

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
 Data File : BP014766.D
 Acq On : 20 Apr 2023 20:45
 Operator : CG/JU
 Sample : 02296-05DL 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCBK4DL

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-BP041823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014766.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.216	3	5	14	rVB	135585	142950	2.36%	0.260%
2	6.652	585	589	593	rBV3	24291	41882	0.69%	0.076%
3	7.128	666	670	676	rVB3	19815	32949	0.55%	0.060%
4	7.287	691	697	702	rVV	19196	30399	0.50%	0.055%
5	7.363	704	710	716	rVB	18492	28270	0.47%	0.051%
6	7.463	721	727	736	rBV	72174	127266	2.11%	0.232%
7	7.634	750	756	758	rVV2	29448	48786	0.81%	0.089%
8	7.669	759	762	773	rVB	54583	97505	1.61%	0.178%
9	7.793	778	783	792	rBV2	28716	49566	0.82%	0.090%
10	8.151	835	844	857	rVB	902331	1428161	23.63%	2.600%
11	8.287	857	867	874	rBV4	31089	72401	1.20%	0.132%
12	8.528	902	908	914	rBV	39576	68125	1.13%	0.124%
13	8.693	932	936	944	rVV2	23552	40221	0.67%	0.073%
14	8.793	944	953	957	rVV	37121	65551	1.08%	0.119%
15	8.846	957	962	973	rVV2	44120	78391	1.30%	0.143%
16	9.263	1028	1033	1037	rBV	29120	41813	0.69%	0.076%
17	9.316	1037	1042	1052	rVV2	64788	131400	2.17%	0.239%
18	9.404	1052	1057	1063	rVB2	14691	24096	0.40%	0.044%
19	9.793	1111	1123	1129	rBV2	30903	63657	1.05%	0.116%
20	9.851	1129	1133	1140	rVB2	29880	44180	0.73%	0.080%
21	10.045	1160	1166	1171	rBV	35411	58748	0.97%	0.107%
22	10.098	1171	1175	1181	rVV3	37145	60833	1.01%	0.111%
23	10.216	1189	1195	1201	rVV	33574	61561	1.02%	0.112%
24	10.334	1210	1215	1218	rVV	59578	103955	1.72%	0.189%
25	10.381	1218	1223	1232	rVV3	103548	237298	3.93%	0.432%
26	10.522	1240	1247	1252	rVV	93918	175004	2.90%	0.319%
27	10.587	1252	1258	1268	rVV	83269	152933	2.53%	0.278%
28	10.981	1312	1325	1328	rVV	1195117	2020704	33.43%	3.679%
29	11.034	1328	1334	1341	rVV	3535333	6044776	100.00%	11.004%
30	11.104	1341	1346	1350	rVV	51803	99440	1.65%	0.181%
31	11.151	1350	1354	1362	rVB	41368	73890	1.22%	0.135%
32	11.328	1378	1384	1391	rVB4	13852	23926	0.40%	0.044%
33	11.634	1432	1436	1442	rBV3	17525	29139	0.48%	0.053%
34	12.345	1551	1557	1566	rVB	67943	110304	1.82%	0.201%
35	12.522	1580	1587	1596	rVV2	46630	97427	1.61%	0.177%
36	12.622	1597	1604	1611	rVV	2490365	3781964	62.57%	6.885%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
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37	12.734	1616	1623	1630	rVV	326331	512345	8.48%	0.933%
38	12.839	1630	1641	1650	rVB	2067114	3146383	52.05%	5.728%
39	13.616	1767	1773	1779	rVV	788900	1115631	18.46%	2.031%
40	13.810	1800	1806	1817	rVB	165075	312563	5.17%	0.569%

41	13.939	1819	1828	1838	rBV	164233	386155	6.39%	0.703%
42	14.098	1846	1855	1861	rBV	411662	732289	12.11%	1.333%
43	14.157	1861	1865	1866	rVV	192187	251901	4.17%	0.459%
44	14.175	1866	1868	1875	rVB	248618	317725	5.26%	0.578%
45	14.339	1890	1896	1899	rVV	121347	199472	3.30%	0.363%

46	14.369	1899	1901	1906	rVB	65789	76617	1.27%	0.139%
47	14.475	1912	1919	1922	rBV3	253988	477537	7.90%	0.869%
48	14.780	1959	1971	1976	rBV	1907451	2930901	48.49%	5.336%
49	14.845	1976	1982	1988	rVV	2637741	3782803	62.58%	6.886%
50	14.904	1988	1992	1996	rVV	52467	73813	1.22%	0.134%

51	14.980	2001	2005	2012	rVV2	21826	55150	0.91%	0.100%
52	15.039	2012	2015	2019	rVV	34837	55088	0.91%	0.100%
53	15.086	2019	2023	2028	rVV	112609	175366	2.90%	0.319%
54	15.175	2032	2038	2046	rVB	1842293	2537659	41.98%	4.620%
55	15.245	2046	2050	2054	rBV	29118	42062	0.70%	0.077%

56	15.392	2068	2075	2079	rBV	31899	56541	0.94%	0.103%
57	15.445	2079	2084	2088	rVV3	16239	26720	0.44%	0.049%
58	15.557	2095	2103	2107	rVV4	36493	72290	1.20%	0.132%
59	15.598	2107	2110	2115	rVV5	15881	24412	0.40%	0.044%
60	15.692	2122	2126	2129	rVV4	18400	27665	0.46%	0.050%

61	15.769	2130	2139	2143	rVV	322477	491082	8.12%	0.894%
62	15.828	2143	2149	2154	rVV	1836769	2685682	44.43%	4.889%
63	15.869	2154	2156	2160	rVV2	43646	54242	0.90%	0.099%
64	15.916	2160	2164	2166	rVV3	58371	88187	1.46%	0.161%
65	15.945	2166	2169	2172	rVV	99999	147612	2.44%	0.269%

66	15.986	2172	2176	2181	rVV3	124872	208412	3.45%	0.379%
67	16.033	2181	2184	2187	rVV2	44520	65902	1.09%	0.120%
68	16.075	2187	2191	2194	rVV3	76148	115691	1.91%	0.211%
69	16.122	2194	2199	2207	rVV2	227566	408480	6.76%	0.744%
70	16.222	2211	2216	2218	rVV4	20423	34021	0.56%	0.062%

71	16.263	2218	2223	2231	rVV	182438	365093	6.04%	0.665%
72	16.327	2231	2234	2236	rVV2	22006	29228	0.48%	0.053%
73	16.357	2236	2239	2244	rVB2	32987	41868	0.69%	0.076%
74	16.498	2252	2263	2269	rBV2	462742	731637	12.10%	1.332%
75	16.580	2272	2277	2280	rVV	36442	59486	0.98%	0.108%

76	16.616	2280	2283	2289	rVV	54049	87101	1.44%	0.159%
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 Title : SVOA CALIBRATION

77	16.698	2294	2297	2301	rVV4	19473	29065	0.48%	0.053%
78	16.792	2308	2313	2316	rVV4	32157	71215	1.18%	0.130%
79	16.827	2316	2319	2324	rVV	65851	92609	1.53%	0.169%
80	16.875	2324	2327	2334	rVV3	19718	32830	0.54%	0.060%
81	16.980	2340	2345	2352	rVB4	38842	69405	1.15%	0.126%
82	17.139	2367	2372	2375	rBV2	58525	88608	1.47%	0.161%
83	17.174	2375	2378	2386	rVB	66840	99349	1.64%	0.181%
84	17.345	2402	2407	2415	rVB	145544	202885	3.36%	0.369%
85	17.527	2426	2438	2442	rBV	2322822	3368912	55.73%	6.133%
86	17.574	2442	2446	2451	rVB	1452445	2006829	33.20%	3.653%
87	17.627	2451	2455	2459	rBV	316488	428688	7.09%	0.780%
88	17.733	2466	2473	2478	rVB5	37417	68528	1.13%	0.125%
89	17.922	2500	2505	2509	rBV	88187	122637	2.03%	0.223%
90	18.380	2579	2583	2587	rBV2	54874	73197	1.21%	0.133%
91	18.427	2587	2591	2596	rVB2	39629	56751	0.94%	0.103%
92	18.574	2610	2616	2621	rBV	116083	166137	2.75%	0.302%
93	18.745	2641	2645	2648	rVB2	18200	24661	0.41%	0.045%
94	18.857	2659	2664	2668	rBV7	11191	24161	0.40%	0.044%
95	19.516	2772	2776	2782	rVB	146242	185856	3.07%	0.338%
96	19.745	2812	2815	2819	rBV4	39185	47427	0.78%	0.086%
97	19.839	2826	2831	2841	rBV2	451581	692501	11.46%	1.261%
98	21.598	3125	3130	3142	rBV	2752631	3761200	62.22%	6.847%
99	24.004	3534	3539	3545	rVB2	238704	434633	7.19%	0.791%
100	24.174	3560	3568	3582	rVB2	1850786	3990833	66.02%	7.265%

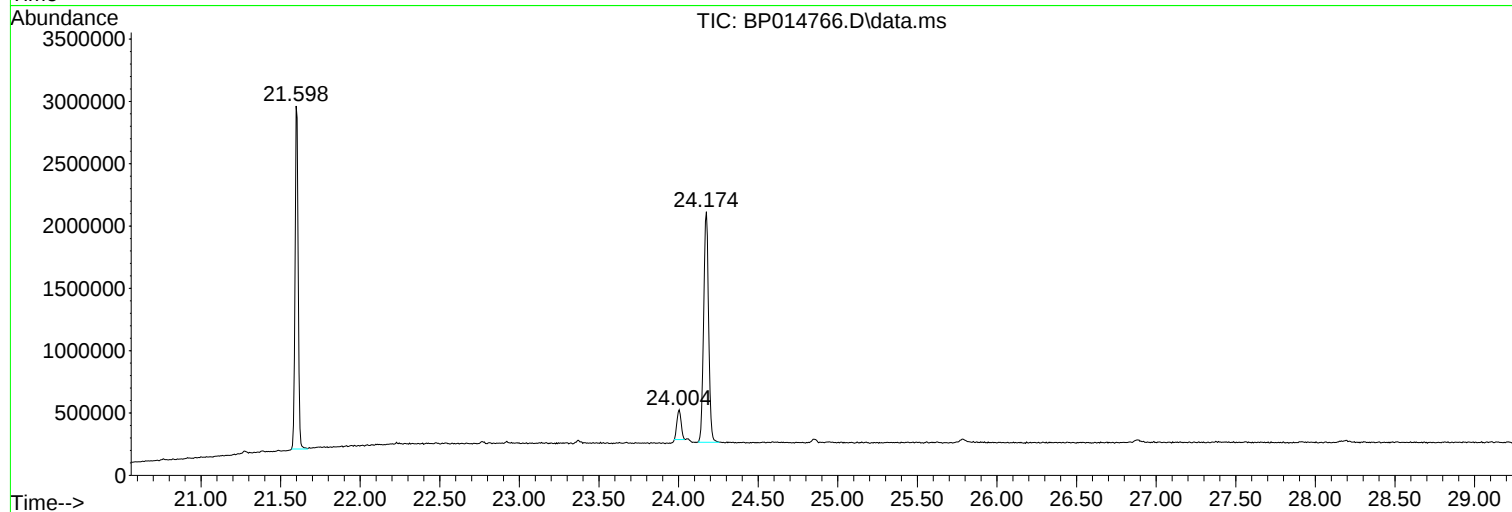
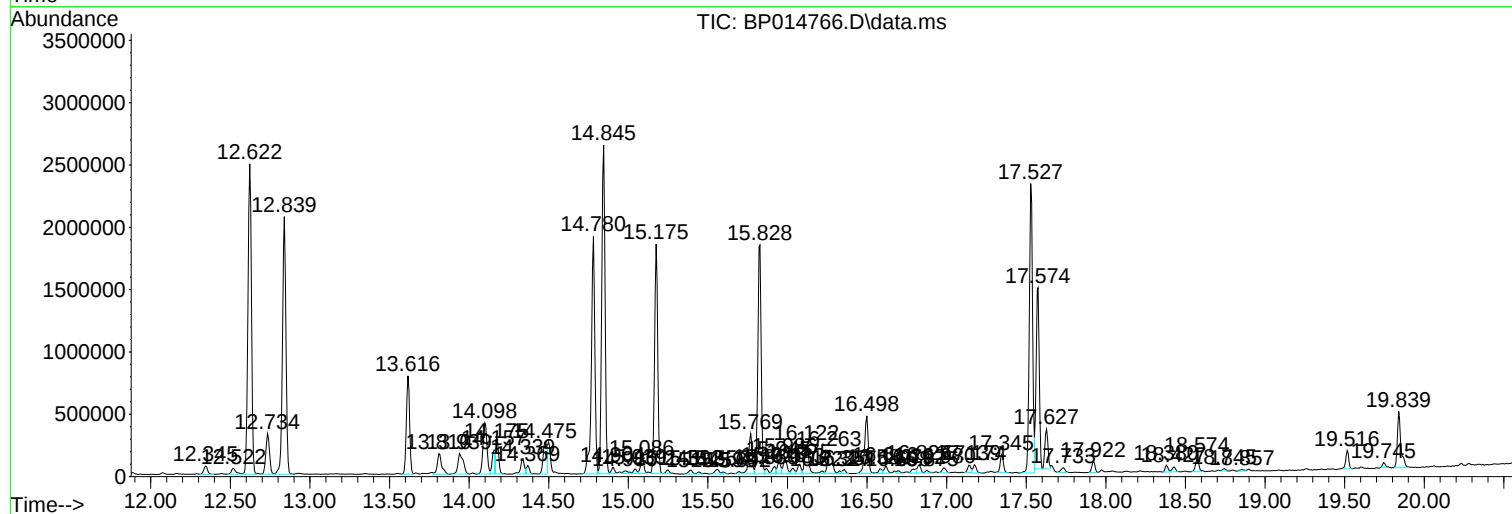
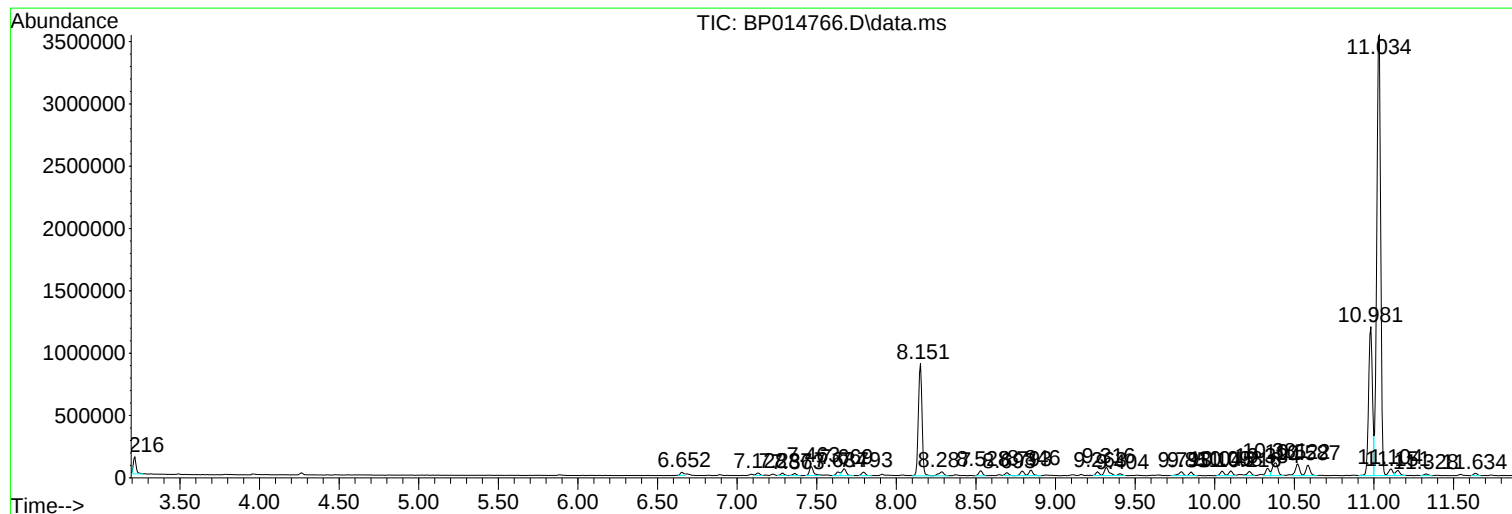
Sum of corrected areas: 54931200

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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



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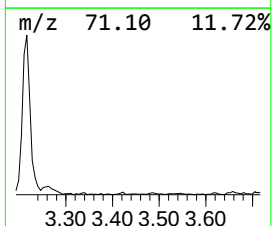
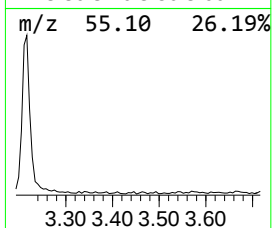
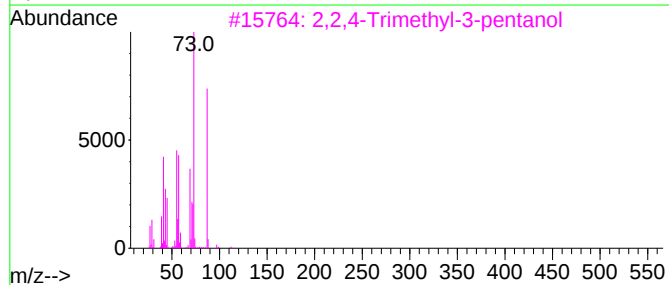
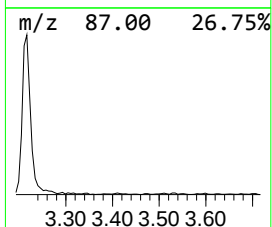
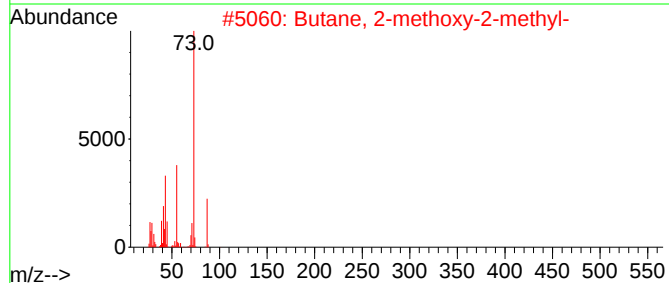
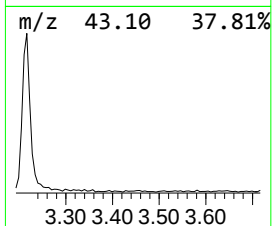
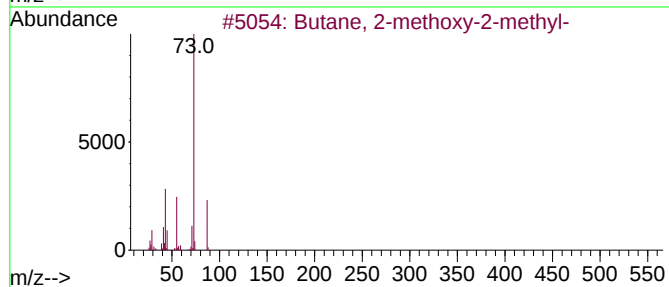
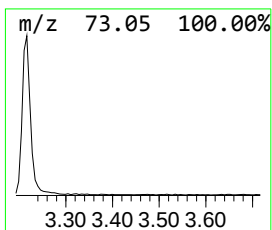
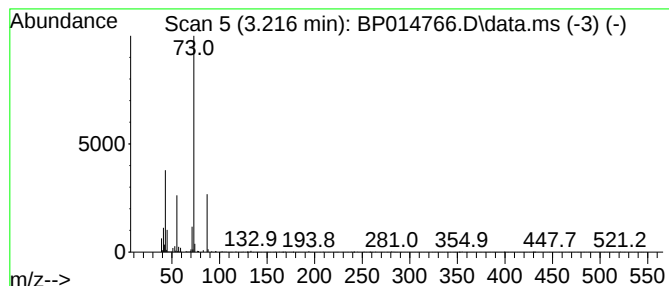
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.216	2.00 ng/ul	142950	1,4-Dichlorobenzene-d4	8.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	17
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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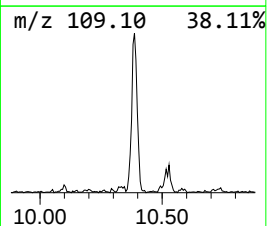
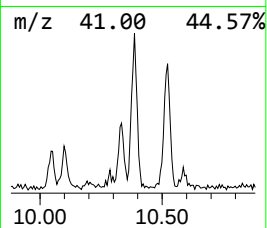
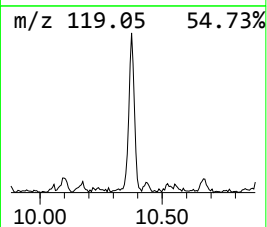
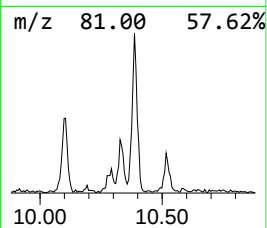
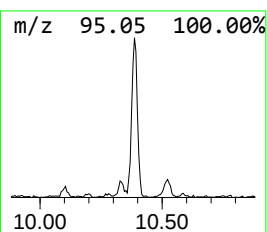
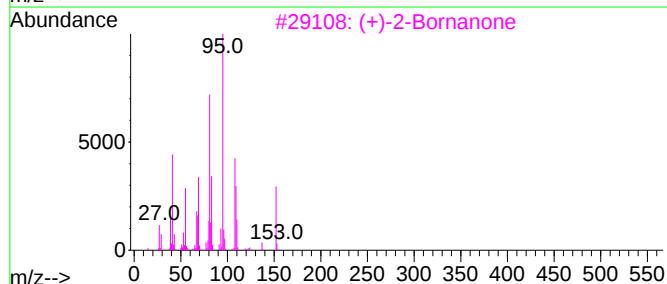
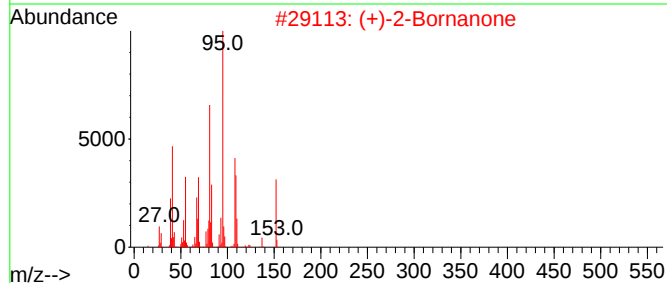
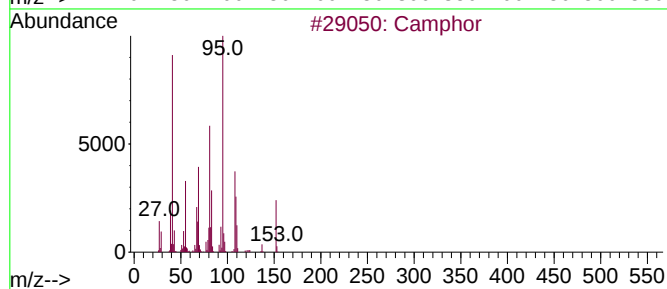
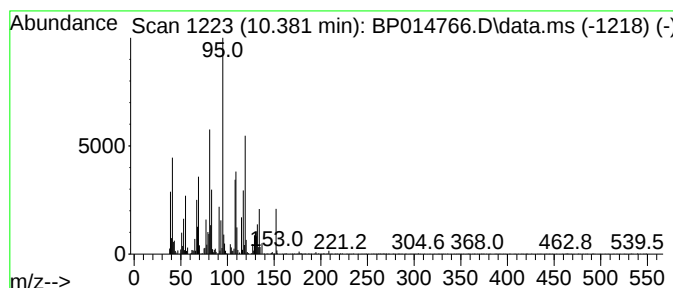
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Camphor Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.381	2.35 ng/ul	237298	Naphthalene-d8	10.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphor	152	C10H16O	000076-22-2	95
2			(+)-2-Bornanone	152	C10H16O	000464-49-3	91
3			(+)-2-Bornanone	152	C10H16O	000464-49-3	91
4			(+)-2-Bornanone	152	C10H16O	000464-49-3	91
5			(+)-2-Bornanone	152	C10H16O	000464-49-3	90



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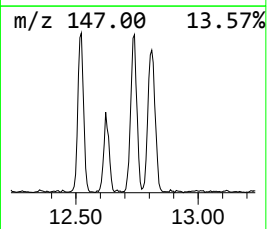
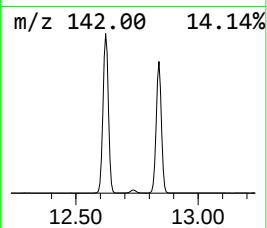
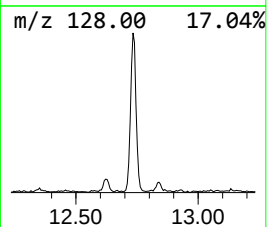
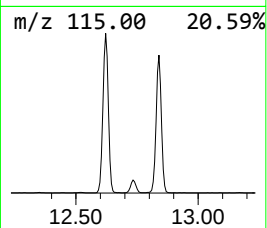
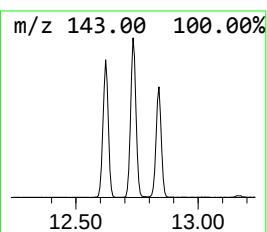
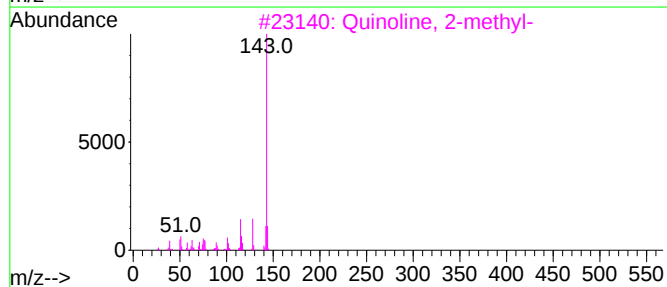
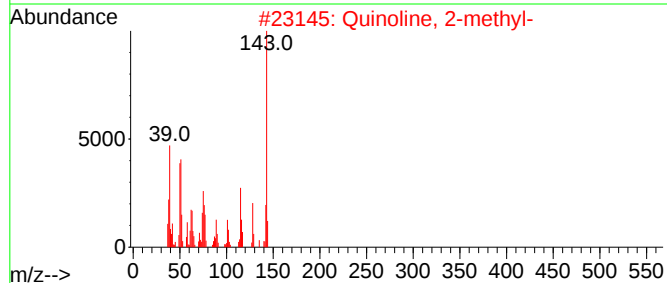
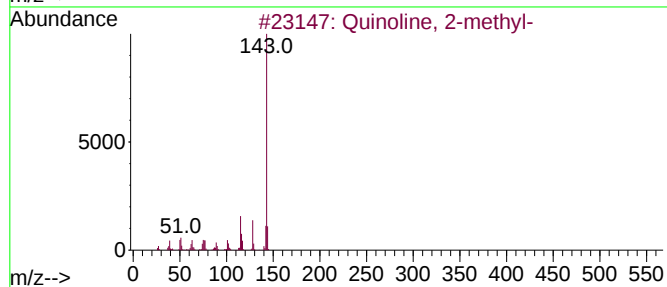
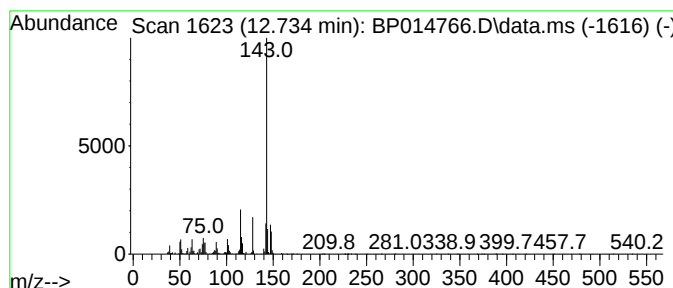
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Quinoline, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	5.07 ng/ul	512345	Naphthalene-d8	10.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Quinoline, 2-methyl-	143	C10H9N	000091-63-4	95
2			Quinoline, 2-methyl-	143	C10H9N	000091-63-4	93
3			Quinoline, 2-methyl-	143	C10H9N	000091-63-4	93
4			Quinoline, 4-methyl-	143	C10H9N	000491-35-0	81
5			Quinoline, 4-methyl-	143	C10H9N	000491-35-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
Operator : CG/JU
Sample : 02296-05DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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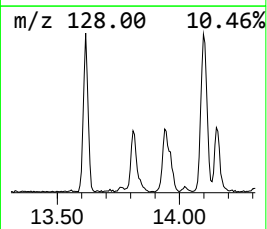
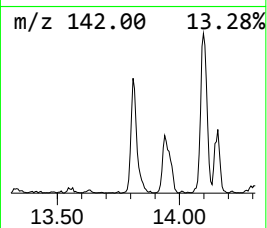
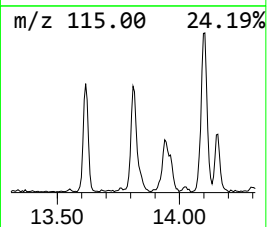
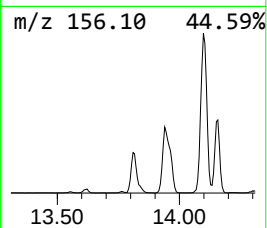
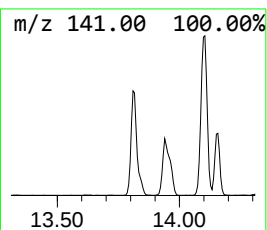
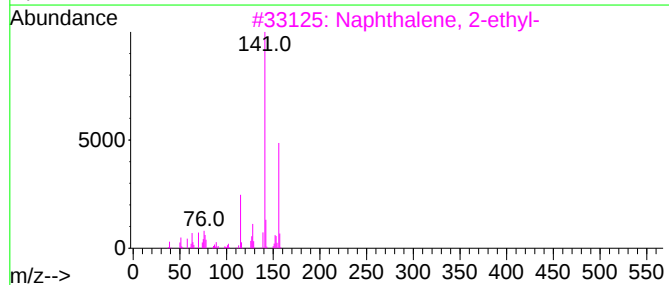
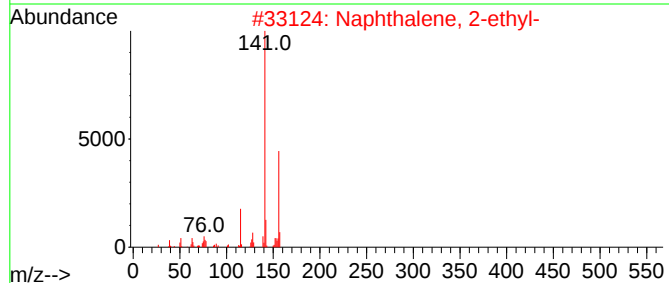
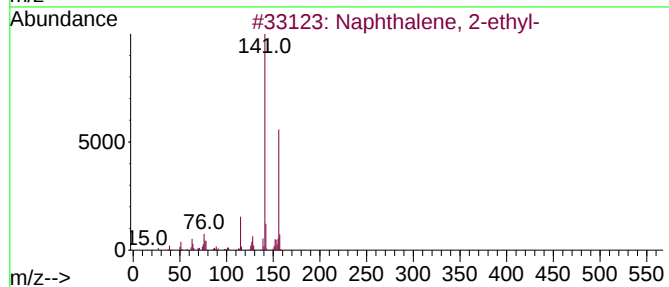
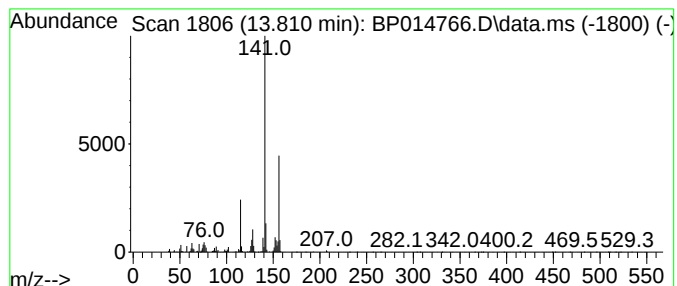
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 2-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	2.13 ng/ul	312563	Acenaphthene-d10	14.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
2			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
3			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
4			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
5			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
Operator : CG/JU
Sample : 02296-05DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCBK4DL

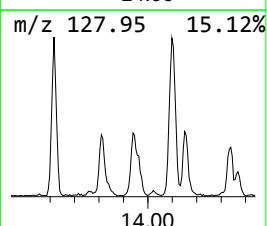
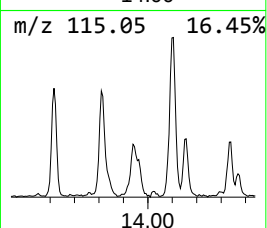
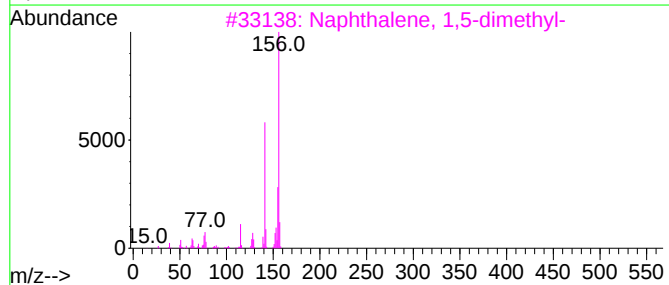
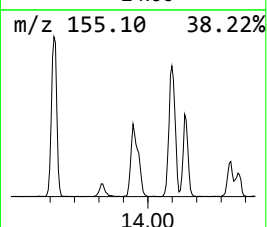
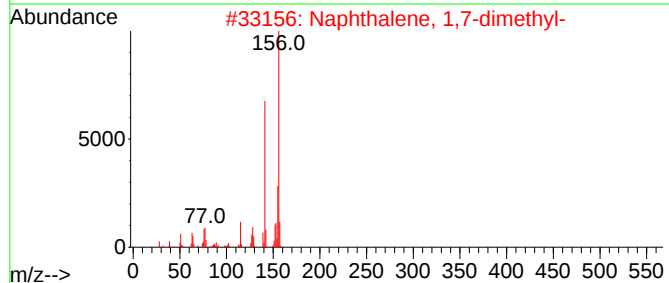
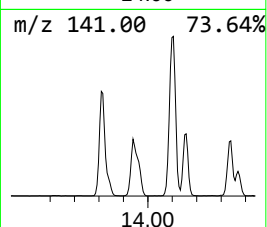
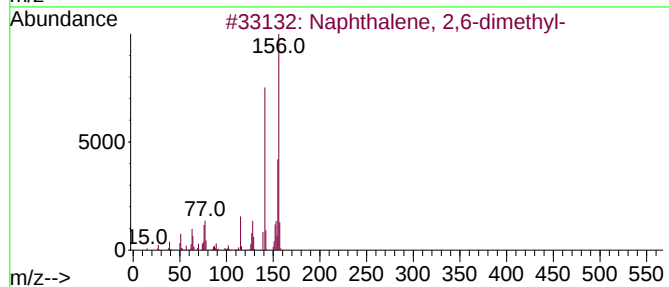
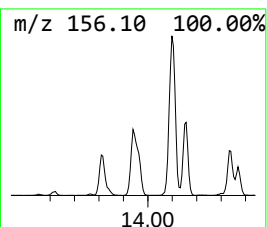
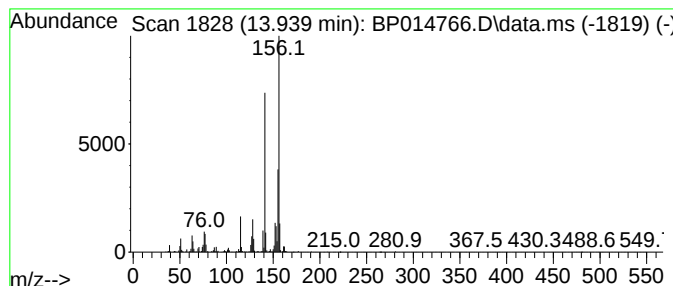
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.939	2.64 ng/ul	386155	Acenaphthene-d10	14.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
Operator : CG/JU
Sample : 02296-05DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

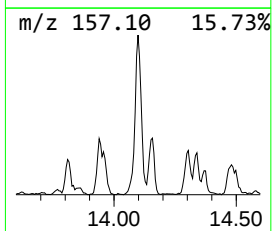
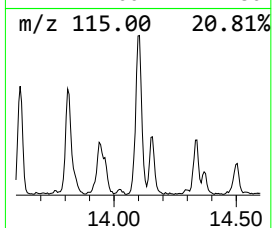
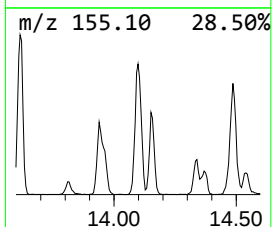
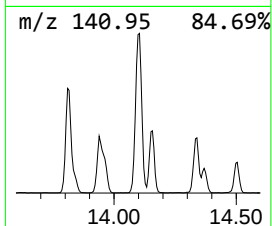
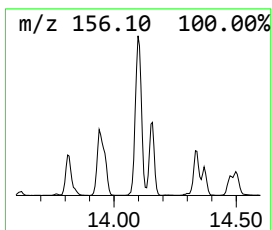
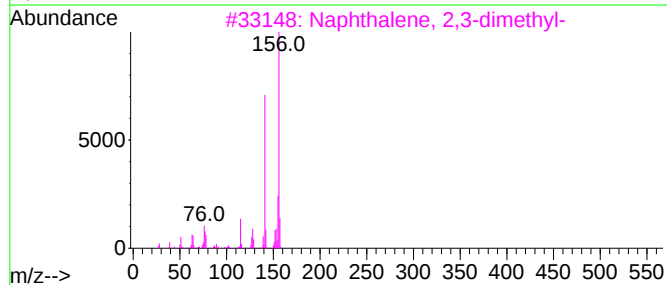
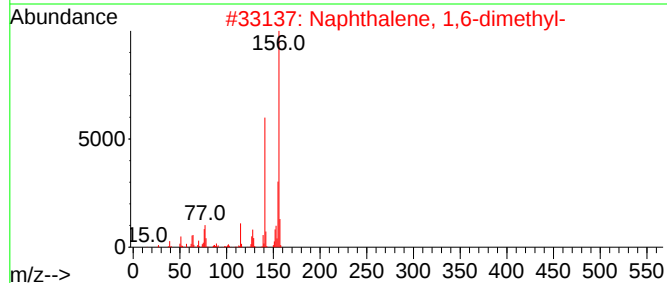
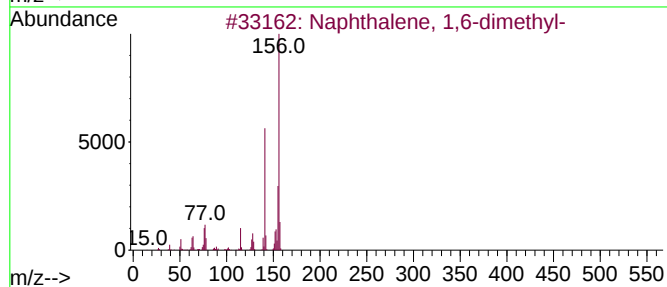
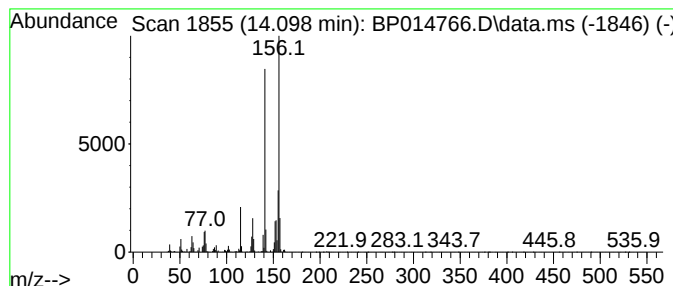
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 1,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.098	5.00 ng/ul	732289	Acenaphthene-d10		14.781	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-		156	C12H12	000575-43-9	97
2	Naphthalene, 1,6-dimethyl-		156	C12H12	000575-43-9	97
3	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	97
4	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	97
5	Naphthalene, 2,3-dimethyl-		156	C12H12	000581-40-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
Operator : CG/JU
Sample : 02296-05DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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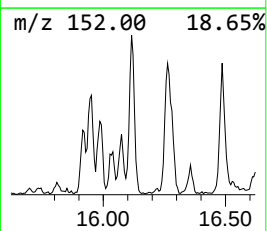
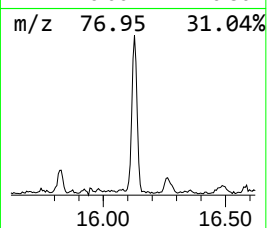
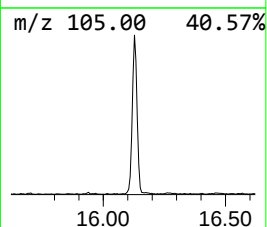
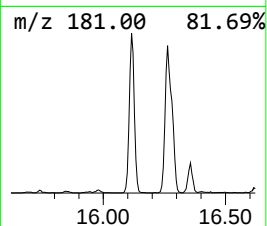
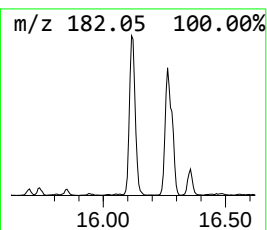
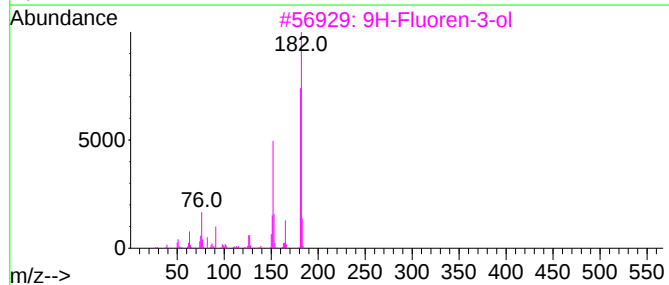
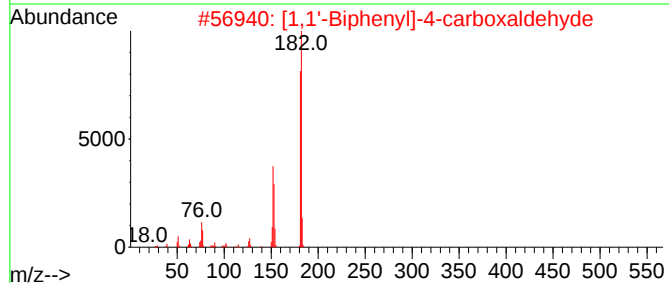
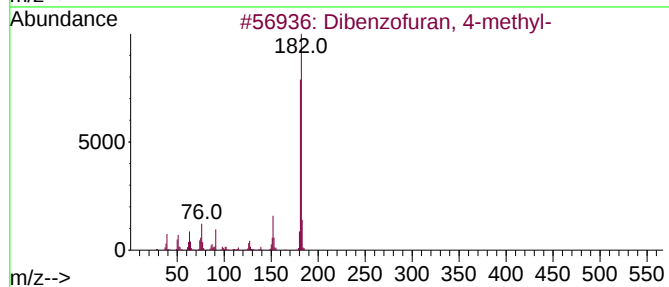
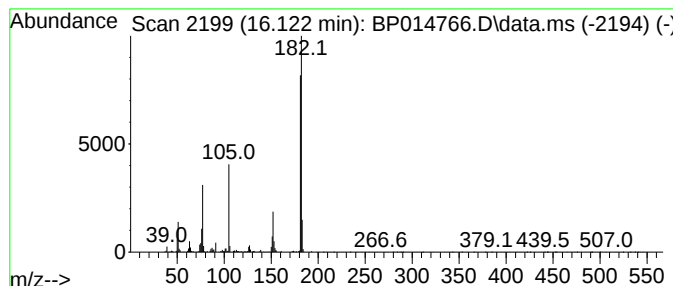
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Dibenzofuran, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.122	2.79 ng/ul	408480	Acenaphthene-d10	14.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	87
2			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	81
3			9H-Fluoren-3-ol	182	C13H10O	006344-67-8	80
4			2-Hydroxyfluorene	182	C13H10O	002443-58-5	80
5			3-(2-Naphthyl)acrylaldehyde	182	C13H10O	020884-05-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
Operator : CG/JU
Sample : 02296-05DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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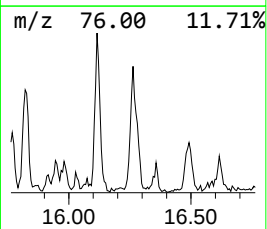
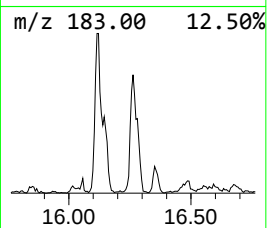
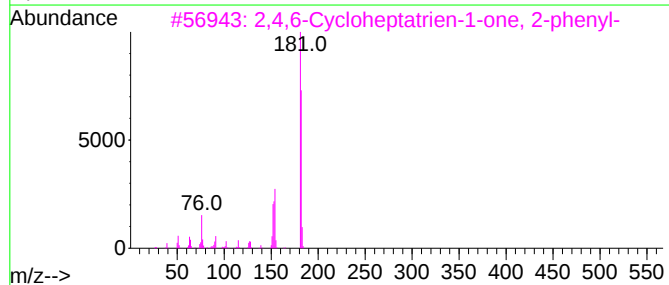
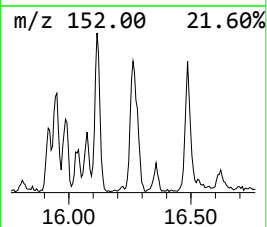
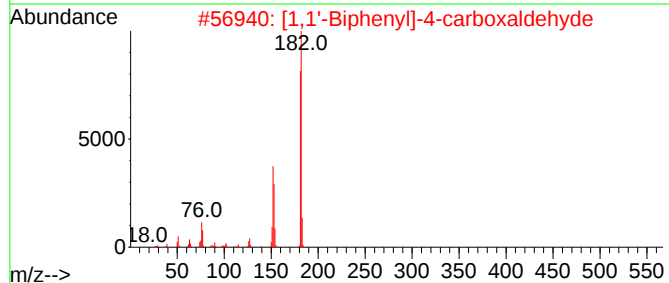
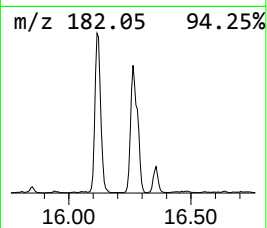
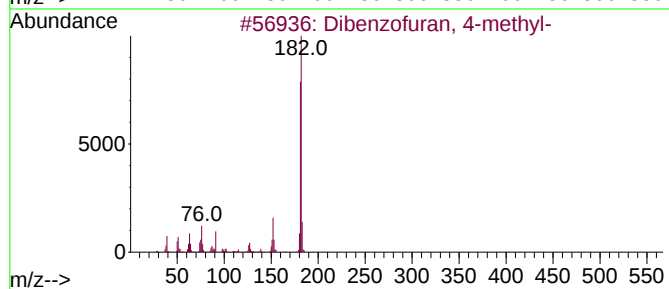
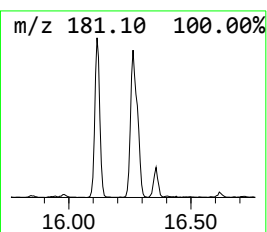
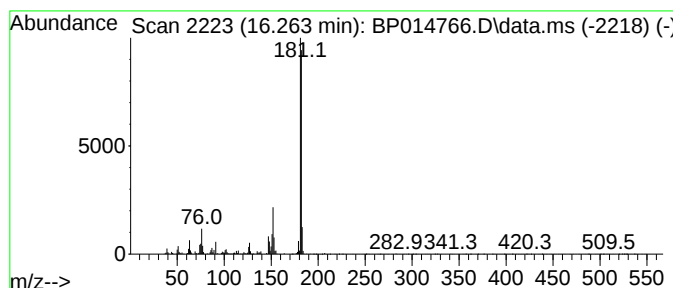
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.263	2.17 ng/ul	365093	Phenanthrene-d10	17.527

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	94
2		[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	91
3		2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86
4		9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64
5		Pyrido[2,3-b]indole, 6-methyl-	182	C12H10N2	108349-67-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
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ALS Vial : 20 Sample Multiplier: 1

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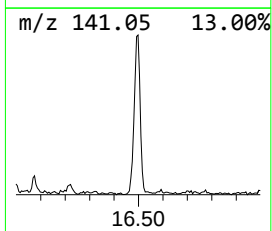
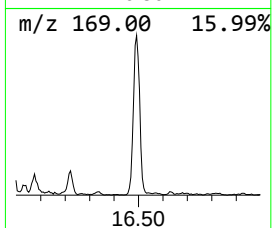
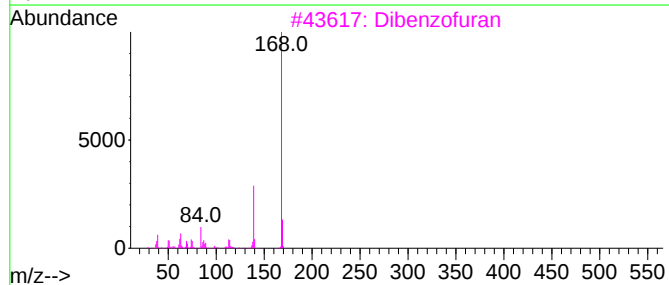
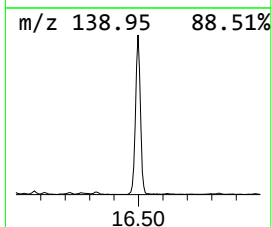
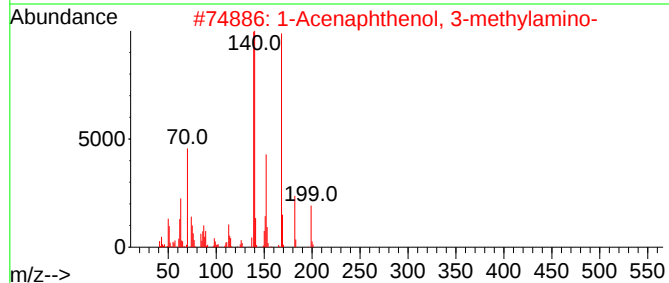
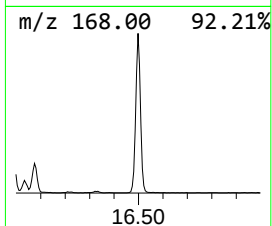
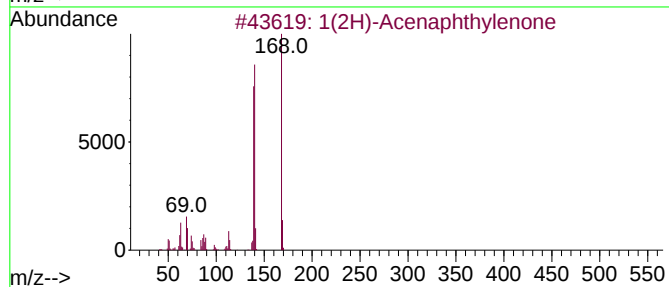
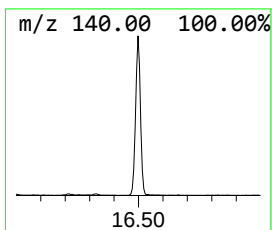
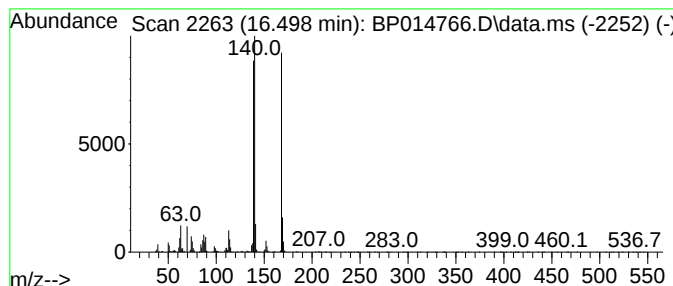
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1(2H)-Acenaphthylenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.498	4.34 ng/ul	731637	Phenanthrene-d10	17.527

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	96
2			1-Acenaphthenol, 3-methylamino-	199	C13H13NO	1000129-22-6	64
3			Dibenzofuran	168	C12H8O	000132-64-9	58
4			Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	58
5			5,6,7,8-Tetrahydro-thiazolo[5,4-...	168	C7H8N2OS	1000317-75-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014766.D
Acq On : 20 Apr 2023 20:45
Operator : CG/JU
Sample : 02296-05DL 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCBK4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.216	2.0	ng/ul	142950	1	8.151	1428160	20.0
Camphor	10.381	2.4	ng/ul	237298	2	10.981	2020700	20.0
Quinoline, 2-me...	12.733	5.1	ng/ul	512345	2	10.981	2020700	20.0
Naphthalene, 2-...	13.810	2.1	ng/ul	312563	3	14.781	2930900	20.0
Naphthalene, 2,...	13.939	2.6	ng/ul	386155	3	14.781	2930900	20.0
Naphthalene, 1,...	14.098	5.0	ng/ul	732289	3	14.781	2930900	20.0
Dibenzofuran, 4...	16.122	2.8	ng/ul	408480	3	14.781	2930900	20.0
[1,1'-Biphenyl]...	16.263	2.2	ng/ul	365093	4	17.527	3368910	20.0
1(2H)-Acenaphth...	16.498	4.3	ng/ul	731637	4	17.527	3368910	20.0