

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
 Data File : BN034654.D
 Acq On : 18 Oct 2024 17:56
 Operator : RC/JU
 Sample : P4380-16
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E27J1

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_N\Methods\SFAM-EPA-BN101624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN034654.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.969	8	14	37	rVB	11725484	17320942	100.00%	16.257%
2	3.211	51	55	60	rVB	88709	103405	0.60%	0.097%
3	3.605	118	122	132	rBV	315602	499162	2.88%	0.469%
4	3.934	173	178	183	rVB	89099	127598	0.74%	0.120%
5	4.440	259	264	273	rVB	48559	78283	0.45%	0.073%
6	4.581	280	288	294	rVB	135330	250869	1.45%	0.235%
7	4.716	301	311	322	rBV	99235	179878	1.04%	0.169%
8	5.187	387	391	394	rVV	50977	80096	0.46%	0.075%
9	5.346	412	418	426	rBV	69109	126954	0.73%	0.119%
10	5.705	472	479	491	rVB2	71393	120545	0.70%	0.113%
11	6.375	587	593	598	rVB2	50844	78857	0.46%	0.074%
12	6.669	633	643	647	rBV3	52583	95000	0.55%	0.089%
13	6.834	665	671	679	rBV	574594	1003981	5.80%	0.942%
14	6.904	679	683	692	rVB	77886	120179	0.69%	0.113%
15	7.004	692	700	707	rBV	1477155	2252857	13.01%	2.114%
16	7.069	707	711	715	rVV2	60290	106804	0.62%	0.100%
17	7.199	726	733	742	rBV	1707019	2595771	14.99%	2.436%
18	7.322	748	754	761	rVB2	126307	221921	1.28%	0.208%
19	7.581	790	798	805	rVB6	40192	88600	0.51%	0.083%
20	7.663	805	812	817	rBV	1285285	2015207	11.63%	1.891%
21	7.740	821	825	829	rVV	272463	371738	2.15%	0.349%
22	7.781	829	832	836	rVB	82844	101708	0.59%	0.095%
23	8.034	870	875	881	rVB	60528	95463	0.55%	0.090%
24	8.098	881	886	891	rBV	101523	153194	0.88%	0.144%
25	8.157	891	896	900	rBV	653168	1082896	6.25%	1.016%
26	8.193	900	902	909	rVB	471152	637407	3.68%	0.598%
27	8.363	924	931	937	rVV	1412062	2259859	13.05%	2.121%
28	8.428	937	942	947	rVV	306833	556151	3.21%	0.522%
29	8.475	947	950	957	rVV2	52912	107021	0.62%	0.100%
30	8.551	957	963	970	rVB2	70297	131917	0.76%	0.124%
31	8.804	1000	1006	1012	rBV	1502469	2335786	13.49%	2.192%
32	8.898	1018	1022	1027	rVB3	52671	89229	0.52%	0.084%
33	8.957	1027	1032	1037	rBV	114435	184882	1.07%	0.174%
34	9.075	1046	1052	1058	rBV7	48319	110328	0.64%	0.104%
35	9.345	1093	1098	1106	rBV2	69344	152238	0.88%	0.143%
36	9.522	1121	1128	1136	rVV	1188367	1918272	11.07%	1.800%

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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_N\Methods\SFAM-EPA-BN101624.MA.M
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37	9.622	1138	1145	1148	rVV2	162635	293813	1.70%	0.276%
38	9.657	1148	1151	1163	rVB3	181916	321357	1.86%	0.302%
39	9.857	1177	1185	1193	rBV4	182245	459146	2.65%	0.431%
40	9.928	1193	1197	1204	rVV	174048	304951	1.76%	0.286%
41	9.998	1204	1209	1212	rVV3	88543	140384	0.81%	0.132%
42	10.057	1212	1219	1229	rVB	1802839	3167627	18.29%	2.973%
43	10.234	1242	1249	1257	rBV2	182877	334045	1.93%	0.314%
44	10.310	1257	1262	1269	rBV	88309	162140	0.94%	0.152%
45	10.440	1277	1284	1288	rBV	1595861	2644001	15.26%	2.482%
46	10.487	1288	1292	1300	rVB	2132248	3543546	20.46%	3.326%
47	10.569	1300	1306	1310	rBV	752800	1314712	7.59%	1.234%
48	10.981	1370	1376	1382	rBV2	44081	100822	0.58%	0.095%
49	11.063	1384	1390	1399	rVB4	44575	105061	0.61%	0.099%
50	11.204	1409	1414	1418	rBV3	47062	88160	0.51%	0.083%
51	11.269	1419	1425	1433	rVB4	104539	208206	1.20%	0.195%
52	11.545	1463	1472	1478	rVV4	84817	228465	1.32%	0.214%
53	11.604	1478	1482	1491	rVV2	53347	96057	0.55%	0.090%
54	11.816	1504	1518	1527	rBV3	149484	400390	2.31%	0.376%
55	11.928	1531	1537	1544	rBV	85437	159709	0.92%	0.150%
56	12.104	1562	1567	1577	rVB	266038	438425	2.53%	0.411%
57	12.228	1581	1588	1591	rBV2	92748	173246	1.00%	0.163%
58	12.322	1599	1604	1613	rVB	283065	492687	2.84%	0.462%
59	12.422	1618	1621	1628	rVB4	82552	153870	0.89%	0.144%
60	13.034	1718	1725	1733	rBV9	44174	141288	0.82%	0.133%
61	13.128	1737	1741	1744	rBV	93222	148504	0.86%	0.139%
62	13.210	1750	1755	1760	rBV	319124	464033	2.68%	0.436%
63	13.457	1793	1797	1805	rBV5	85448	183271	1.06%	0.172%
64	13.522	1805	1808	1813	rBV6	51323	83114	0.48%	0.078%
65	13.622	1819	1825	1830	rVB	167232	274709	1.59%	0.258%
66	13.669	1830	1833	1836	rBV	68180	94360	0.54%	0.089%
67	13.716	1836	1841	1848	rVB	2452332	3321525	19.18%	3.118%
68	13.798	1851	1855	1857	rBV2	51723	80179	0.46%	0.075%
69	13.986	1881	1887	1897	rVV	3190260	4986756	28.79%	4.680%
70	14.157	1911	1916	1925	rVB	534516	758587	4.38%	0.712%
71	14.292	1933	1939	1946	rBV	1870771	2911538	16.81%	2.733%
72	14.357	1946	1950	1955	rVB	203871	297230	1.72%	0.279%
73	14.492	1968	1973	1978	rBV2	419665	592471	3.42%	0.556%
74	14.575	1984	1987	1992	rVB3	90194	114492	0.66%	0.107%
75	14.645	1995	1999	2003	rVB4	86630	120567	0.70%	0.113%
76	14.698	2003	2008	2015	rVB	185146	318733	1.84%	0.299%

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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 Title : SVOA CALIBRATION

77	14.775	2016	2021	2026	rVB6	74602	154014	0.89%	0.145%
78	14.833	2027	2031	2035	rBV2	66498	103348	0.60%	0.097%
79	15.216	2093	2096	2101	rVB2	98713	127692	0.74%	0.120%
80	15.292	2103	2109	2114	rVV	3596406	5073569	29.29%	4.762%
81	15.345	2114	2118	2123	rVB	288491	393238	2.27%	0.369%
82	15.404	2123	2128	2135	rBV	834107	1161367	6.70%	1.090%
83	15.557	2151	2154	2163	rVB	150938	212811	1.23%	0.200%
84	15.745	2181	2186	2191	rVB	900988	1131835	6.53%	1.062%
85	15.963	2220	2223	2225	rBV	98133	115173	0.66%	0.108%
86	16.133	2249	2252	2255	rBV	83221	103009	0.59%	0.097%
87	16.345	2284	2288	2293	rVB	612920	783706	4.52%	0.736%
88	16.663	2339	2342	2344	rBV	85842	107489	0.62%	0.101%
89	17.039	2401	2406	2410	rBV	1906139	2670284	15.42%	2.506%
90	17.080	2410	2413	2416	rVV	278255	345182	1.99%	0.324%
91	17.139	2417	2423	2435	rVB	3599522	5196705	30.00%	4.878%
92	17.439	2470	2474	2480	rVB	348139	511043	2.95%	0.480%
93	17.598	2498	2501	2506	rVB	176281	221192	1.28%	0.208%
94	18.933	2725	2728	2732	rVB	201068	239411	1.38%	0.225%
95	19.433	2807	2813	2826	rVB	4006885	5623683	32.47%	5.278%
96	20.221	2944	2947	2951	rBV	268369	310963	1.80%	0.292%
97	20.939	3065	3069	3085	rVB	885888	1318346	7.61%	1.237%
98	21.262	3119	3124	3136	rVB2	2081928	2910940	16.81%	2.732%
99	23.956	3574	3582	3595	rVB	2720923	6439645	37.18%	6.044%
100	24.151	3607	3615	3625	rVB	1356857	3312079	19.12%	3.109%

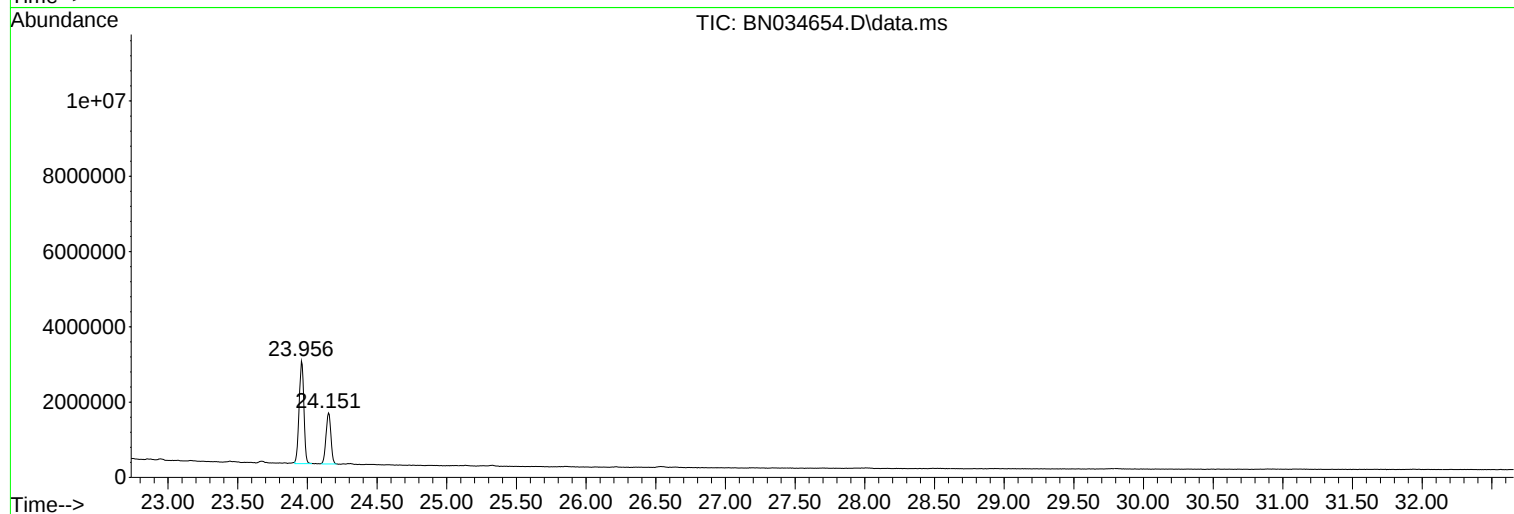
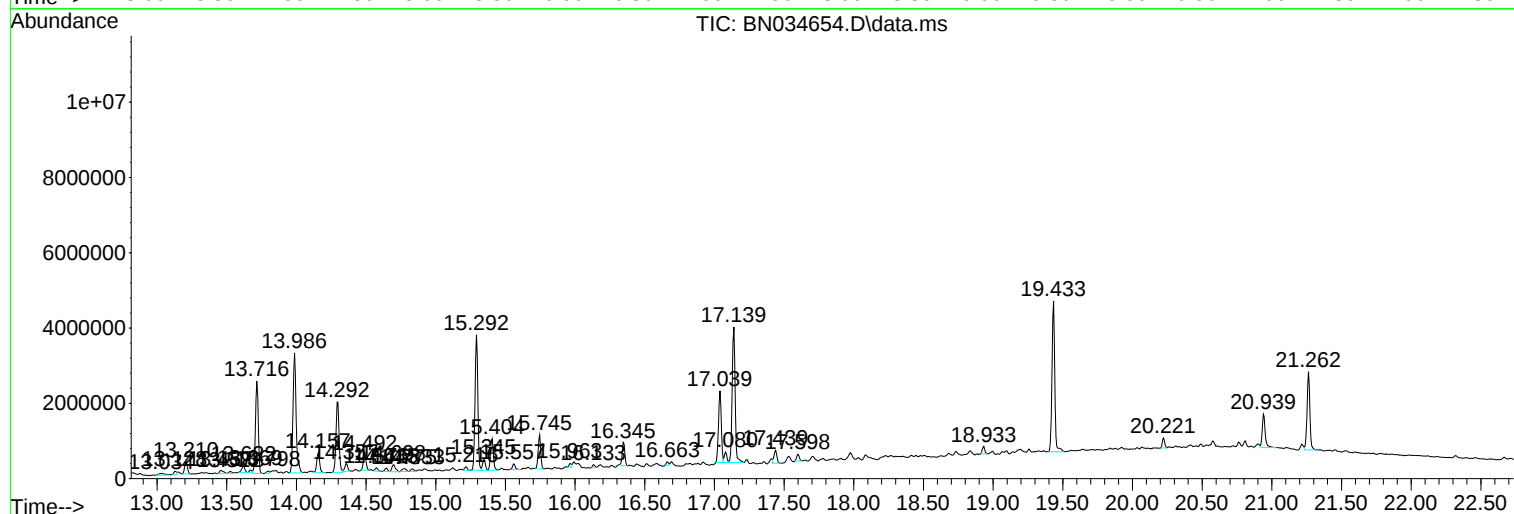
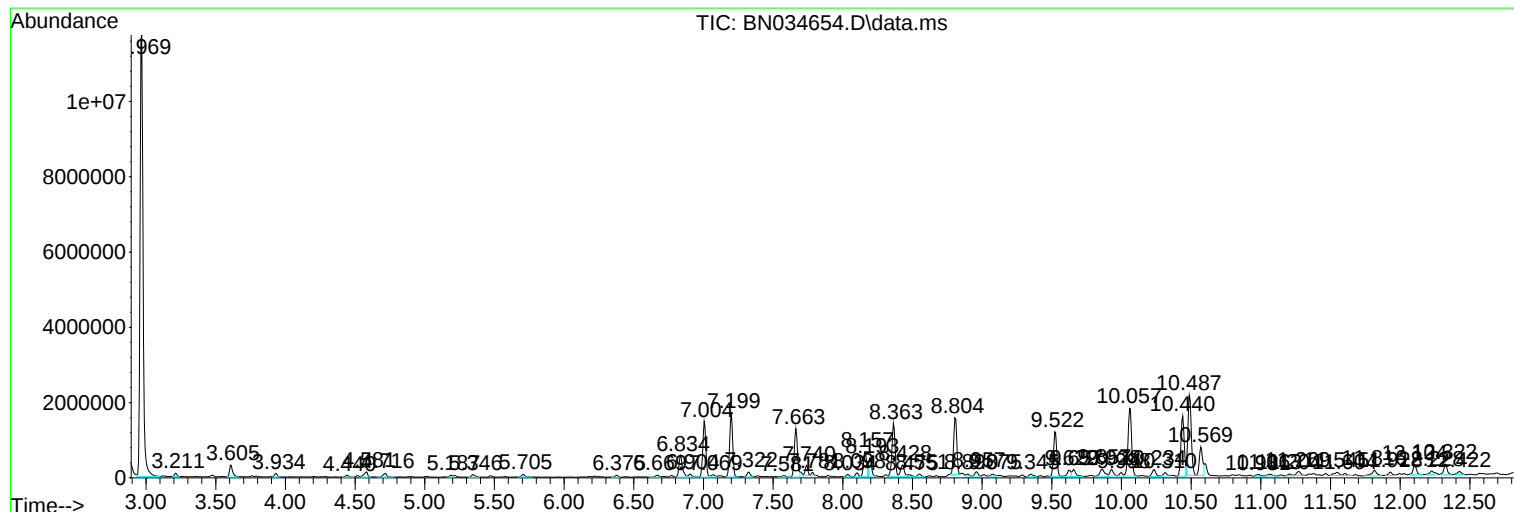
Sum of corrected areas: 106543899

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
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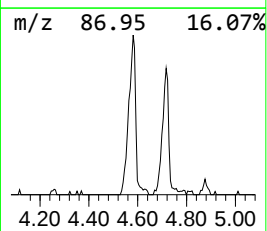
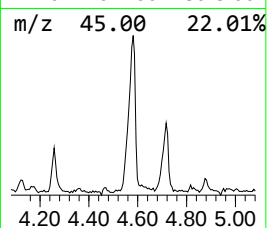
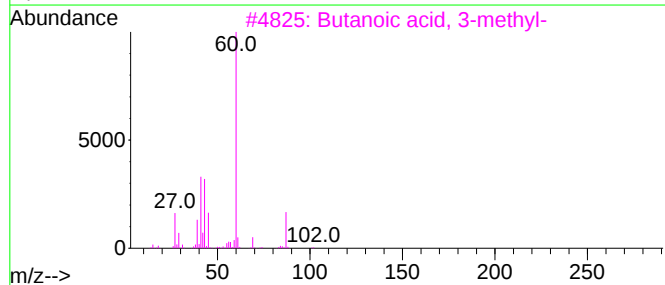
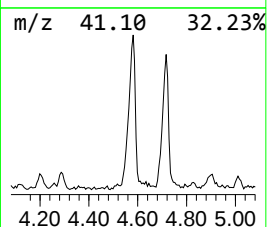
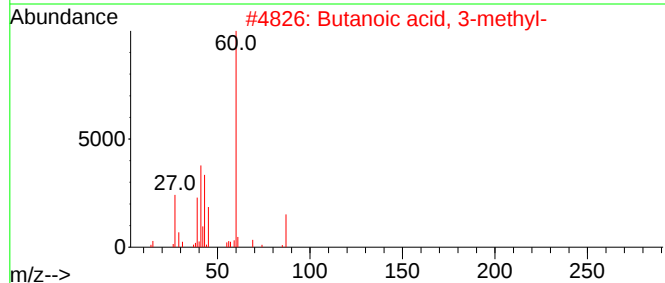
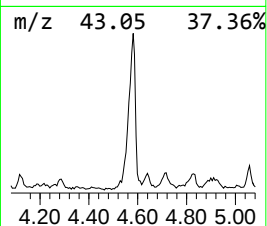
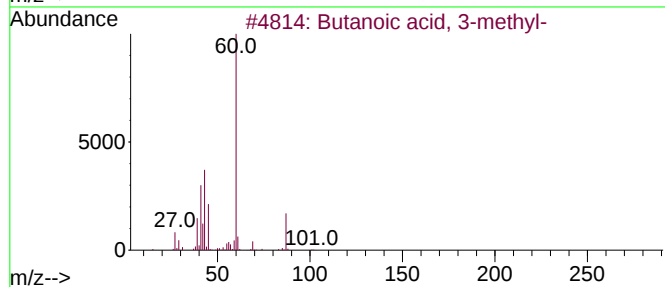
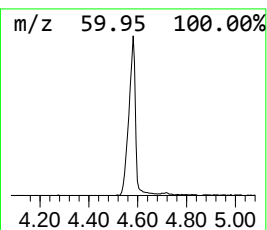
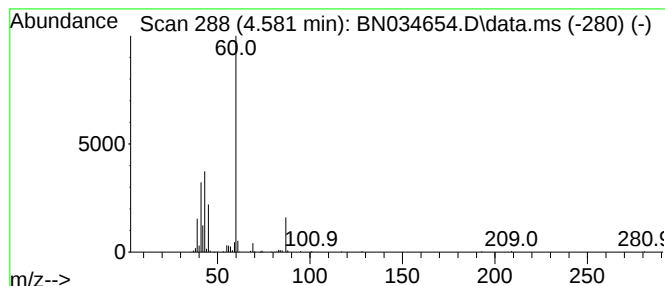
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butanoic acid, 3-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.581	2.49 ng/ul	250869	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	90
2			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	83
3			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
4			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64
5			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	56



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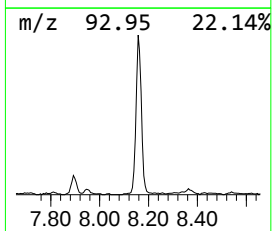
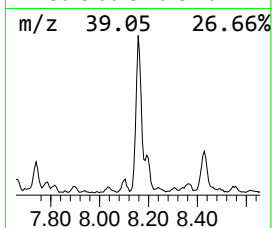
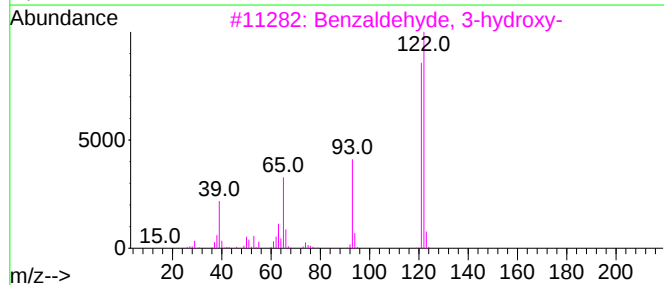
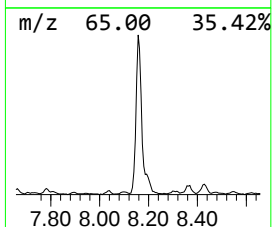
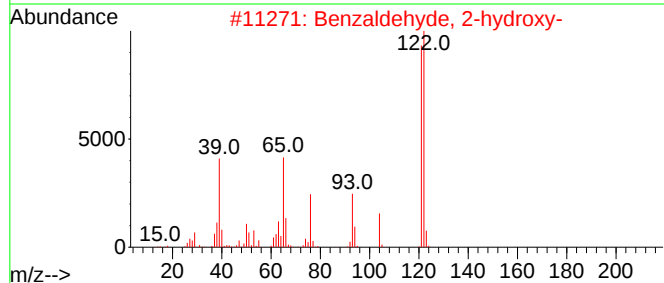
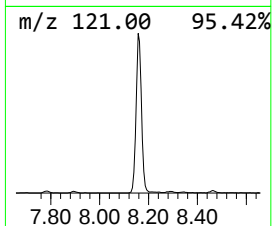
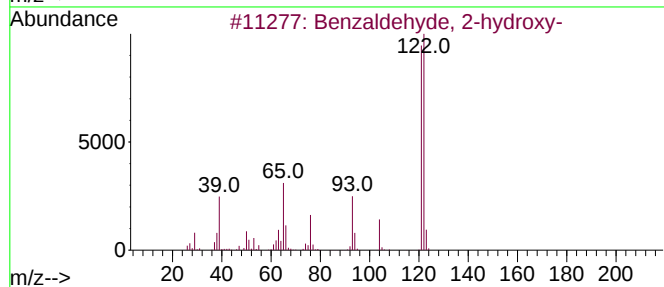
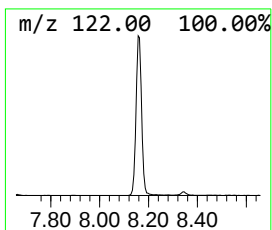
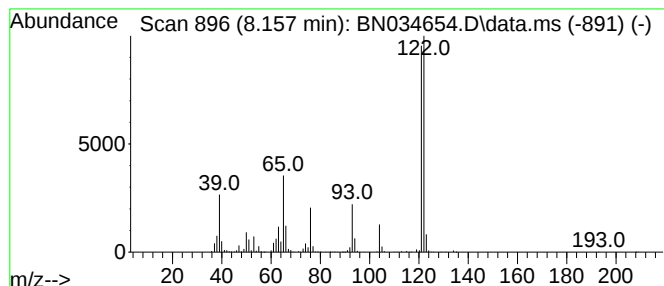
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzaldehyde, 2-hydroxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.157	10.75 ng/ul	1082900	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 2-hydroxy-	122	C7H6O2	000090-02-8	96
2			Benzaldehyde, 2-hydroxy-	122	C7H6O2	000090-02-8	95
3			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	90
4			Benzaldehyde, 4-hydroxy-	122	C7H6O2	000123-08-0	90
5			Benzaldehyde, 3-hydroxy-	122	C7H6O2	000100-83-4	90



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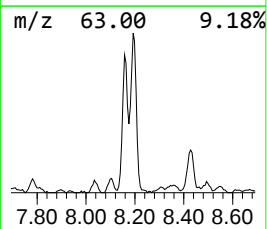
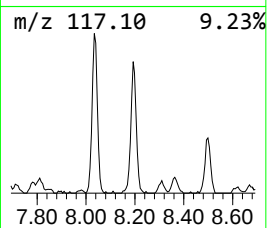
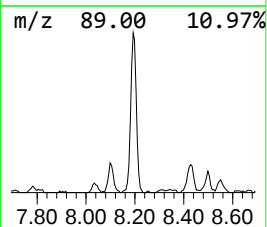
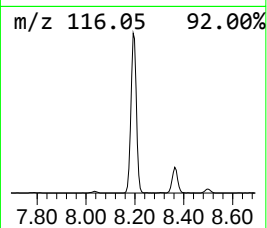
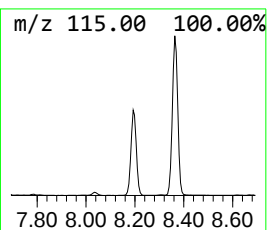
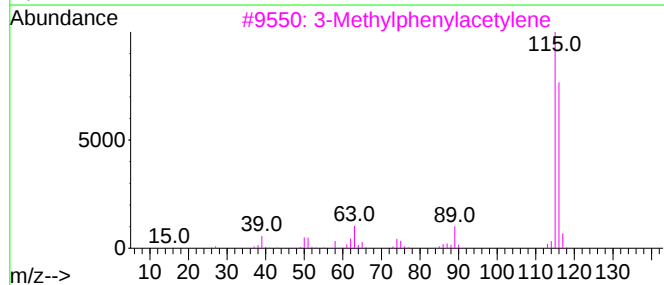
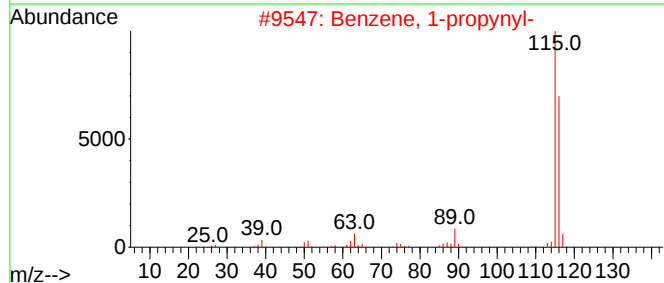
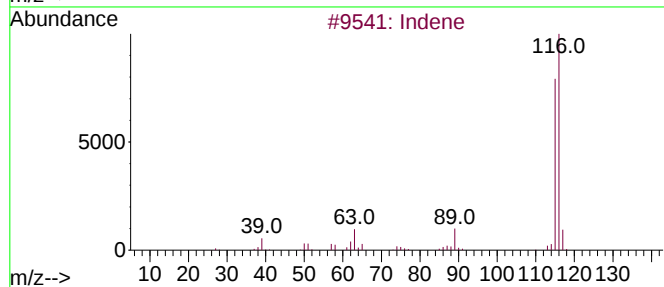
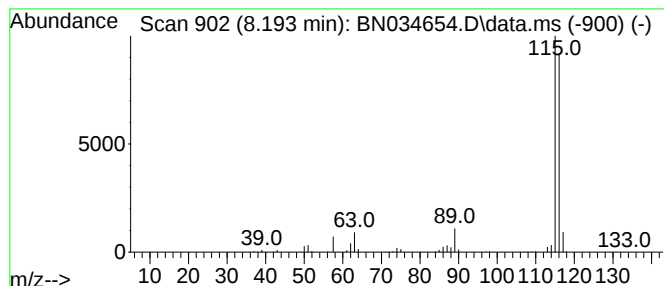
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Indene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.193	6.33 ng/ul	637407	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	95
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
3			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
Acq On : 18 Oct 2024 17:56
Operator : RC/JU
Sample : P4380-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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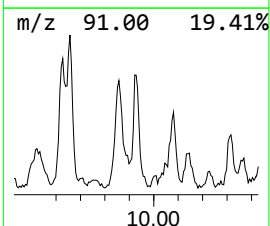
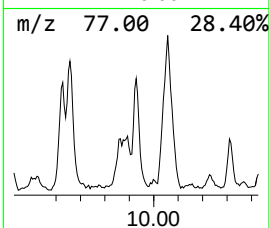
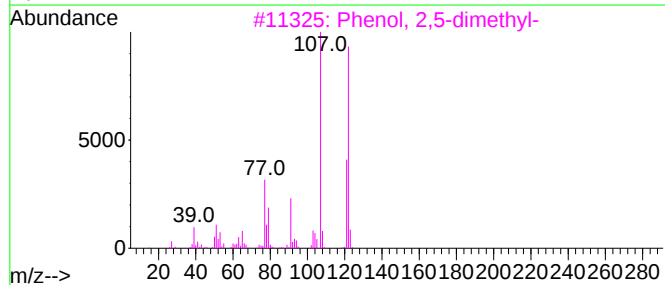
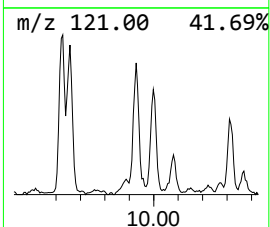
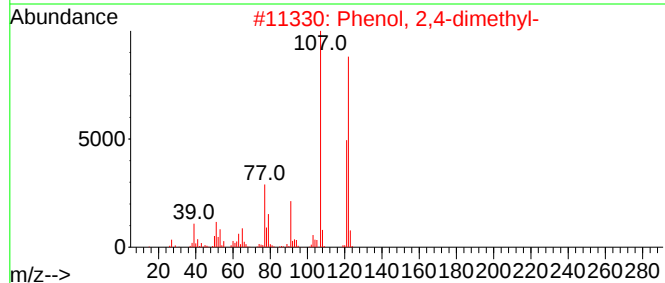
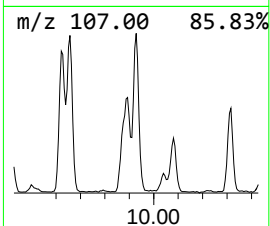
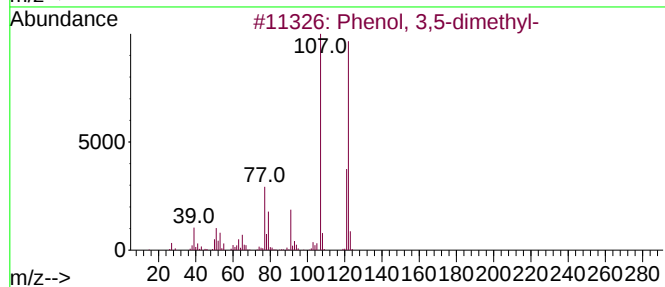
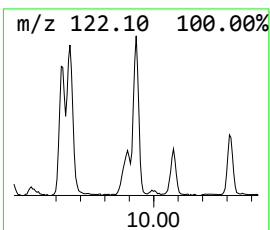
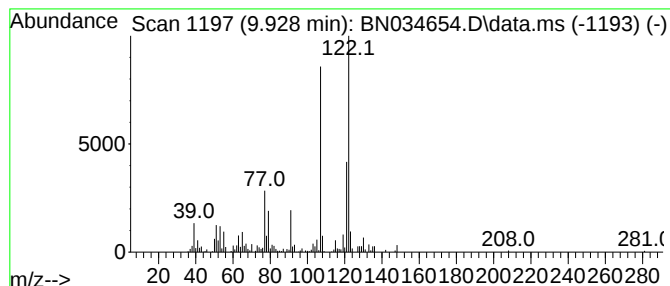
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenol, 3,5-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.928	2.31 ng/ul	304951	Naphthalene-d8	10.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	95
2			Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	93
3			Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	93
4			Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	90
5			Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
Acq On : 18 Oct 2024 17:56
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Misc :
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ClientSampleId :
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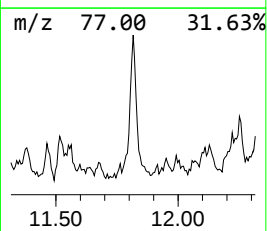
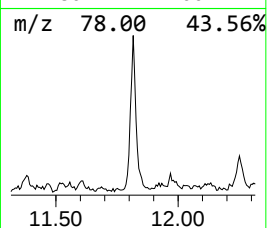
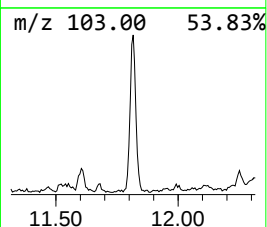
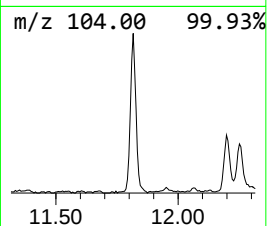
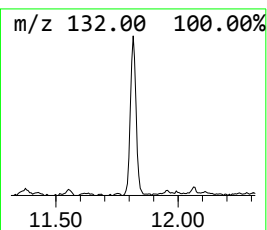
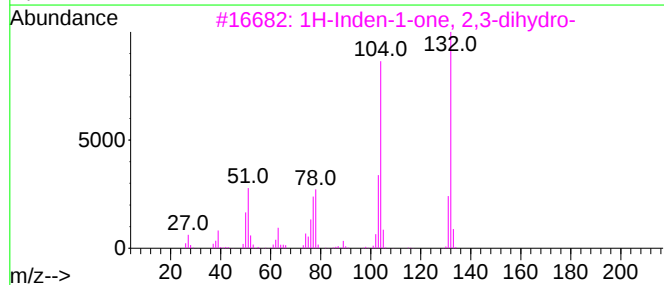
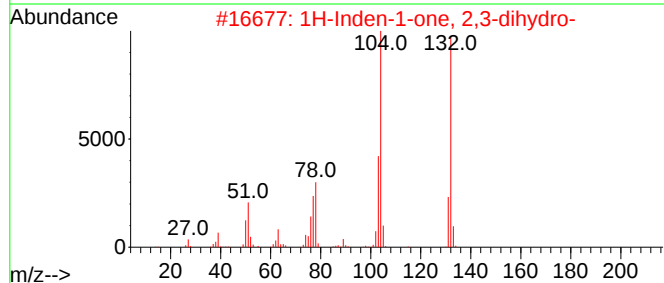
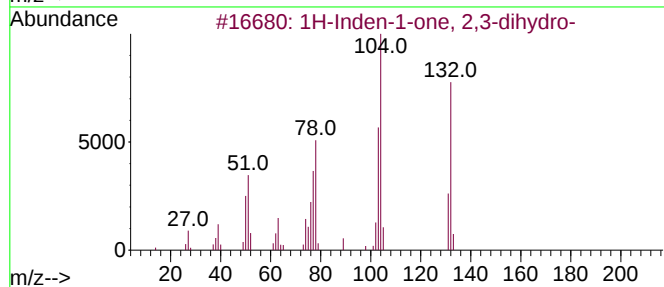
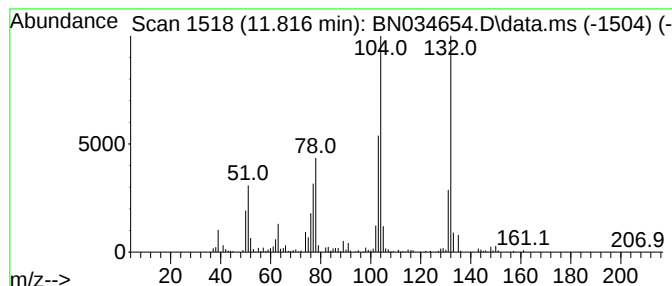
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.816	3.03 ng/ul	400390	Naphthalene-d8	10.440

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	94
2		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	94
3		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	91
4		Styrene	104	C8H8	000100-42-5	70
5		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
Acq On : 18 Oct 2024 17:56
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Sample : P4380-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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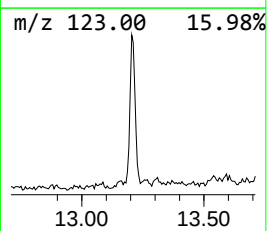
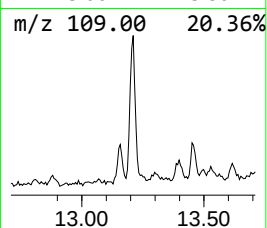
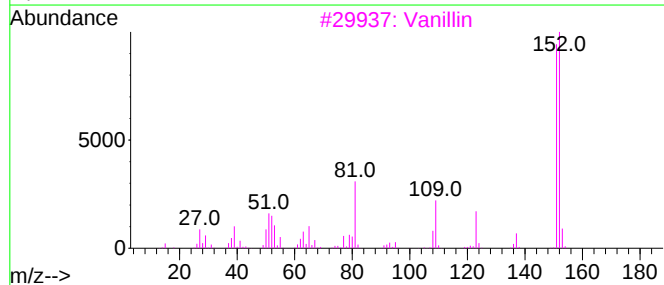
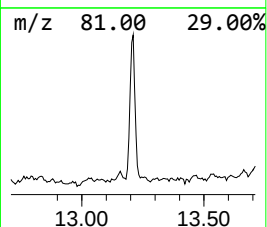
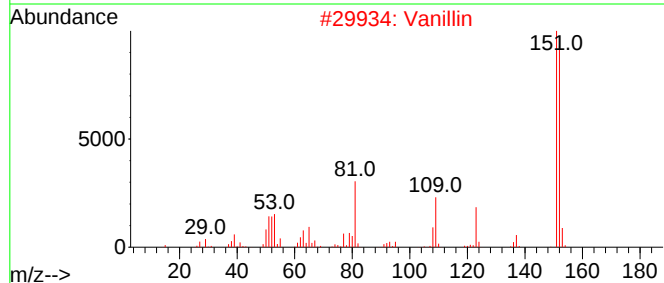
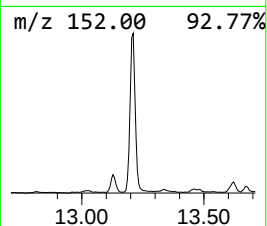
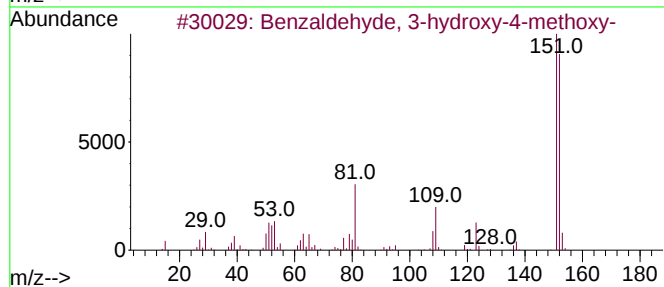
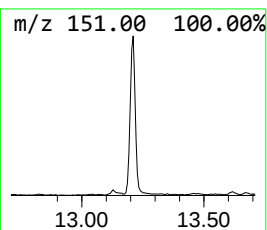
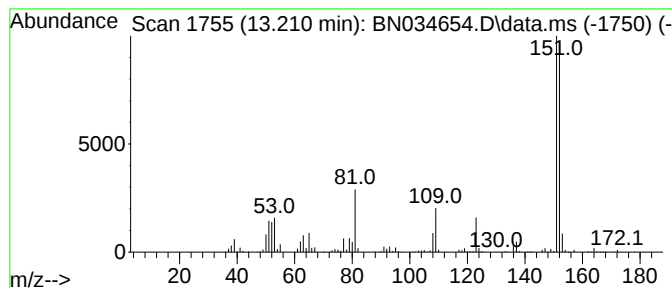
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzaldehyde, 3-hydroxy-4-m... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.210	3.19 ng/ul	464033	Acenaphthene-d10	14.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	97
2			Vanillin	152	C8H8O3	000121-33-5	97
3			Vanillin	152	C8H8O3	000121-33-5	97
4			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96
5			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
Acq On : 18 Oct 2024 17:56
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Sample : P4380-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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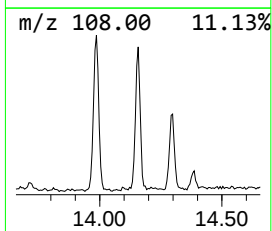
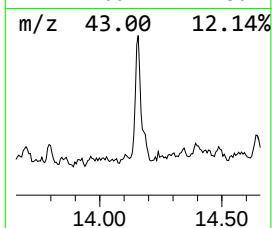
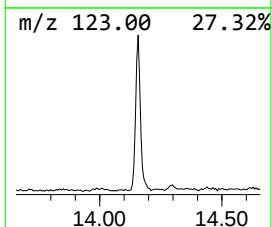
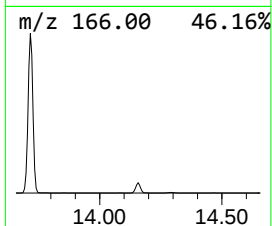
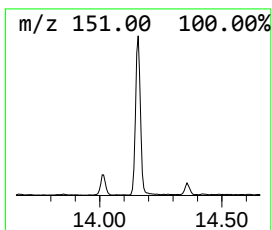
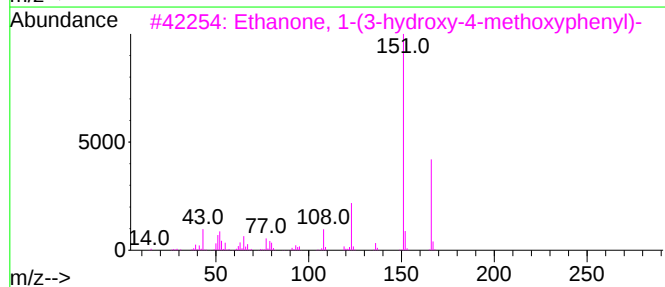
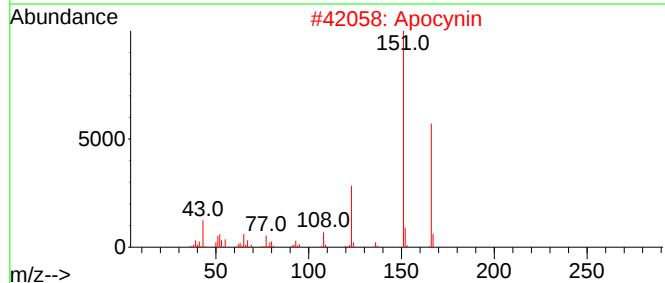
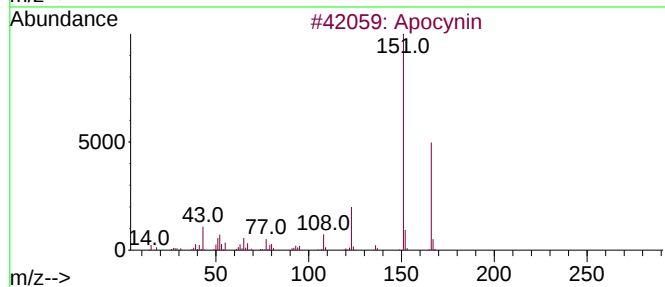
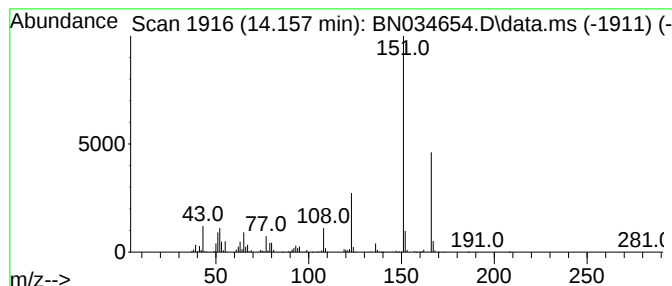
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Apocynin Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.157	5.21 ng/ul	758587	Acenaphthene-d10	14.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Apocynin			166	C9H10O3	000498-02-2	95
2	Apocynin			166	C9H10O3	000498-02-2	95
3	Ethanone, 1-(3-hydroxy-4-methoxy...			166	C9H10O3	006100-74-9	94
4	Apocynin			166	C9H10O3	000498-02-2	94
5	Ethanone, 1-(3-hydroxy-4-methoxy...			166	C9H10O3	006100-74-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
Acq On : 18 Oct 2024 17:56
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Misc :
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Instrument :
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ClientSampleId :
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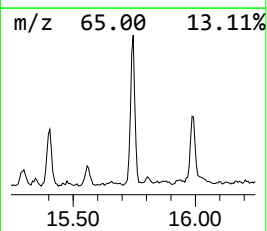
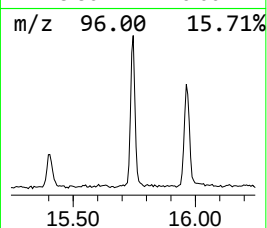
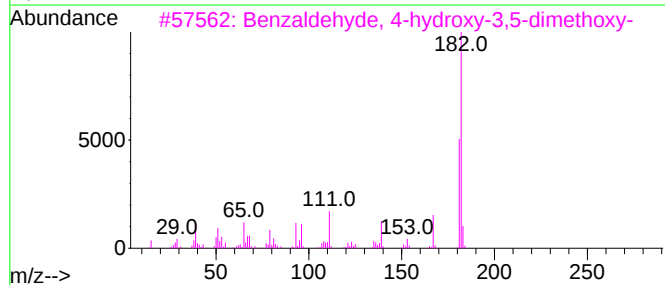
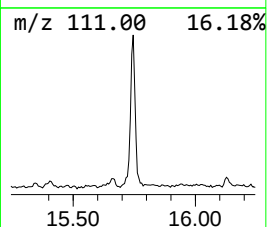
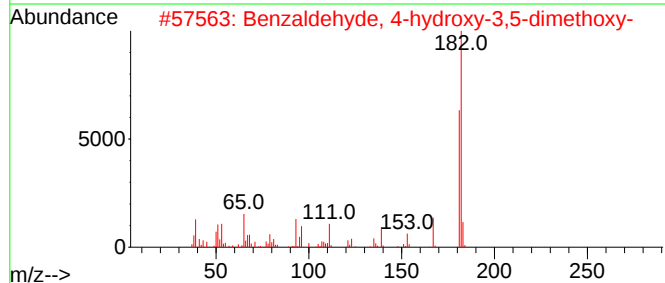
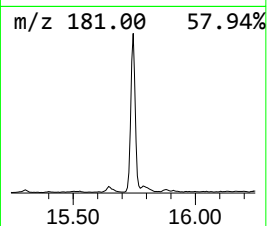
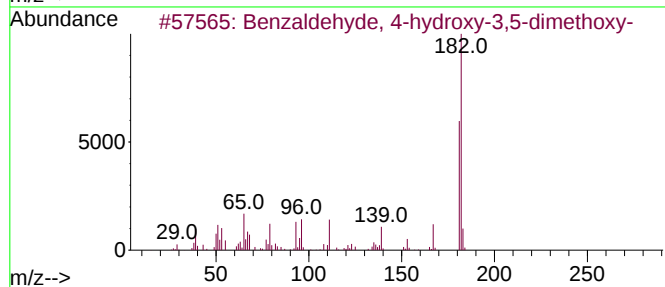
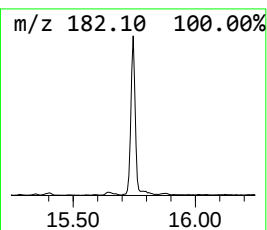
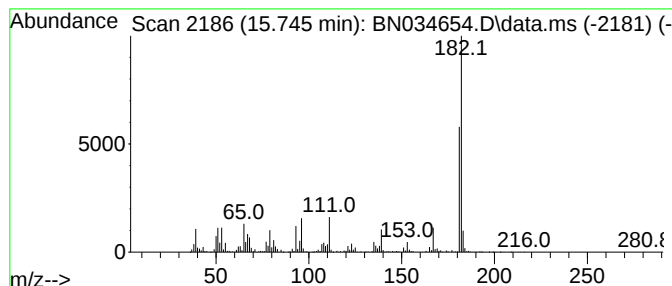
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzaldehyde, 4-hydroxy-3,5... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.745	8.48 ng/ul	1131840	Phenanthrene-d10	17.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	98
2			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	97
3			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	96
4			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	96
5			Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
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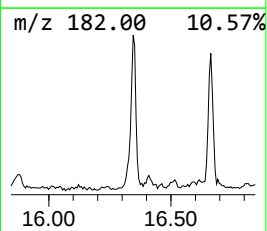
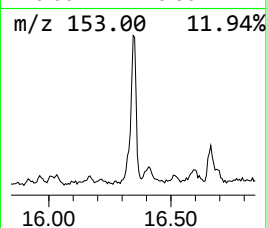
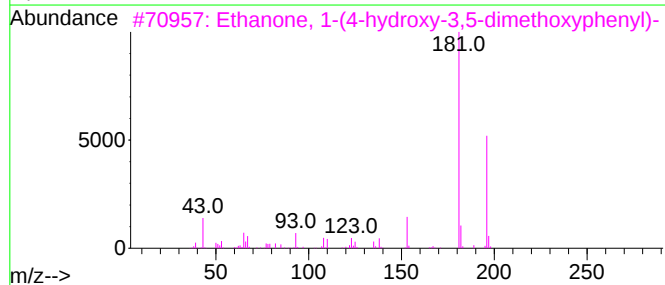
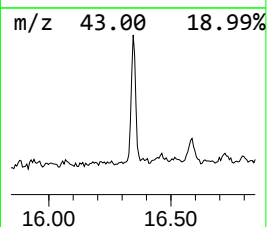
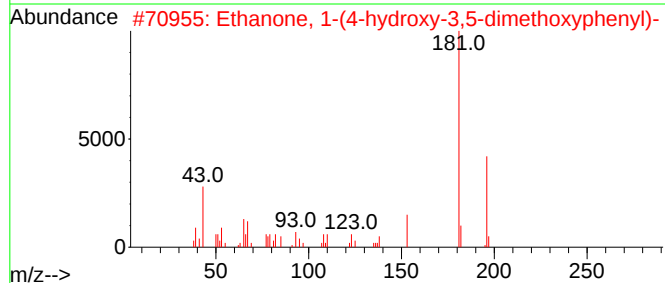
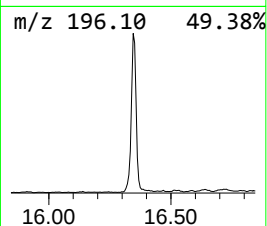
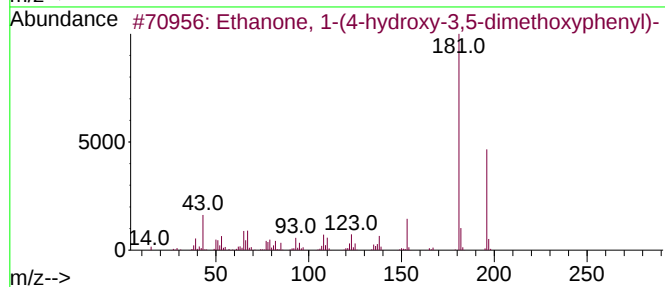
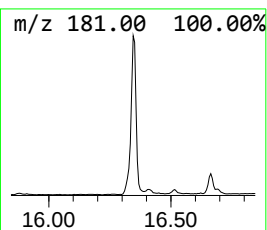
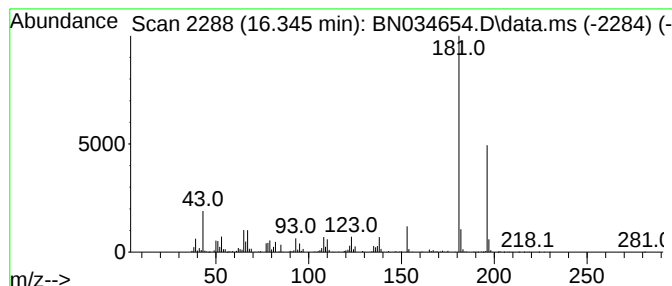
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Ethanone, 1-(4-hydroxy-3,5-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.345	5.87 ng/ul	783706	Phenanthrene-d10	17.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	98
2			Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	97
3			Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	97
4			Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	93
5			4,5-Dimethoxy-2-hydroxyacetophenone	196	C10H12O4	020628-06-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
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ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
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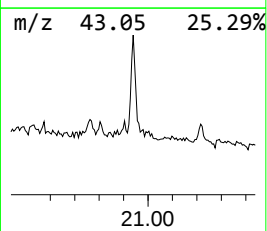
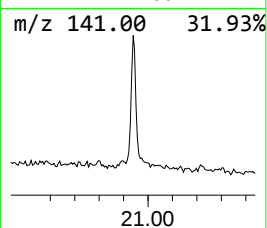
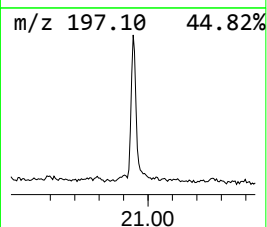
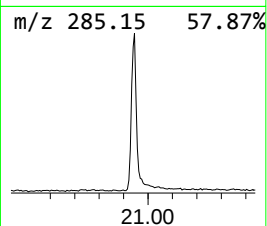
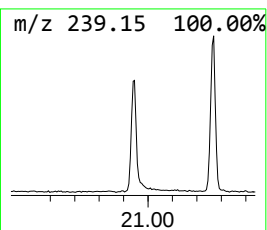
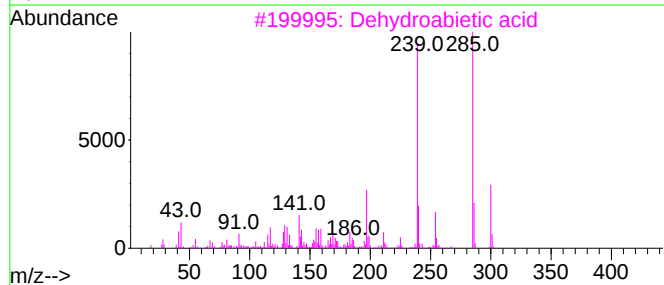
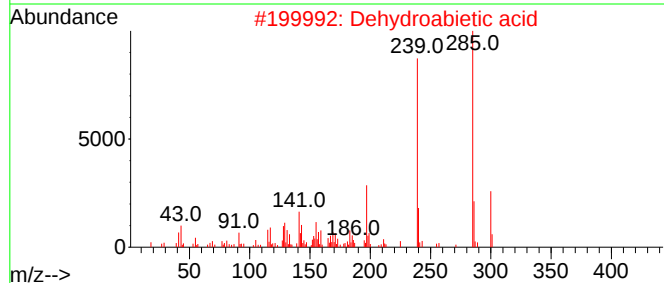
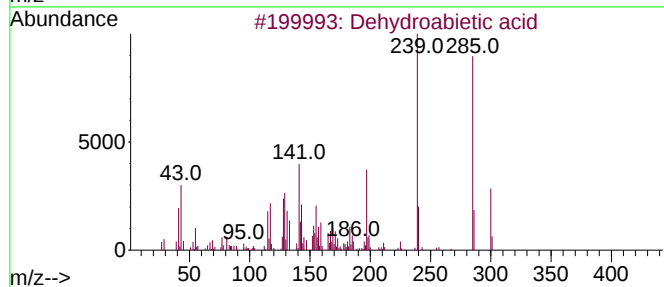
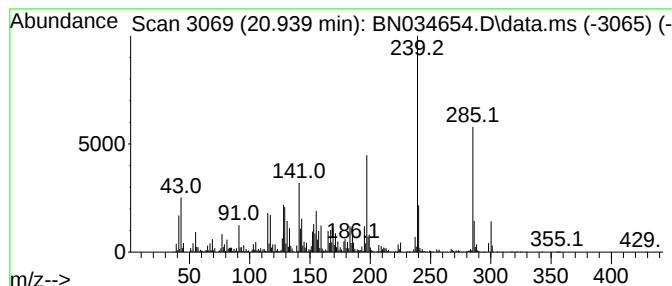
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Dehydroabiatic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.939	9.06 ng/ul	1318350	Chrysene-d12	21.262

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabiatic acid	300	C20H28O2	001740-19-8	99
2			Dehydroabiatic acid	300	C20H28O2	001740-19-8	96
3			Dehydroabiatic acid	300	C20H28O2	001740-19-8	93
4			Dehydroabiatic acid	300	C20H28O2	001740-19-8	90
5			1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034654.D
Acq On : 18 Oct 2024 17:56
Operator : RC/JU
Sample : P4380-16
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
E27J1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.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
Butanoic acid, ...	4.581	2.5	ng/ul	250869	1	7.663	2015210	20.0
Benzaldehyde, 2...	8.157	10.8	ng/ul	1082900	1	7.663	2015210	20.0
Indene	8.193	6.3	ng/ul	637407	1	7.663	2015210	20.0
Phenol, 3,5-dim...	9.928	2.3	ng/ul	304951	2	10.440	2644000	20.0
1H-Inden-1-one,...	11.816	3.0	ng/ul	400390	2	10.440	2644000	20.0
Benzaldehyde, 3...	13.210	3.2	ng/ul	464033	3	14.292	2911540	20.0
Apocynin	14.157	5.2	ng/ul	758587	3	14.292	2911540	20.0
Benzaldehyde, 4...	15.745	8.5	ng/ul	1131840	4	17.039	2670280	20.0
Ethanone, 1-(4-...	16.345	5.9	ng/ul	783706	4	17.039	2670280	20.0
Dehydroabietic ...	20.939	9.1	ng/ul	1318350	5	21.262	2910940	20.0