

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
 Data File : BP017652.D
 Acq On : 11 Oct 2023 18:02
 Operator : MA/JU
 Sample : 04744-06RX
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH583RX

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-BP100923.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017652.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	7	12	30	rVB	1324679	1781860	46.32%	5.046%
2	3.281	30	33	42	rVB2	35393	50330	1.31%	0.143%
3	3.446	57	61	62	rVB4	11551	12766	0.33%	0.036%
4	3.723	104	108	116	rBV	170747	280440	7.29%	0.794%
5	3.905	136	139	144	rVB	20977	26824	0.70%	0.076%
6	3.981	149	152	156	rBV6	8942	14574	0.38%	0.041%
7	4.446	228	231	237	rVB4	21419	35325	0.92%	0.100%
8	4.558	243	250	262	rVB	1497167	2126858	55.28%	6.023%
9	4.840	289	298	311	rVB	398821	668244	17.37%	1.892%
10	4.940	311	315	322	rVB5	30052	53938	1.40%	0.153%
11	5.070	335	337	341	rVB3	15601	20717	0.54%	0.059%
12	5.658	433	437	438	rBV3	5715	6593	0.17%	0.019%
13	6.081	507	509	511	rVV2	7328	7591	0.20%	0.021%
14	6.111	511	514	515	rVV2	7156	5909	0.15%	0.017%
15	6.152	517	521	528	rVB4	37391	60830	1.58%	0.172%
16	6.205	528	530	533	rBV	5305	5911	0.15%	0.017%
17	6.317	544	549	550	rBV2	6723	8125	0.21%	0.023%
18	6.593	594	596	604	rVB	5114	9927	0.26%	0.028%
19	7.264	704	710	722	rBV	462186	765388	19.89%	2.167%
20	7.399	730	733	735	rBV2	7709	10346	0.27%	0.029%
21	7.452	737	742	749	rVV3	35008	77691	2.02%	0.220%
22	7.564	760	761	764	rVB	9047	6061	0.16%	0.017%
23	7.828	800	806	812	rVB2	43009	74089	1.93%	0.210%
24	7.922	815	822	831	rVV	439656	730738	18.99%	2.069%
25	7.999	831	835	842	rVV3	39824	69322	1.80%	0.196%
26	8.187	863	867	868	rBV	4796	6374	0.17%	0.018%
27	8.293	883	885	889	rVV3	10050	13648	0.35%	0.039%
28	8.411	903	905	906	rBV2	8123	6585	0.17%	0.019%
29	8.422	906	907	912	rVB	8260	7336	0.19%	0.021%
30	8.552	927	929	932	rVB	5903	7088	0.18%	0.020%
31	8.852	979	980	985	rVB	8095	8979	0.23%	0.025%
32	8.905	985	989	990	rBV2	7115	9554	0.25%	0.027%
33	8.922	990	992	997	rVV2	5609	9146	0.24%	0.026%
34	9.134	1021	1028	1037	rBV	530488	933726	24.27%	2.644%
35	9.222	1037	1043	1049	rVV5	14647	40744	1.06%	0.115%
36	9.875	1147	1154	1157	rBV6	15689	31086	0.81%	0.088%

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Integrator: RTE

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Stop Thrs : 0

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.MA.M
 Title : SVOA CALIBRATION

37	10.240	1214	1216	1219	rBV	4772	5867	0.15%	0.017%
38	10.399	1238	1243	1251	rBV4	45917	114767	2.98%	0.325%
39	10.587	1273	1275	1281	rVB4	4652	6856	0.18%	0.019%
40	10.699	1291	1294	1295	rBV	10037	7844	0.20%	0.022%

41	10.758	1298	1304	1313	rVB	610948	1010234	26.26%	2.861%
42	10.887	1320	1326	1328	rBV2	4765	6923	0.18%	0.020%
43	11.299	1395	1396	1402	rVB4	4630	6584	0.17%	0.019%
44	11.793	1477	1480	1481	rBV3	7257	6236	0.16%	0.018%
45	11.834	1481	1487	1492	rVV4	16379	31761	0.83%	0.090%

46	12.363	1571	1577	1580	rBV6	8708	15339	0.40%	0.043%
47	12.452	1590	1592	1595	rBV	6217	5713	0.15%	0.016%
48	12.687	1630	1632	1637	rVB2	5136	7306	0.19%	0.021%
49	12.799	1646	1651	1652	rBV4	5281	6558	0.17%	0.019%
50	13.293	1733	1735	1737	rBV2	5411	5807	0.15%	0.016%

51	13.375	1744	1749	1750	rBV	6171	6482	0.17%	0.018%
52	13.451	1760	1762	1763	rVV2	7243	6767	0.18%	0.019%
53	13.463	1763	1764	1768	rVB4	8758	6559	0.17%	0.019%
54	13.622	1787	1791	1794	rVV3	3686	5947	0.15%	0.017%
55	13.799	1819	1821	1823	rVV3	5706	6711	0.17%	0.019%

56	13.840	1823	1828	1831	rVV3	18033	32549	0.85%	0.092%
57	13.863	1831	1832	1835	rVV2	8729	8209	0.21%	0.023%
58	13.887	1835	1836	1839	rVV2	7017	7390	0.19%	0.021%
59	13.951	1844	1847	1849	rVB2	4966	6169	0.16%	0.017%
60	14.034	1854	1861	1869	rBV	1436864	1971054	51.23%	5.581%

61	14.316	1902	1909	1919	rBV	771413	1120366	29.12%	3.173%
62	14.616	1954	1960	1968	rVV	965676	1472496	38.28%	4.170%
63	14.687	1970	1972	1975	rVV3	7141	8918	0.23%	0.025%
64	14.716	1975	1977	1980	rVV2	7188	8555	0.22%	0.024%
65	14.746	1980	1982	1989	rVV5	9328	13033	0.34%	0.037%

66	14.857	1999	2001	2005	rBV3	4224	5744	0.15%	0.016%
67	14.981	2020	2022	2027	rVV	4784	8243	0.21%	0.023%
68	15.028	2028	2030	2032	rVV3	6148	6159	0.16%	0.017%
69	15.322	2076	2080	2082	rBV	6884	9325	0.24%	0.026%
70	15.369	2085	2088	2090	rBV2	6719	6796	0.18%	0.019%

71	15.416	2090	2096	2097	rBV3	9627	11562	0.30%	0.033%
72	15.622	2125	2131	2141	rBV	2299624	3167003	82.32%	8.968%
73	16.134	2215	2218	2222	rVV4	15934	20898	0.54%	0.059%
74	16.204	2227	2230	2232	rVV3	5136	6498	0.17%	0.018%
75	16.245	2234	2237	2238	rBV2	4700	6283	0.16%	0.018%

76	16.345	2248	2254	2266	rBV2	184661	328382	8.54%	0.930%
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Integrator: RTE

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 Title : SVOA CALIBRATION

77	16.769	2322	2326	2330	rVB4	4991	6945	0.18%	0.020%
78	16.934	2352	2354	2356	rBV	10468	6613	0.17%	0.019%
79	17.051	2367	2374	2379	rBV2	63900	111307	2.89%	0.315%
80	17.122	2384	2386	2388	rVB2	8888	6863	0.18%	0.019%
81	17.140	2388	2389	2393	rBV	12139	10226	0.27%	0.029%
82	17.292	2412	2415	2422	rVB	11396	15874	0.41%	0.045%
83	17.410	2429	2435	2443	rBV	1382506	1967042	51.13%	5.570%
84	17.510	2446	2452	2466	rVV	1675227	2414659	62.76%	6.838%
85	17.675	2477	2480	2482	rBV2	7298	8282	0.22%	0.023%
86	17.816	2499	2504	2506	rBV4	6511	10620	0.28%	0.030%
87	18.051	2543	2544	2547	rVB3	8999	6230	0.16%	0.018%
88	18.239	2571	2576	2583	rBV4	33469	52544	1.37%	0.149%
89	18.586	2633	2635	2636	rBV	8704	7061	0.18%	0.020%
90	18.657	2644	2647	2649	rBV3	6414	8373	0.22%	0.024%
91	18.798	2666	2671	2679	rBV2	129035	208096	5.41%	0.589%
92	19.028	2708	2710	2712	rBV2	8437	6405	0.17%	0.018%
93	19.334	2759	2762	2768	rVB4	33057	45048	1.17%	0.128%
94	19.404	2771	2774	2780	rVB	39699	54320	1.41%	0.154%
95	19.639	2812	2814	2816	rBV3	14546	13702	0.36%	0.039%
96	19.763	2829	2835	2844	rBV	2757005	3847146	100.00%	10.894%
97	20.963	3035	3039	3046	rVB	356073	463707	12.05%	1.313%
98	21.551	3134	3139	3145	rBV	1818676	2412466	62.71%	6.831%
99	23.951	3540	3547	3557	rVB2	1644744	3577720	93.00%	10.131%
100	24.122	3570	3576	3588	rVB	1232477	2572288	66.86%	7.284%

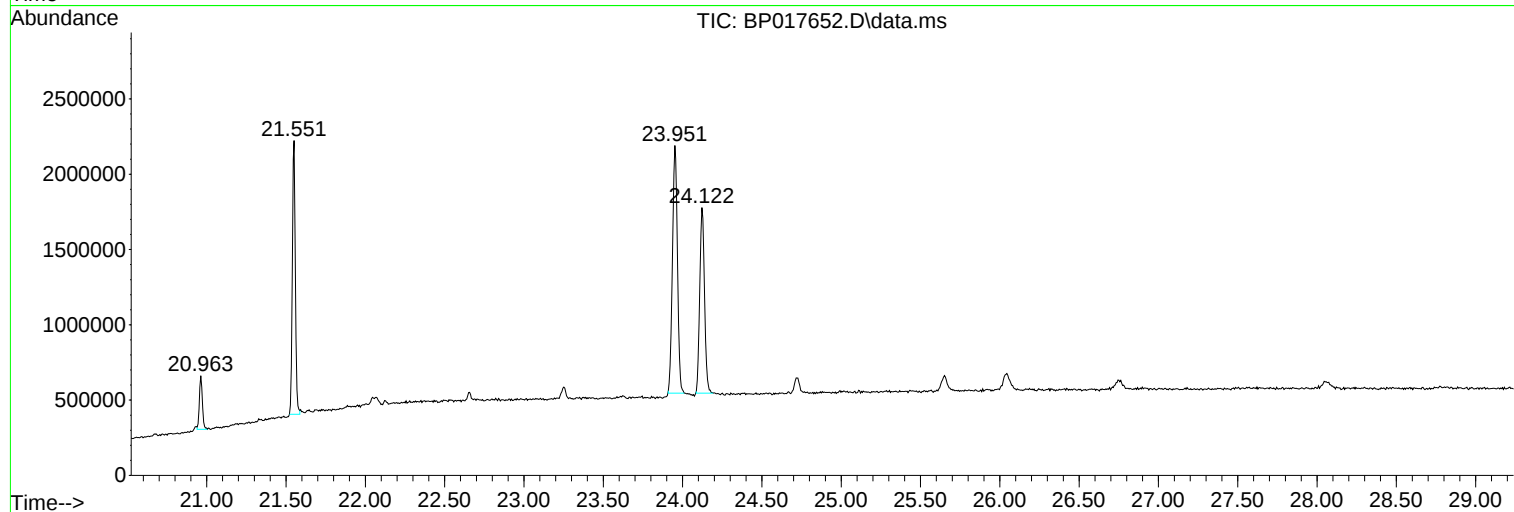
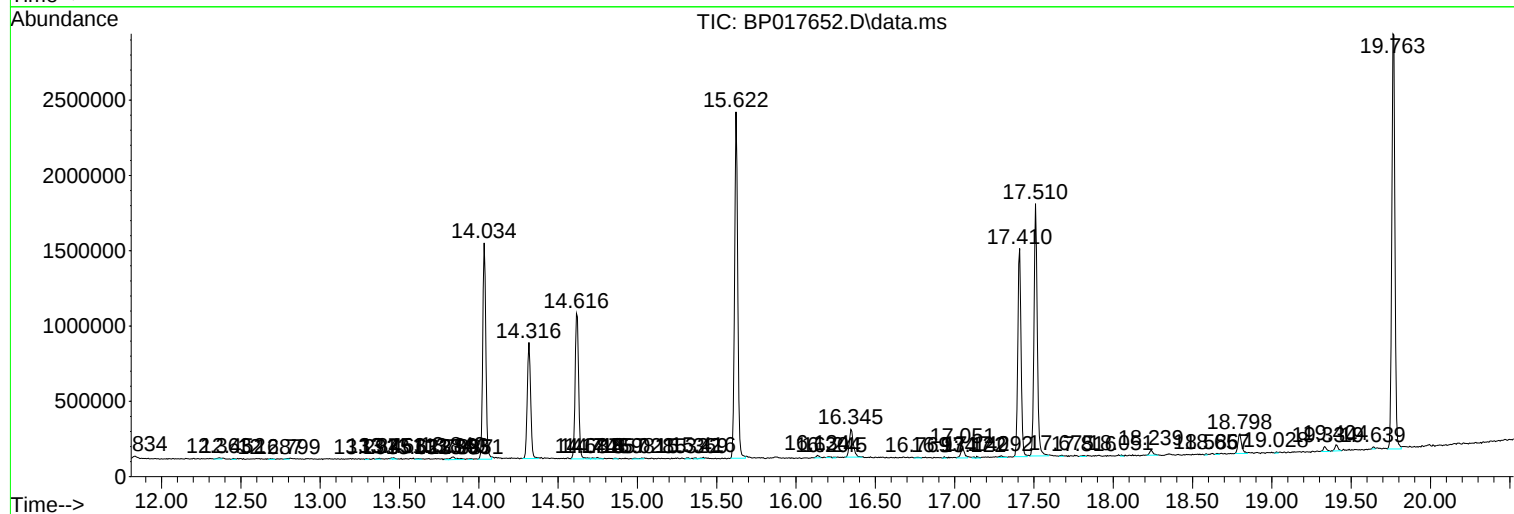
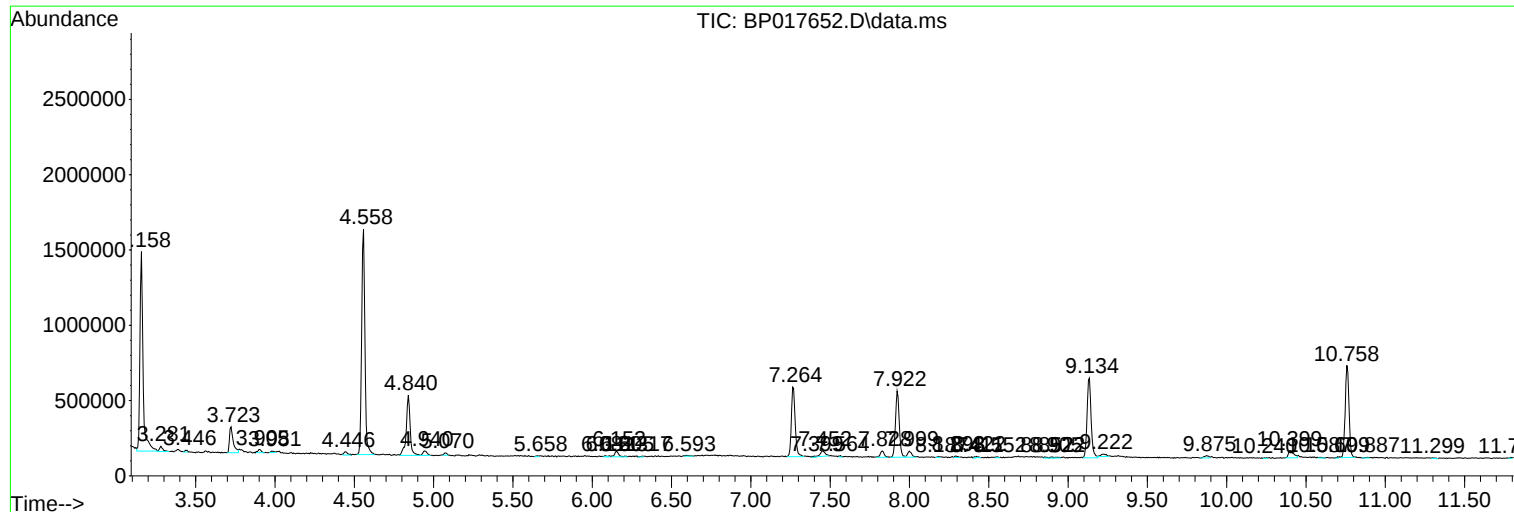
Sum of corrected areas: 35314093

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH583RX

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.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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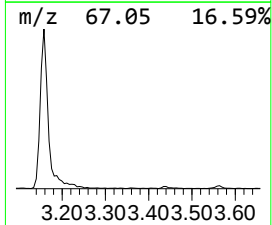
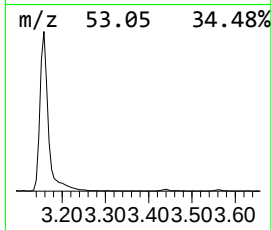
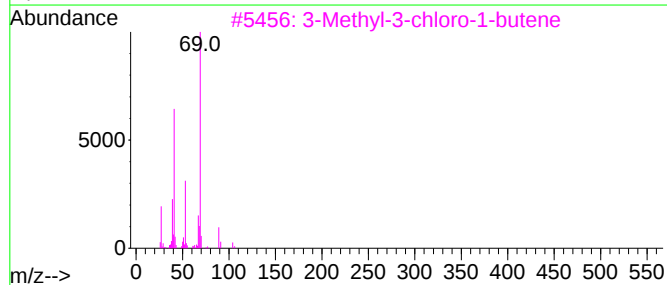
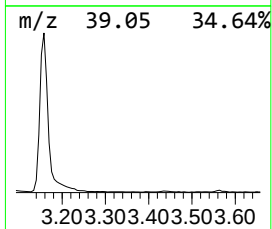
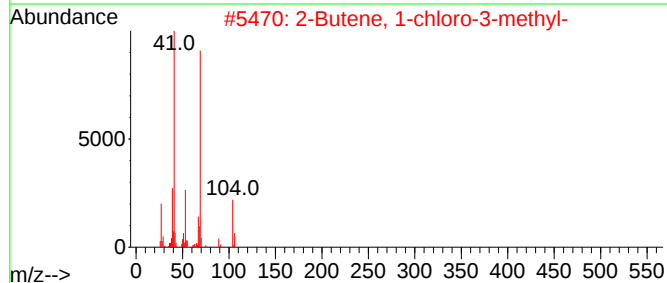
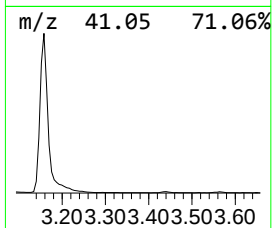
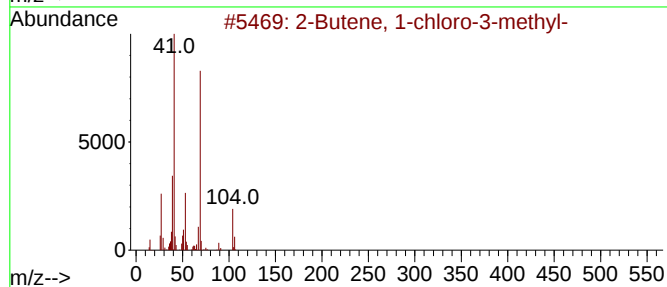
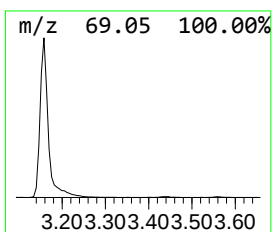
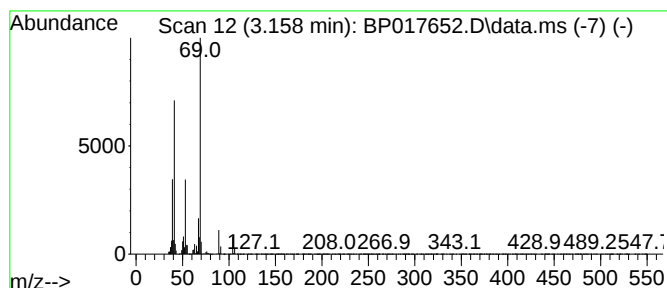
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Butene, 1-chloro-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.158	48.77 ng/ul	1781860	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 1-chloro-3-methyl-		104	C5H9Cl	000503-60-6	80	
2	2-Butene, 1-chloro-3-methyl-		104	C5H9Cl	000503-60-6	80	
3	3-Methyl-3-chloro-1-butene		104	C5H9Cl	002190-48-9	64	
4	2-Butene, 1-chloro-3-methyl-		104	C5H9Cl	000503-60-6	59	
5	2-Butene, 1-chloro-2-methyl-		104	C5H9Cl	013417-43-1	56	



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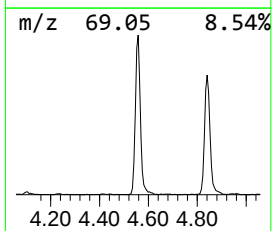
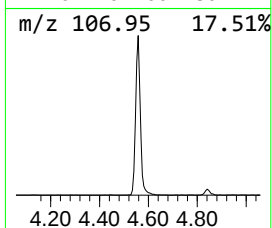
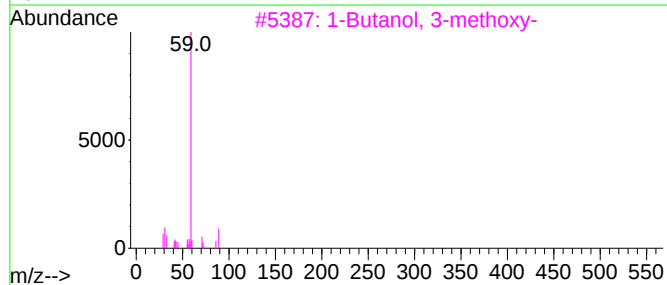
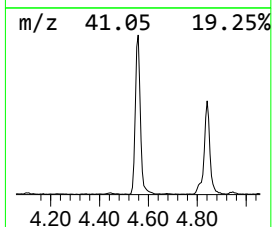
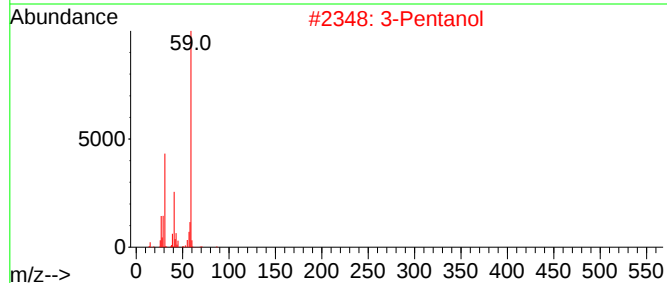
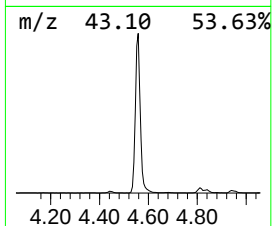
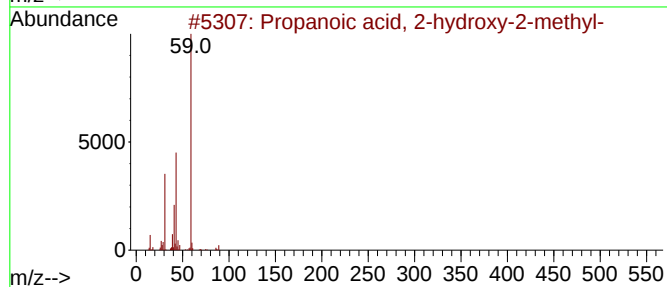
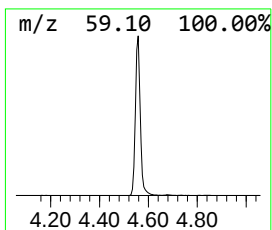
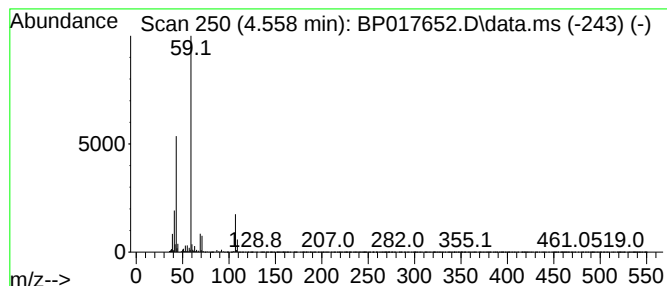
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propanoic acid, 2-hydroxy-2... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.558	58.21 ng/ul	2126860	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	53
2			3-Pentanol	88	C5H12O	000584-02-1	42
3			1-Butanol, 3-methoxy-	104	C5H12O2	002517-43-3	38
4			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	36
5			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36



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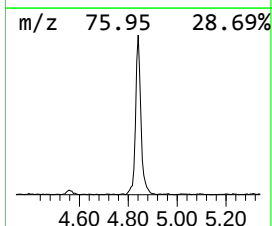
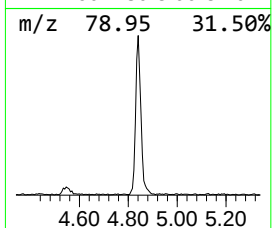
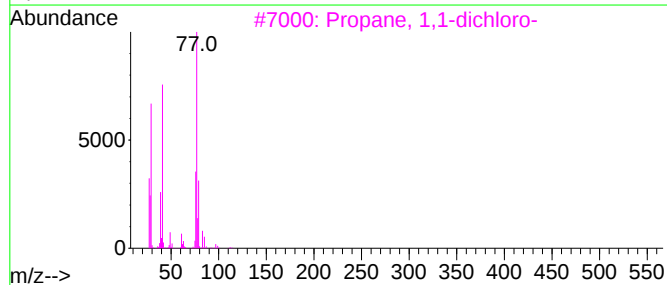
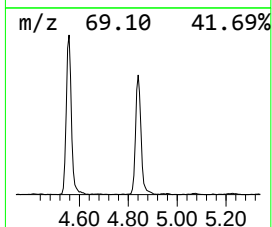
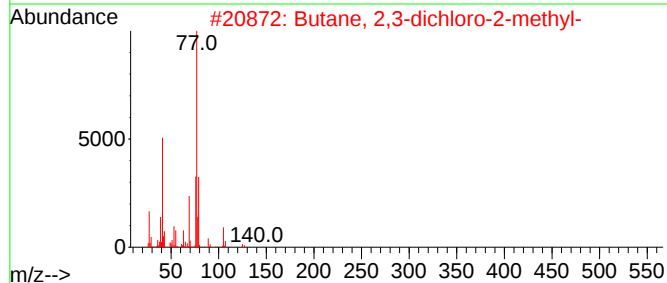
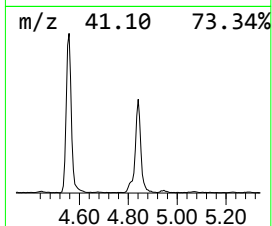
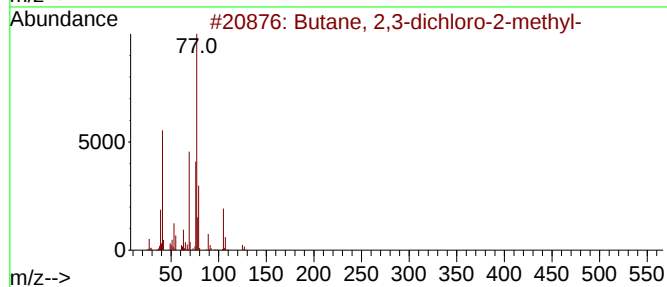
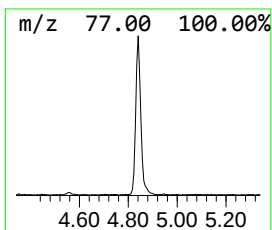
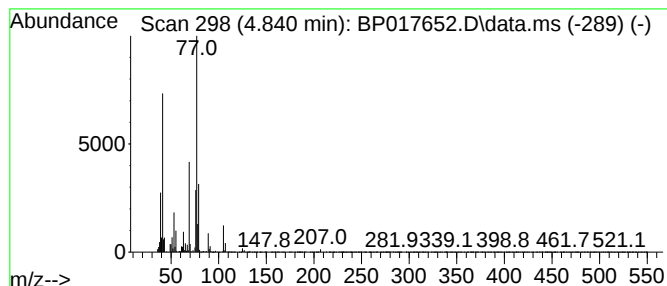
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Butane, 2,3-dichloro-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.840	18.29 ng/ul	668244	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	83
2			Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	83
3			Propane, 1,1-dichloro-	112	C3H6Cl2	000078-99-9	38
4			Propane, 1,1,2-trichloro-2-methyl-	160	C4H7Cl3	029559-52-2	33
5			Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	28



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
Data File : BP017652.D
Acq On : 11 Oct 2023 18:02
Operator : MA/JU
Sample : 04744-06RX
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH583RX

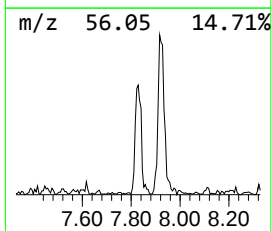
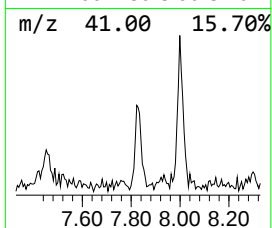
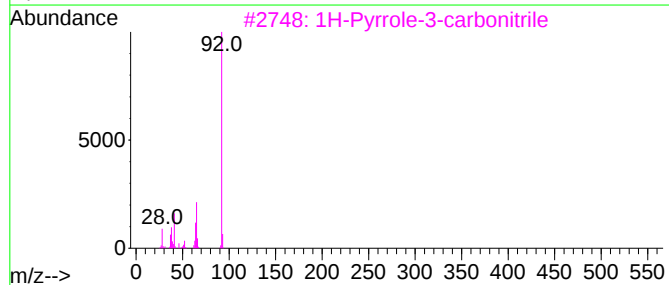
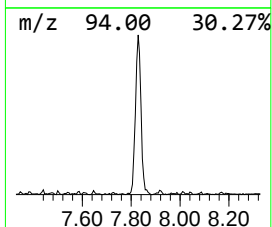
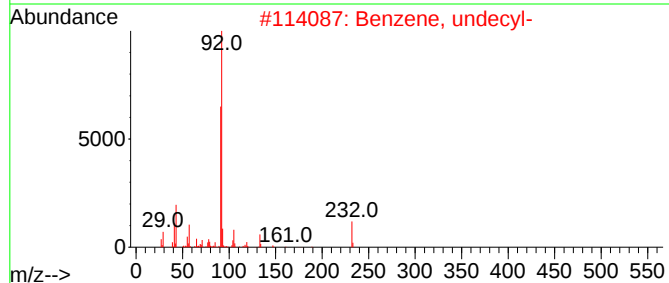
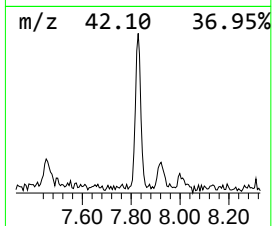
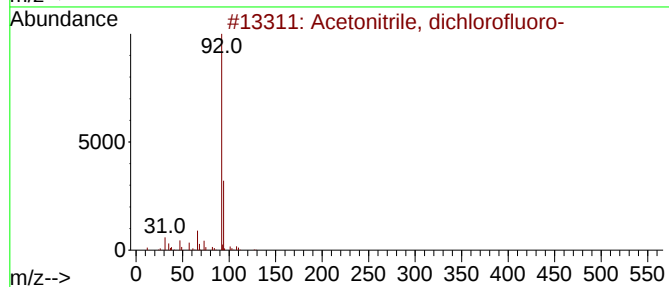
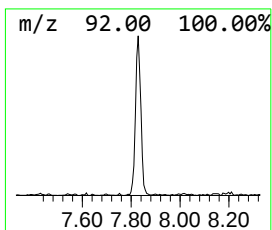
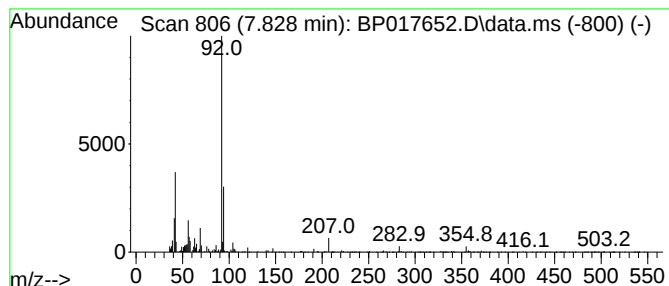
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	2.03 ng/ul	74089	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetonitrile, dichlorofluoro-	127	C2Cl2FN	000353-82-2	28
2			Benzene, undecyl-	232	C17H28	006742-54-7	23
3			1H-Pyrrole-3-carbonitrile	92	C5H4N2	007126-38-7	9
4			Benzeneethanol, .alpha.-(2-methy...	178	C12H18O	007779-78-4	9
5			1,1,3,3,5,5-Hexamethyl-5-(pyridi...	331	C12H25NO4Si3	1000334-15-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
Data File : BP017652.D
Acq On : 11 Oct 2023 18:02
Operator : MA/JU
Sample : 04744-06RX
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH583RX

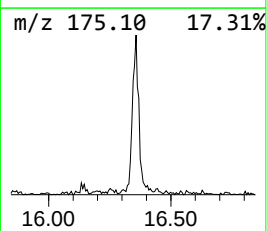
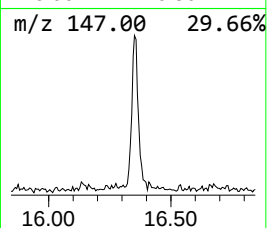
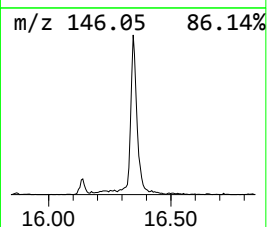
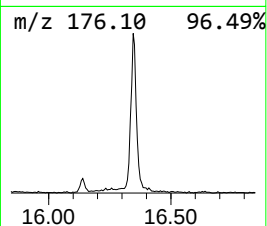
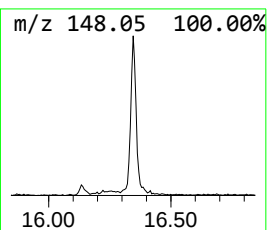
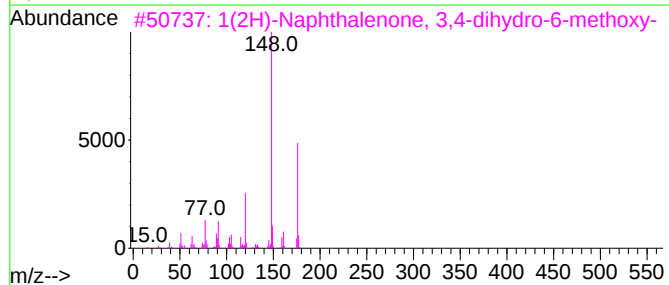
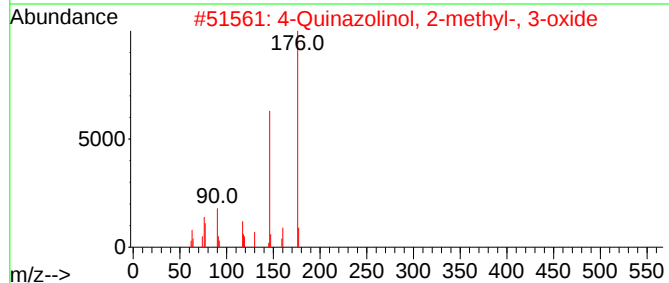
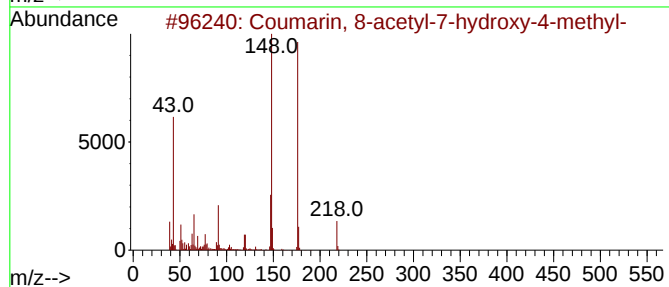
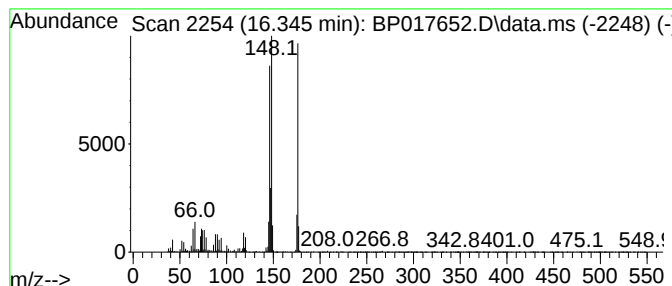
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Coumarin, 8-acetyl-7-hydrox... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.345	3.34 ng/ul	328382	Phenanthrene-d10	17.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Coumarin, 8-acetyl-7-hydroxy-4-m...	218	C12H10O4	002555-29-5	50
2			4-Quinazolinol, 2-methyl-, 3-oxide	176	C9H8N2O2	054518-07-9	47
3			1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47
4			1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	47
5			Furan, 3-bromo-	146	C4H3BrO	022037-28-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
Data File : BP017652.D
Acq On : 11 Oct 2023 18:02
Operator : MA/JU
Sample : 04744-06RX
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH583RX

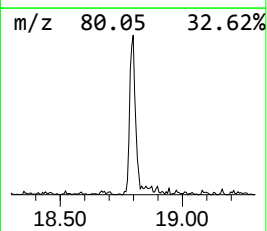
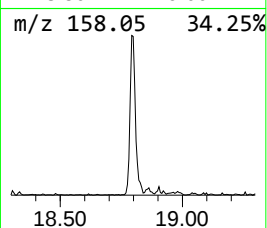
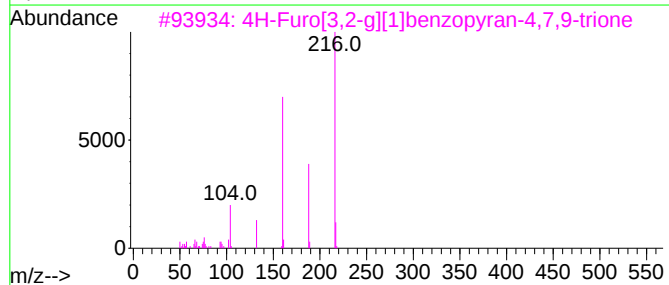
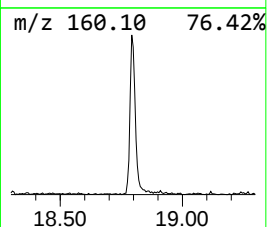
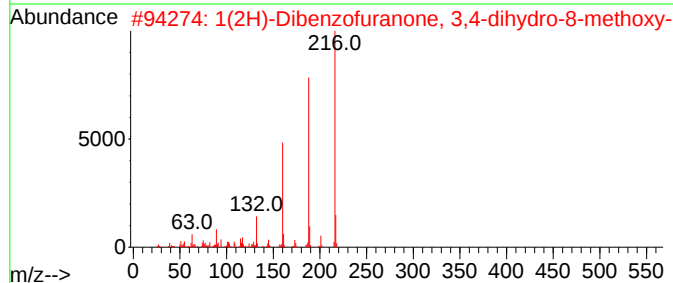
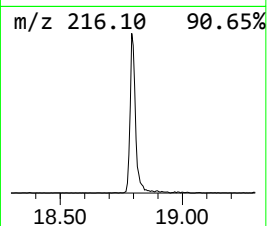
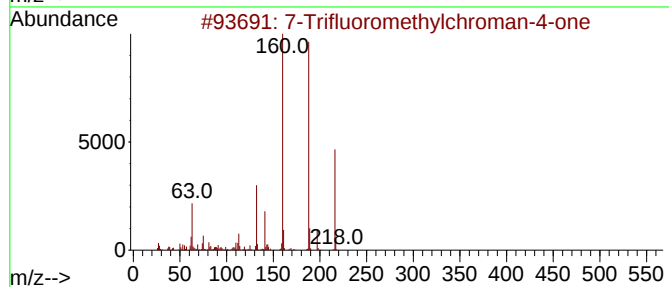
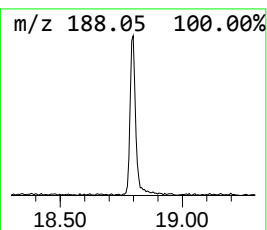
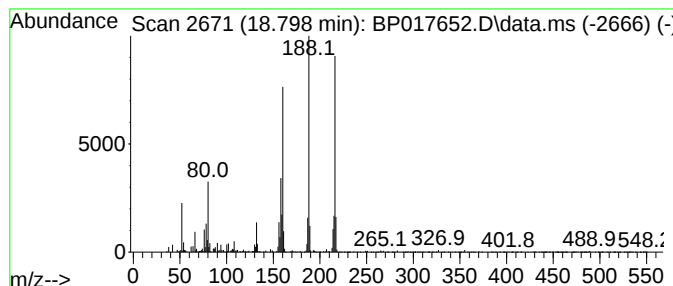
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.798	2.12 ng/ul	208096	Phenanthrene-d10	17.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Trifluoromethylchroman-4-one	216	C10H7F3O2	111141-02-7	49		
2	1(2H)-Dibenzofuranone, 3,4-dihyd...	216	C13H12O3	095338-40-2	43		
3	4H-Furo[3,2-g][1]benzopyran-4,7,...	216	C11H4O5	000483-36-3	35		
4	1,2-Dihydro-2,7-naphthyridin-1-o...	160	C9H8N2O	1000501-27-1	11		
5	4-Fluoro-6-methyl-2-phenylpyrimi...	188	C11H9FN2	051421-92-2	11		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
Data File : BP017652.D
Acq On : 11 Oct 2023 18:02
Operator : MA/JU
Sample : 04744-06RX
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH583RX

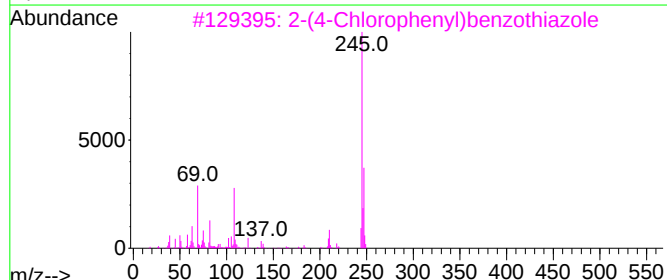
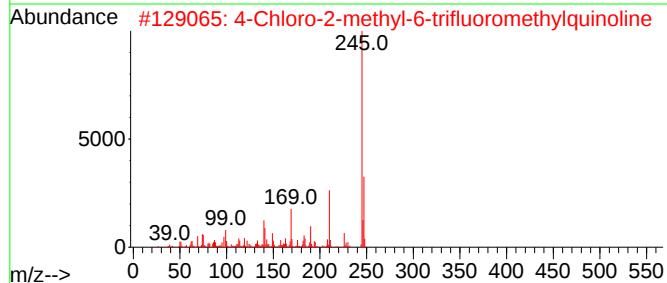
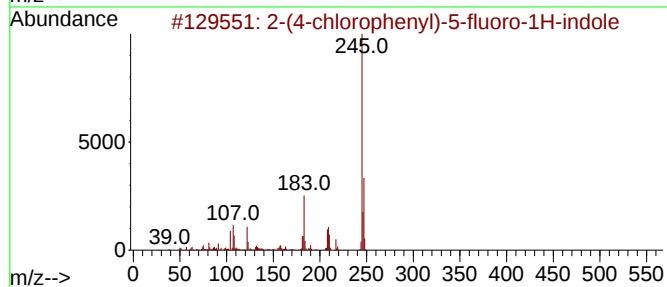
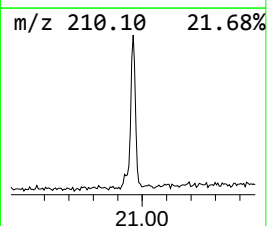
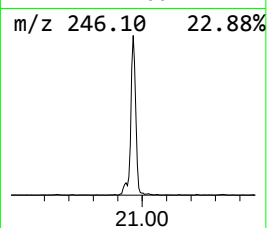
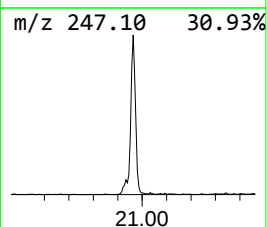
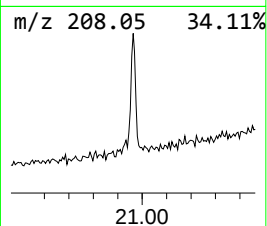
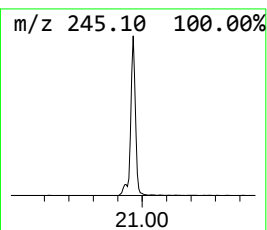
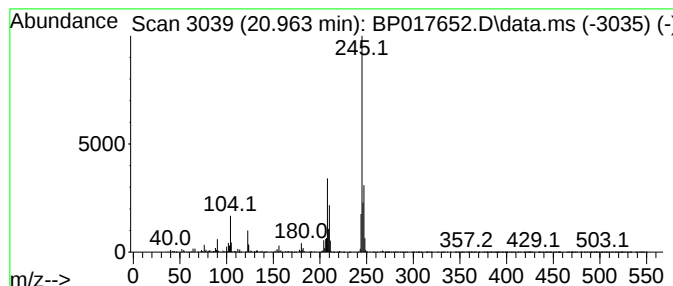
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-(4-chlorophenyl)-5-fluoro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.963	3.84 ng/ul	463707	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-chlorophenyl)-5-fluoro-1H-i...	245	C14H9ClFN	881040-32-0	76		
2	4-Chloro-2-methyl-6-trifluoromet...	245	C11H7ClF3N	867167-05-3	53		
3	2-(4-Chlorophenyl)benzothiazole	245	C13H8ClNS	006265-91-4	53		
4	2-Chloro-4-methylpyridazino(6,1-...	245	C12H8ClN3O	001703-04-4	50		
5	4-Chloro-2-methyl-7-(trifluorome...	245	C11H7ClF3N	018529-09-4	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101123\
Data File : BP017652.D
Acq On : 11 Oct 2023 18:02
Operator : MA/JU
Sample : 04744-06RX
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH583RX

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100923.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
2-Butene, 1-chl...	3.158	48.8	ng/ul	1781860	1	7.922	730738	20.0
Propanoic acid,...	4.558	58.2	ng/ul	2126860	1	7.922	730738	20.0
Butane, 2,3-dic...	4.840	18.3	ng/ul	668244	1	7.922	730738	20.0
unknown-01	7.828	2.0	ng/ul	74089	1	7.922	730738	20.0
Coumarin, 8-ace...	16.345	3.3	ng/ul	328382	4	17.410	1967040	20.0
unknown-02	18.798	2.1	ng/ul	208096	4	17.410	1967040	20.0
2-(4-chlorophen...	20.963	3.8	ng/ul	463707	5	21.551	2412470	20.0