670	61	65	74	rVB3	27198	54704	0.47%	0.049%
4	3.828	88	92	115	rVB	634528	1005456	8.65%	0.892%
5	4.034	122	127	136	rVB	44769	70390	0.61%	0.062%
6	4.146	142	146	153	rVB	19382	32602	0.28%	0.029%
7	4.552	209	215	222	rVB7	17896	41526	0.36%	0.037%
8	4.675	230	236	251	rVB	729653	1126060	9.69%	0.999%
9	5.399	351	359	365	rBV7	24472	48185	0.41%	0.043%
10	5.875	435	440	445	rVB	108248	160898	1.38%	0.143%
11	5.922	445	448	453	rVB2	28019	41115	0.35%	0.036%
12	5.993	455	460	466	rVB2	17851	32224	0.28%	0.029%
13	6.228	496	500	505	rBV5	19114	36349	0.31%	0.032%
14	6.299	507	512	518	rVB	54408	90932	0.78%	0.081%
15	6.687	576	578	582	rVV4	18403	28718	0.25%	0.025%
16	6.799	590	597	602	rVB5	22151	42460	0.37%	0.038%
17	6.987	624	629	635	rBV5	37237	70787	0.61%	0.063%
18	7.134	648	654	658	rBV2	81287	125847	1.08%	0.112%
19	7.187	658	663	672	rVB	442660	663514	5.71%	0.589%
20	7.387	690	697	707	rVB	1542959	2439938	20.99%	2.165%
21	7.481	709	713	720	rVB8	26190	45771	0.39%	0.041%
22	7.569	721	728	739	rBV	1432136	2373529	20.42%	2.106%
23	7.746	750	758	763	rBV4	31790	81015	0.70%	0.072%
24	7.916	779	787	793	rBV	96037	174276	1.50%	0.155%
25	7.975	793	797	802	rVV	34347	53789	0.46%	0.048%
26	8.052	802	810	824	rVV	1171120	1985582	17.08%	1.762%
27	8.163	824	829	835	rVB4	18633	33069	0.28%	0.029%
28	8.316	853	855	860	rVV2	17256	29582	0.25%	0.026%
29	8.522	880	890	894	rBV3	22481	60800	0.52%	0.054%
30	8.563	894	897	907	rVB9	19837	44243	0.38%	0.039%
31	8.752	922	929	938	rVB	1227946	1947117	16.75%	1.728%
32	9.128	985	993	999	rVB3	128338	237206	2.04%	0.210%
33	9.246	1006	1013	1019	rBV	1840099	2993869	25.76%	2.656%
34	9.387	1025	1037	1041	rBV	579974	974386	8.38%	0.865%
35	9.422	1041	1043	1048	rVV2	84203	121713	1.05%	0.108%
36	9.481	1048	1053	1059	rVB2	43797	86033	0.74%	0.076%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
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37	9.599	1064	1073	1079	rBV	130512	218665	1.88%	0.194%
38	9.716	1089	1093	1096	rBV	35152	53414	0.46%	0.047%
39	9.840	1108	1114	1120	rVB	274998	427649	3.68%	0.379%
40	9.899	1120	1124	1128	rBV2	18035	35863	0.31%	0.032%
41	9.963	1128	1135	1141	rBV	1375616	2223260	19.13%	1.973%
42	10.122	1156	1162	1169	rVB	386466	598723	5.15%	0.531%
43	10.228	1169	1180	1186	rBV	110192	197238	1.70%	0.175%
44	10.416	1205	1212	1216	rBV2	42325	73040	0.63%	0.065%
45	10.487	1216	1224	1231	rVV	2066216	3422782	29.45%	3.037%
46	10.546	1231	1234	1243	rVB6	42497	85891	0.74%	0.076%
47	10.775	1269	1273	1279	rVB	47063	77117	0.66%	0.068%
48	10.881	1279	1291	1300	rBV	1823943	3671388	31.59%	3.257%
49	10.963	1300	1305	1310	rVV	175257	304477	2.62%	0.270%
50	11.034	1310	1317	1332	rVV	1999563	3472407	29.88%	3.081%
51	11.146	1332	1336	1342	rVV	66976	116912	1.01%	0.104%
52	11.216	1343	1348	1360	rVB3	48282	97780	0.84%	0.087%
53	11.363	1369	1373	1380	rVV3	13911	29971	0.26%	0.027%
54	11.628	1412	1418	1421	rBV5	29033	52811	0.45%	0.047%
55	11.657	1421	1423	1427	rVV5	26456	41535	0.36%	0.037%
56	11.698	1427	1430	1434	rVV3	20033	32135	0.28%	0.029%
57	11.751	1434	1439	1445	rVV6	20827	43906	0.38%	0.039%
58	11.922	1458	1468	1472	rBV6	13938	46424	0.40%	0.041%
59	11.993	1472	1480	1484	rVV	47832	98709	0.85%	0.088%
60	12.028	1484	1486	1494	rVB9	15227	28128	0.24%	0.025%
61	12.140	1500	1505	1510	rBV2	44966	72362	0.62%	0.064%
62	12.228	1515	1520	1525	rBV3	26085	45379	0.39%	0.040%
63	12.428	1550	1554	1556	rBV	47180	62365	0.54%	0.055%
64	12.457	1556	1559	1563	rVV	38915	57856	0.50%	0.051%
65	12.522	1563	1570	1578	rVB3	136702	271889	2.34%	0.241%
66	12.663	1587	1594	1599	rBV3	31633	60813	0.52%	0.054%
67	12.757	1606	1610	1616	rVB5	29570	46068	0.40%	0.041%
68	12.869	1625	1629	1634	rVB4	25810	40781	0.35%	0.036%
69	12.922	1634	1638	1641	rBV3	20365	31583	0.27%	0.028%
70	12.981	1641	1648	1654	rVV5	59320	140016	1.20%	0.124%
71	13.063	1658	1662	1666	rVV6	22681	40862	0.35%	0.036%
72	13.104	1666	1669	1675	rVV7	17131	28455	0.24%	0.025%
73	13.163	1675	1679	1685	rVB	28915	48563	0.42%	0.043%
74	13.340	1702	1709	1718	rBV8	49300	151024	1.30%	0.134%
75	13.469	1727	1731	1736	rVB7	21800	32777	0.28%	0.029%
76	13.545	1740	1744	1748	rVV2	34959	54333	0.47%	0.048%

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 Title : SVOA CALIBRATION

77	13.693	1766	1769	1773	rVB2	24004	36105	0.31%	0.032%
78	13.810	1783	1789	1796	rVB2	126631	214193	1.84%	0.190%
79	13.975	1810	1817	1822	rBV7	35616	68188	0.59%	0.060%
80	14.110	1834	1840	1848	rVV	4332996	5882774	50.61%	5.219%
81	14.357	1878	1882	1883	rVV	66533	74128	0.64%	0.066%
82	14.398	1883	1889	1899	rVB2	4747094	6979036	60.05%	6.192%
83	14.698	1934	1940	1953	rBV	2686308	3995462	34.38%	3.545%
84	14.904	1969	1975	1979	rBV	353902	560786	4.82%	0.498%
85	15.239	2028	2032	2039	rVB4	63908	120340	1.04%	0.107%
86	15.375	2050	2055	2058	rBV	410808	606534	5.22%	0.538%
87	15.404	2058	2060	2071	rVB	364444	528822	4.55%	0.469%
88	15.692	2103	2109	2118	rBV	6090974	8743330	75.23%	7.757%
89	15.810	2123	2129	2145	rVV	1966645	2816761	24.23%	2.499%
90	15.928	2147	2149	2158	rVB4	33993	53966	0.46%	0.048%
91	17.469	2404	2411	2421	rBV	3375962	4822505	41.49%	4.279%
92	17.569	2421	2428	2439	rBV	7147928	10063088	86.58%	8.928%
93	18.304	2549	2553	2563	rBV	458651	772187	6.64%	0.685%
94	19.222	2706	2709	2712	rVB2	73998	76498	0.66%	0.068%
95	19.369	2731	2734	2739	rBV3	97141	125862	1.08%	0.112%
96	19.539	2759	2763	2772	rBV	294940	470233	4.05%	0.417%
97	19.798	2802	2807	2816	rBV	8498168	11622818	100.00%	10.312%
98	21.563	3102	3107	3115	rBV2	3817761	4901491	42.17%	4.349%
99	23.980	3510	3518	3533	rVB	4770369	10241530	88.12%	9.087%
100	24.151	3540	3547	3559	rVB	2042482	4532972	39.00%	4.022%

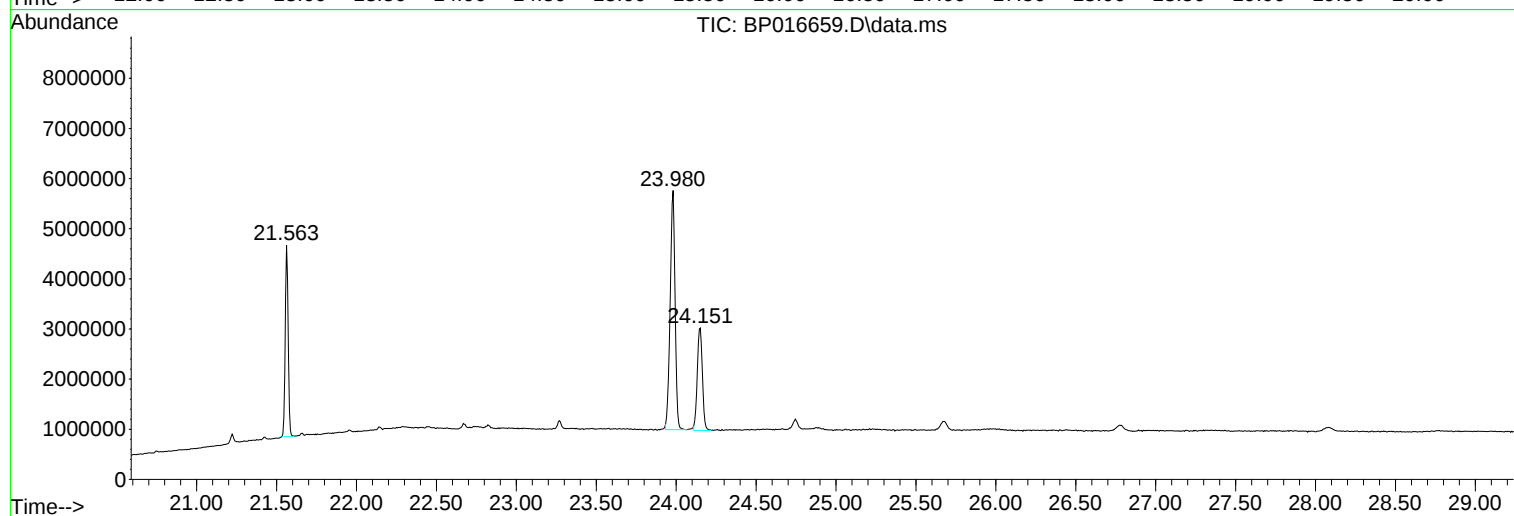
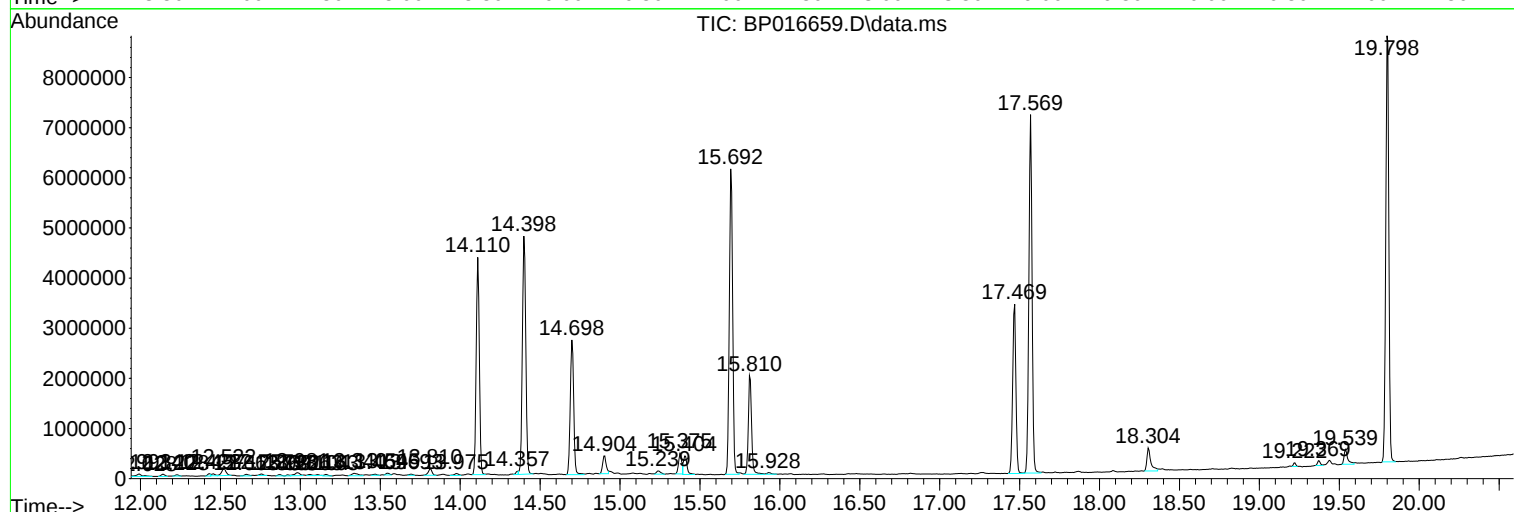
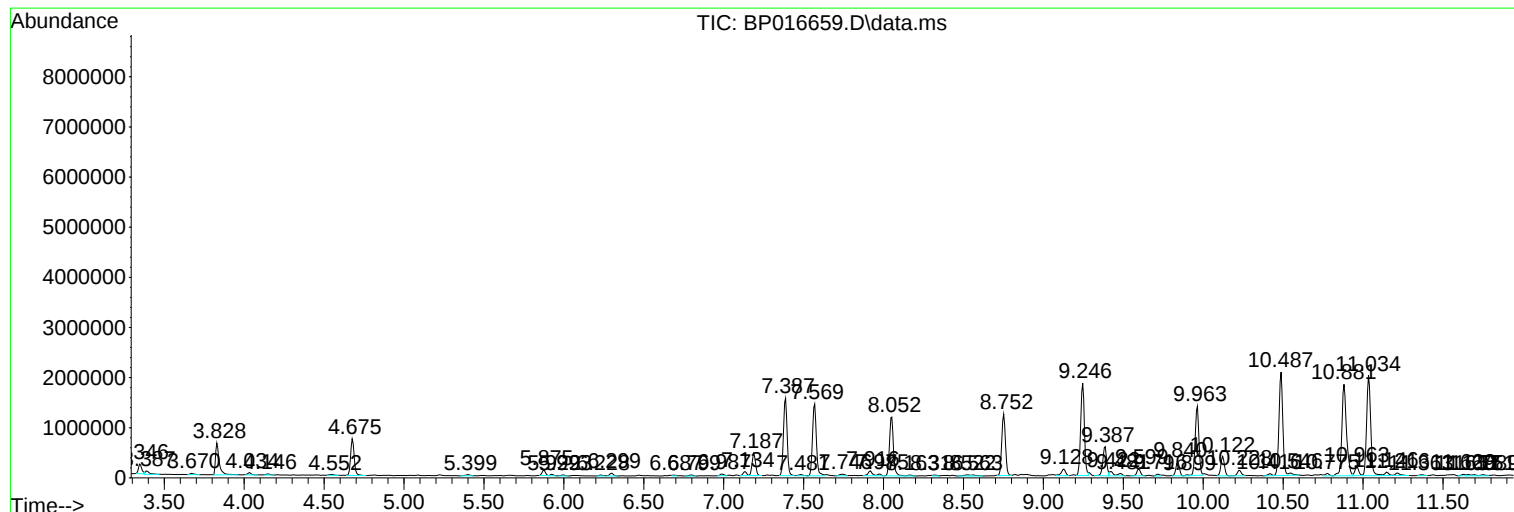
Sum of corrected areas: 112709094

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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



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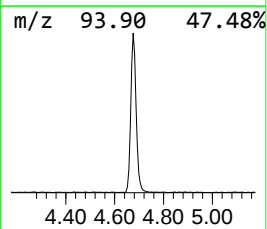
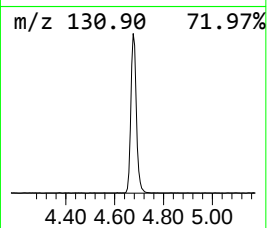
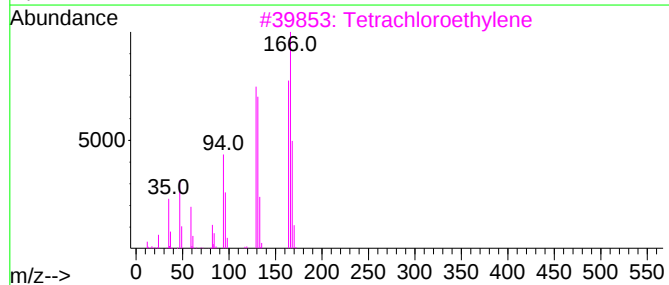
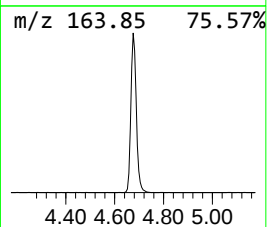
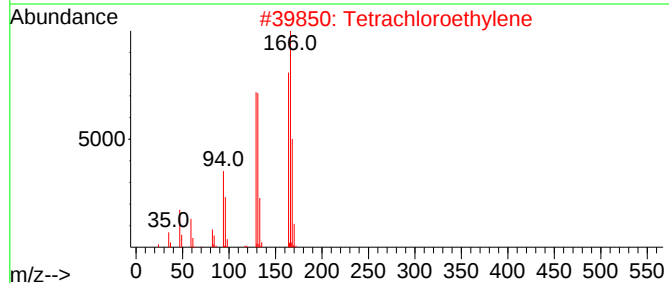
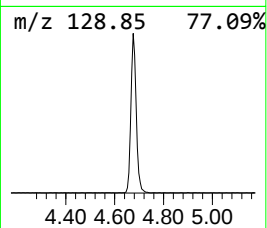
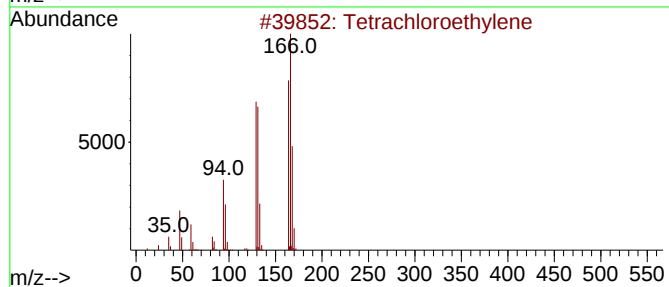
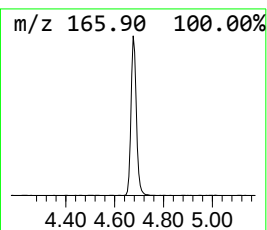
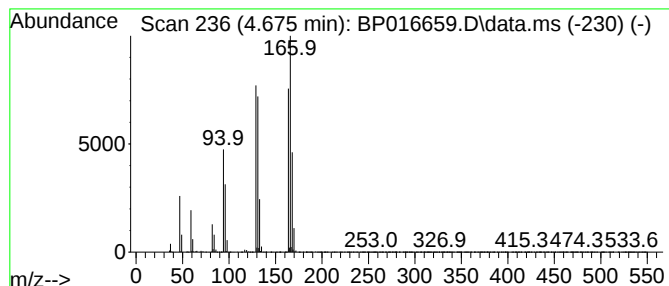
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.675	11.34 ng/ul	1126060	1,4-Dichlorobenzene-d4	8.052

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	25



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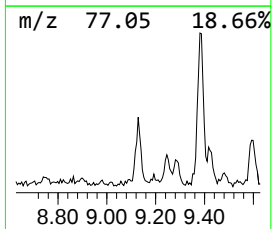
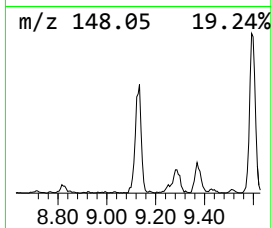
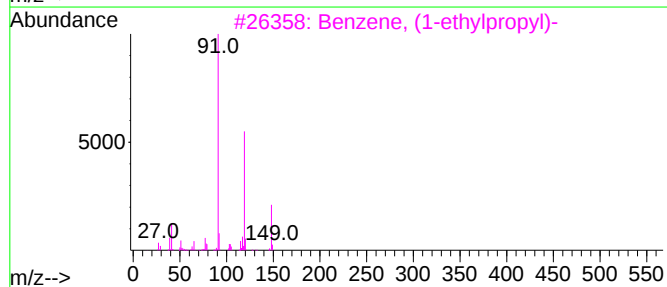
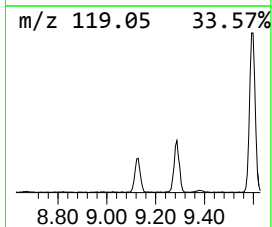
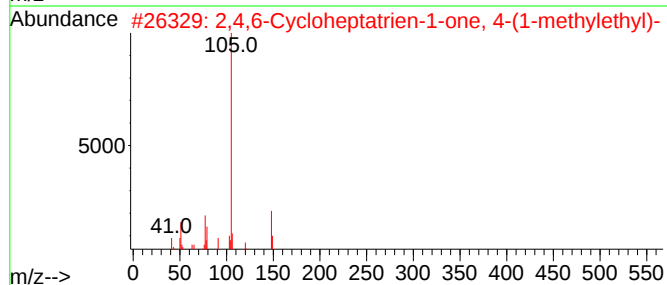
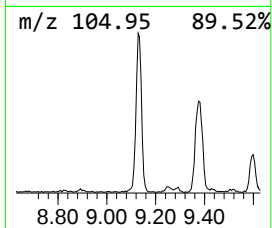
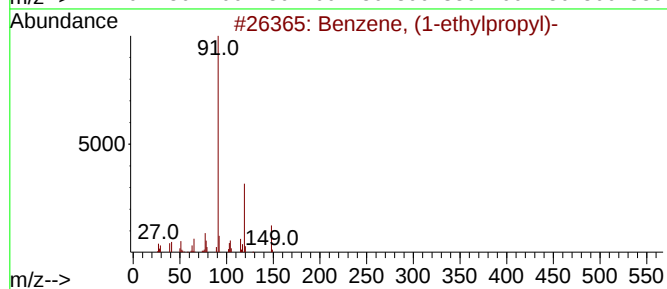
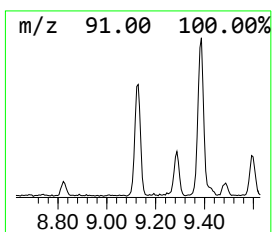
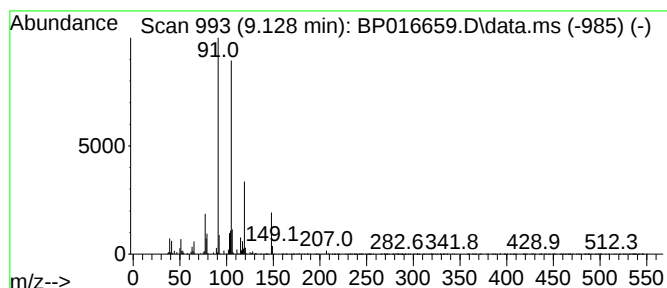
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, (1-ethylpropyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.128	2.39 ng/ul	237206	1,4-Dichlorobenzene-d4	8.052

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethylpropyl)-	148	C11H16	001196-58-3	70
2			2,4,6-Cycloheptatrien-1-one, 4-(...	148	C10H12O	013656-81-0	47
3			Benzene, (1-ethylpropyl)-	148	C11H16	001196-58-3	47
4			Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	47
5			Hexane, 3,4-diphenyl-	238	C18H22	005789-31-1	38



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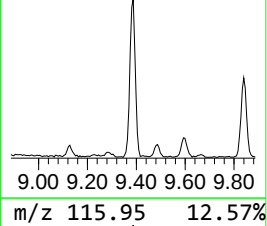
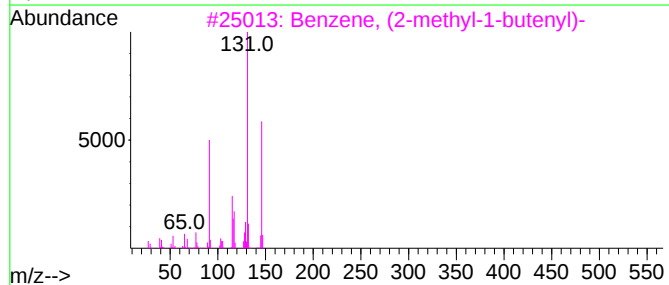
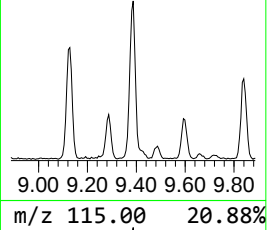
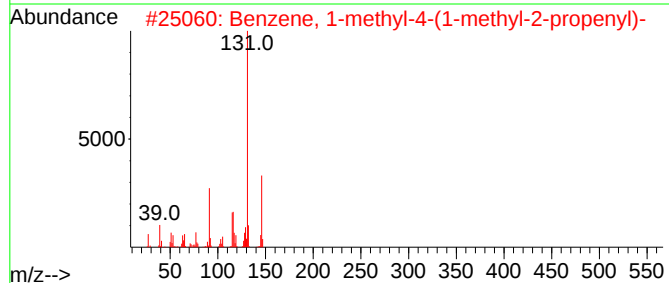
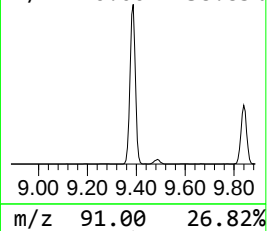
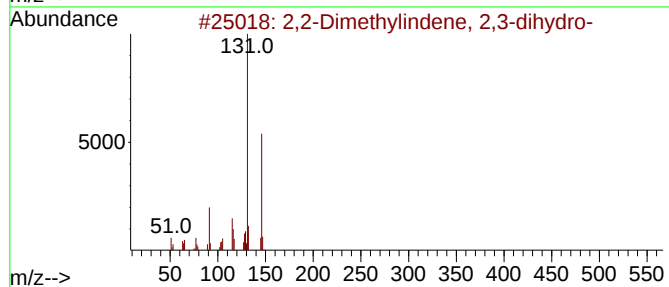
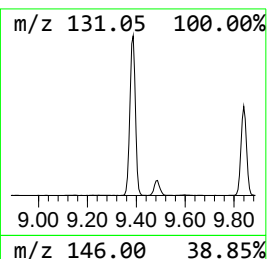
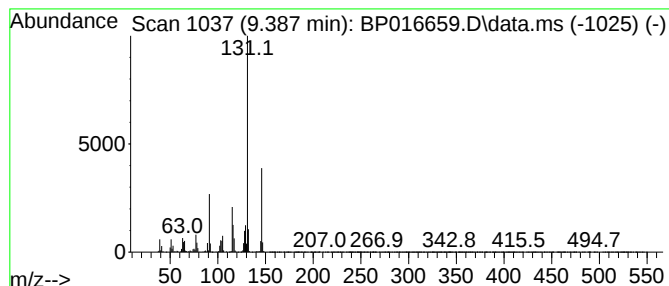
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,2-Dimethylindene, 2,3-dih... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.387	9.81 ng/ul	974386	1,4-Dichlorobenzene-d4	8.052

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	97		
2	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	94		
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94		
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	92		
5	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
Data File : BP016659.D
Acq On : 19 Aug 2023 02:37
Operator : MA/JU
Sample : 04049-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH8T3

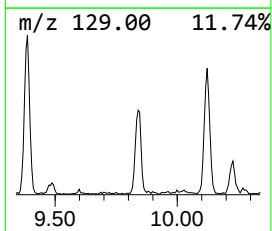
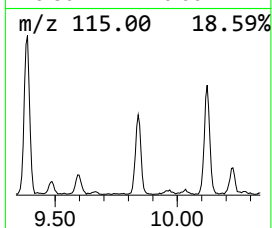
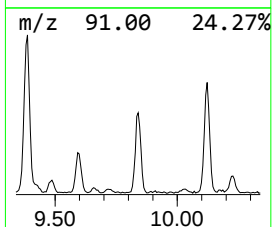
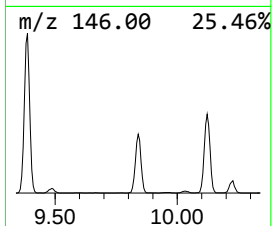
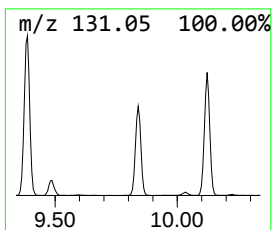
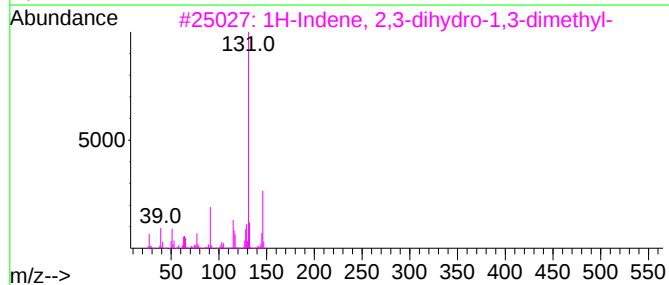
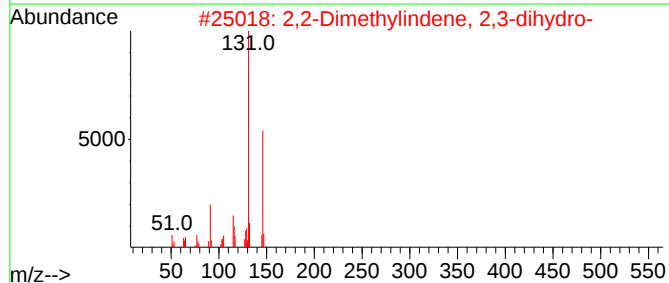
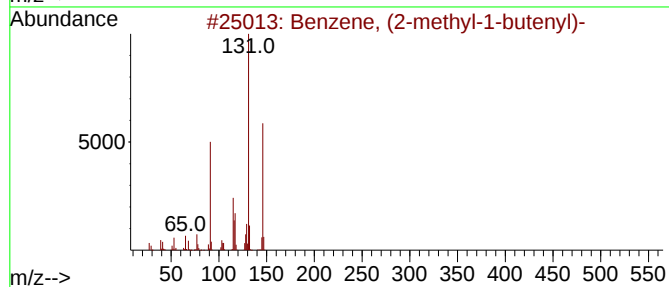
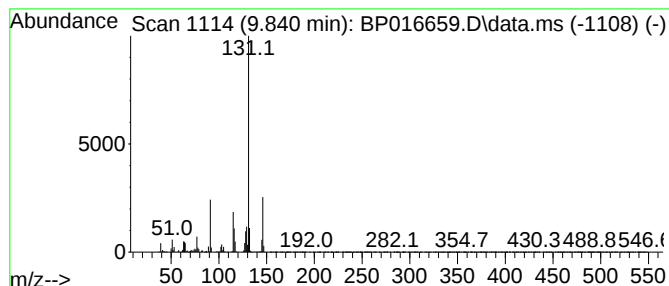
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, (2-methyl-1-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.840	2.33 ng/ul	427649	Naphthalene-d8	10.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
5			Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
Data File : BP016659.D
Acq On : 19 Aug 2023 02:37
Operator : MA/JU
Sample : 04049-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH8T3

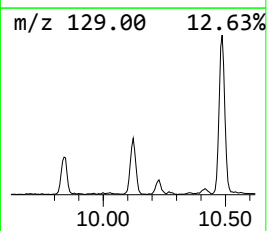
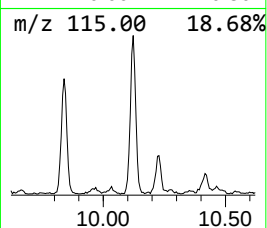
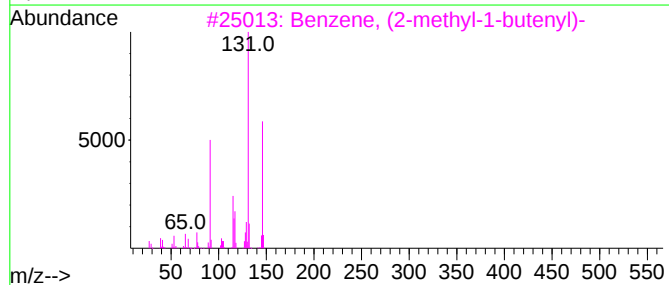
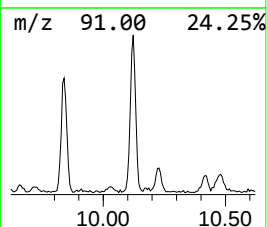
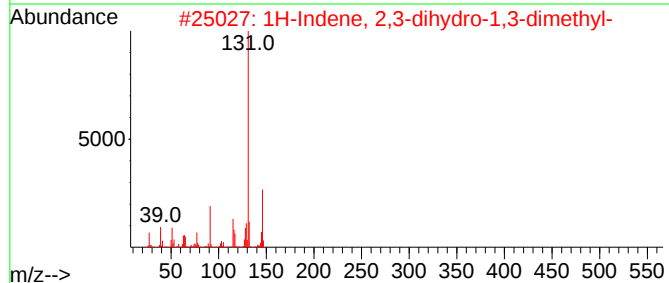
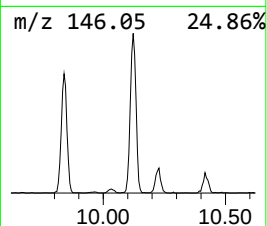
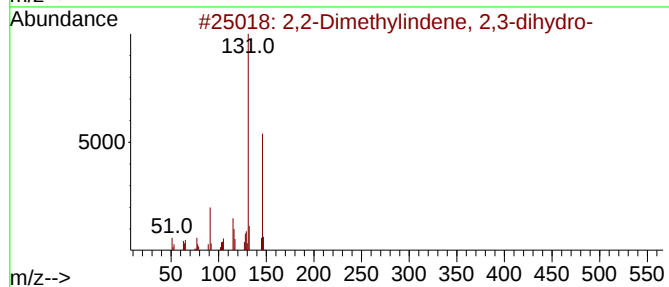
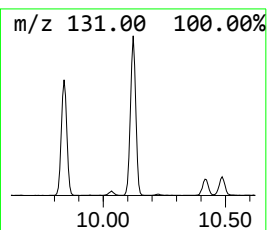
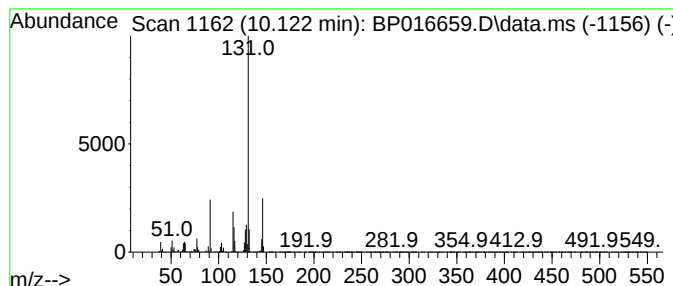
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.122	3.26 ng/ul	598723	Naphthalene-d8	10.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94		
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94		
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94		
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	91		
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
Data File : BP016659.D
Acq On : 19 Aug 2023 02:37
Operator : MA/JU
Sample : 04049-01
Misc :
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Instrument :
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ClientSampleId :
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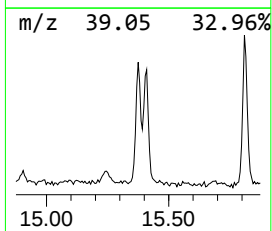
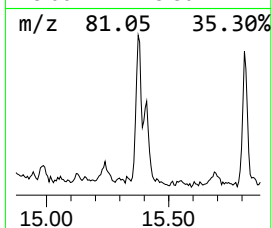
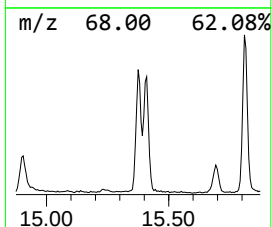
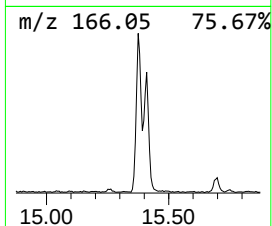
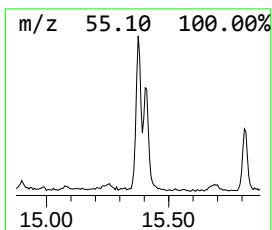
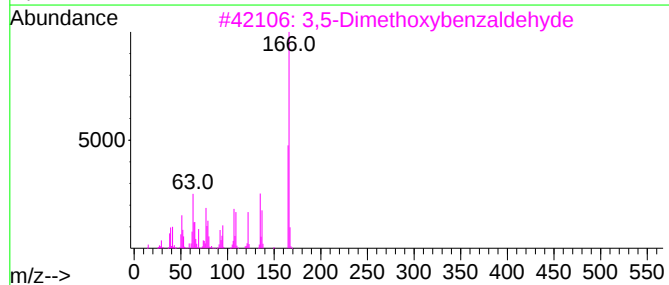
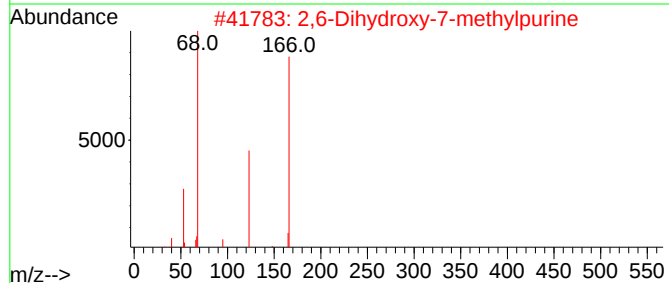
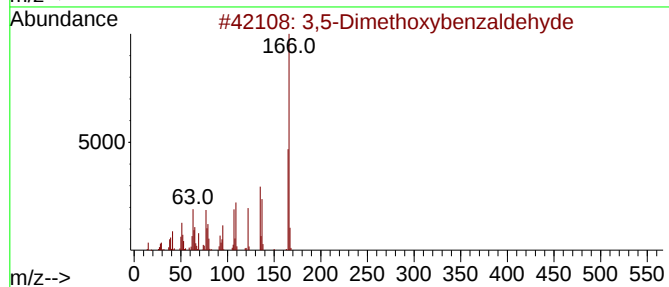
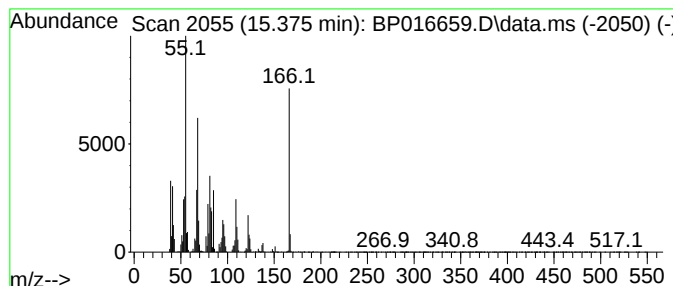
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.375	3.04 ng/ul	606534	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Dimethoxybenzaldehyde	166	C9H10O3	007311-34-4	43
2			2,6-Dihydroxy-7-methylpurine	166	C6H6N4O2	000552-62-5	38
3			3,5-Dimethoxybenzaldehyde	166	C9H10O3	007311-34-4	38
4			1,E-11,Z-13-Pentadecatriene	206	C15H26	080625-34-9	27
5			Thiazolo[3,2-a]pyridinium, 3-car...	211	C9H9N03S	023003-41-0	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
Data File : BP016659.D
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Operator : MA/JU
Sample : 04049-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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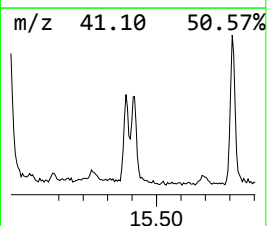
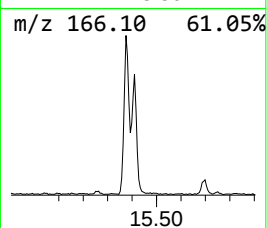
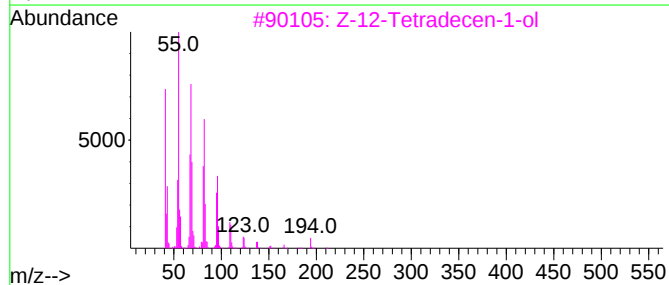
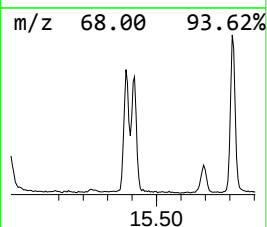
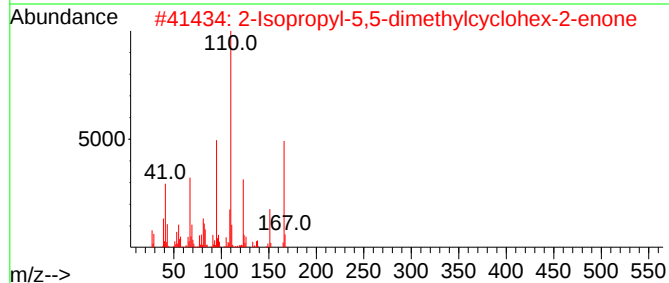
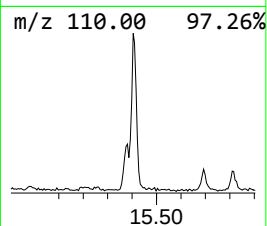
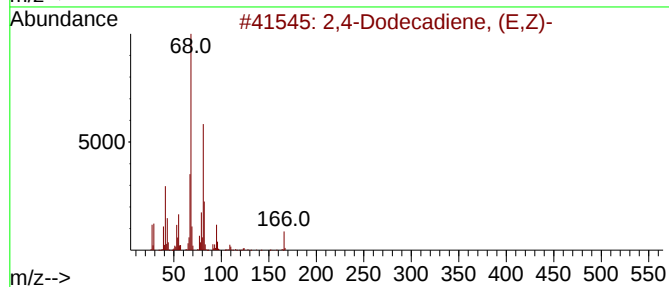
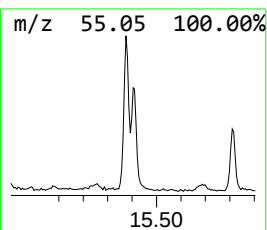
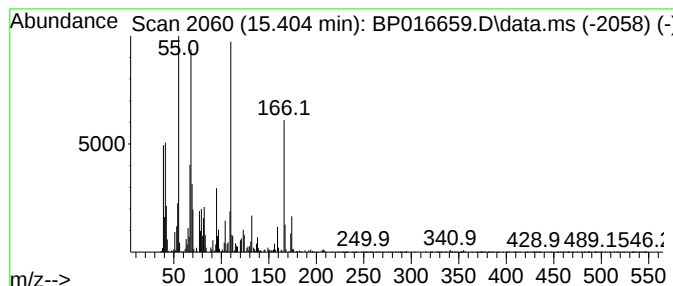
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,4-Dodecadiene, (E,Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.404	2.65 ng/ul	528822	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dodecadiene, (E,Z)-		166	C12H22	074685-27-1	53	
2	2-Isopropyl-5,5-dimethylcyclohex...		166	C11H18O	1000191-19-3	46	
3	Z-12-Tetradecen-1-ol		212	C14H28O	1000130-84-0	45	
4	2-Oxaspiro[5.5]undecane-3,5-dion...		346	C20H26O5	1000277-11-3	38	
5	11-Tetradecen-1-ol, (E)-		212	C14H28O	035153-18-5	38	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
Data File : BP016659.D
Acq On : 19 Aug 2023 02:37
Operator : MA/JU
Sample : 04049-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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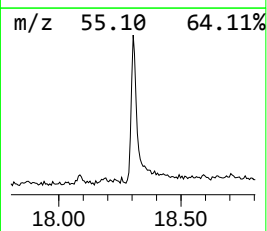
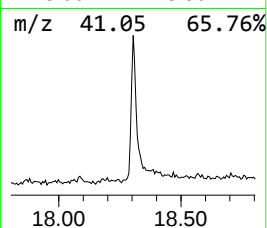
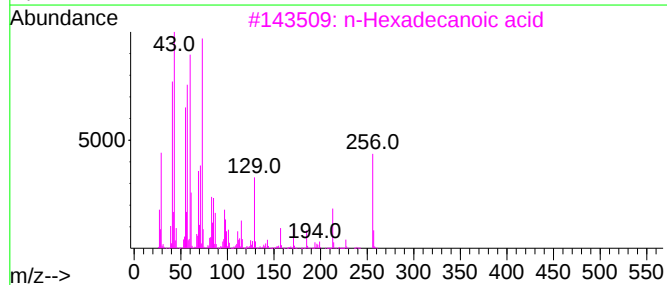
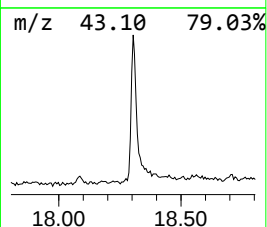
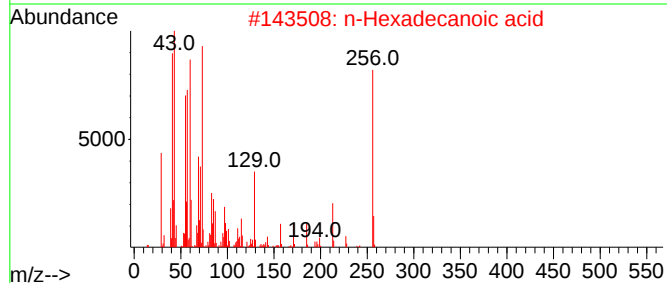
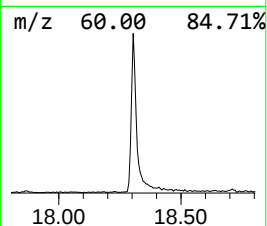
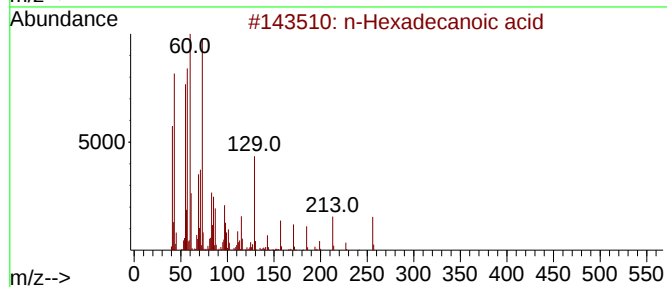
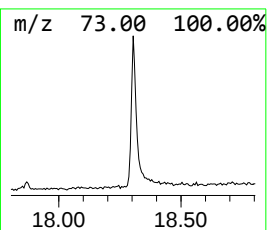
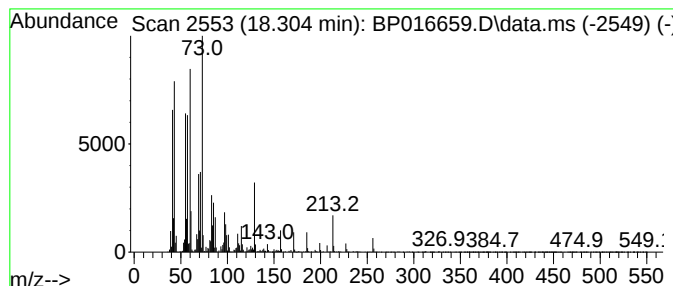
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.304	3.20 ng/ul	772187	Phenanthrene-d10	17.469

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	93
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081823\
Data File : BP016659.D
Acq On : 19 Aug 2023 02:37
Operator : MA/JU
Sample : 04049-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Benzene, (1-eth...	9.128	2.4	ng/ul	237206	1	8.052	1985580	20.0
2,2-Dimethylind...	9.387	9.8	ng/ul	974386	1	8.052	1985580	20.0
Benzene, (2-met...	9.840	2.3	ng/ul	427649	2	10.881	3671390	20.0
1H-Indene, 2,3-...	10.122	3.3	ng/ul	598723	2	10.881	3671390	20.0
unknown-01	15.375	3.0	ng/ul	606534	3	14.698	3995460	20.0
2,4-Dodecadiene...	15.404	2.6	ng/ul	528822	3	14.698	3995460	20.0
n-Hexadecanoic ...	18.304	3.2	ng/ul	772187	4	17.469	4822510	20.0