

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
 Data File : BN034669.D
 Acq On : 19 Oct 2024 04:24
 Operator : RC/JU
 Sample : P4429-14
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4EB8

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: BN034669.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.964	8	13	35	rVB	11077539	17483703	100.00%	14.945%
2	3.211	51	55	63	rVB	83030	112404	0.64%	0.096%
3	3.605	118	122	139	rBV	467274	728565	4.17%	0.623%
4	3.928	173	177	185	rVB	29975	48251	0.28%	0.041%
5	4.205	219	224	228	rBV	28226	38131	0.22%	0.033%
6	4.281	233	237	243	rVB2	42705	60413	0.35%	0.052%
7	4.734	305	314	322	rVB4	24119	56919	0.33%	0.049%
8	4.828	322	330	335	rBV2	24062	40834	0.23%	0.035%
9	5.040	358	366	372	rVB	31859	47202	0.27%	0.040%
10	5.205	390	394	398	rVV	35411	51500	0.29%	0.044%
11	5.734	476	484	487	rBV	92792	150231	0.86%	0.128%
12	5.775	487	491	498	rVB	646964	915069	5.23%	0.782%
13	6.258	568	573	578	rVV4	24140	47707	0.27%	0.041%
14	6.340	578	587	598	rVB2	311397	542703	3.10%	0.464%
15	6.681	640	645	652	rVB	60568	102451	0.59%	0.088%
16	6.840	657	672	680	rBV	542319	999940	5.72%	0.855%
17	7.005	688	700	707	rBV	1583681	2437621	13.94%	2.084%
18	7.199	725	733	742	rBV	1766831	2718105	15.55%	2.323%
19	7.458	769	777	786	rVB	126749	200179	1.14%	0.171%
20	7.663	804	812	818	rBV	1413144	2162028	12.37%	1.848%
21	7.740	820	825	834	rVV	375646	555240	3.18%	0.475%
22	7.822	834	839	846	rVV2	40170	68572	0.39%	0.059%
23	7.899	846	852	860	rVV	294779	474614	2.71%	0.406%
24	7.969	860	864	871	rVB	33263	57654	0.33%	0.049%
25	8.240	905	910	913	rVV2	32989	52270	0.30%	0.045%
26	8.281	913	917	924	rVV5	39443	78471	0.45%	0.067%
27	8.369	924	932	943	rVV	1541969	2418706	13.83%	2.067%
28	8.469	943	949	957	rVB2	227399	425899	2.44%	0.364%
29	8.540	957	961	965	rBV3	28309	42999	0.25%	0.037%
30	8.752	991	997	1001	rBV2	206381	355317	2.03%	0.304%
31	8.805	1001	1006	1019	rVV	1761537	2979238	17.04%	2.547%
32	8.916	1019	1025	1033	rVV	620650	962752	5.51%	0.823%
33	8.993	1033	1038	1047	rVV	693592	1006415	5.76%	0.860%
34	9.193	1061	1072	1081	rVV	69423	136988	0.78%	0.117%
35	9.287	1083	1088	1098	rVV	49001	80718	0.46%	0.069%
36	9.381	1098	1104	1112	rVB4	42954	79961	0.46%	0.068%

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

37	9.522	1121	1128	1135	rBV	1382655	2263881	12.95%	1.935%
38	9.575	1135	1137	1144	rVV4	40710	81846	0.47%	0.070%
39	9.716	1148	1161	1167	rVV2	489322	1083679	6.20%	0.926%
40	9.775	1168	1171	1179	rVV	51586	93650	0.54%	0.080%

41	9.922	1190	1196	1200	rVV	611010	955074	5.46%	0.816%
42	9.957	1200	1202	1211	rVV	309928	416113	2.38%	0.356%
43	10.063	1211	1220	1231	rVV	2152426	3513927	20.10%	3.004%
44	10.234	1242	1249	1255	rBV2	129393	209647	1.20%	0.179%
45	10.310	1257	1262	1266	rBV2	44808	78126	0.45%	0.067%

46	10.352	1266	1269	1277	rVB5	43765	79896	0.46%	0.068%
47	10.440	1277	1284	1289	rBV	1879302	3156129	18.05%	2.698%
48	10.481	1289	1291	1300	rVB2	184549	342108	1.96%	0.292%
49	10.569	1300	1306	1314	rBV	1960301	3216626	18.40%	2.750%
50	10.740	1329	1335	1342	rBV	383707	616582	3.53%	0.527%

51	10.840	1347	1352	1361	rVB	103366	179382	1.03%	0.153%
52	10.999	1371	1379	1384	rVB	107065	169079	0.97%	0.145%
53	11.063	1384	1390	1402	rVB2	142105	339338	1.94%	0.290%
54	11.187	1406	1411	1414	rBV3	20979	36967	0.21%	0.032%
55	11.246	1414	1421	1432	rVB6	68050	141275	0.81%	0.121%

56	11.616	1480	1484	1489	rVV	22163	31382	0.18%	0.027%
57	11.781	1504	1512	1517	rBV3	66088	134054	0.77%	0.115%
58	11.840	1517	1522	1526	rVV	40054	62872	0.36%	0.054%
59	12.028	1548	1554	1560	rBV2	44870	77125	0.44%	0.066%
60	12.428	1615	1622	1627	rBV3	17731	33802	0.19%	0.029%

61	12.663	1656	1662	1672	rVV	403113	602485	3.45%	0.515%
62	12.746	1672	1676	1685	rVB	40017	65904	0.38%	0.056%
63	12.863	1691	1696	1700	rVB	42045	56010	0.32%	0.048%
64	12.922	1700	1706	1717	rVB	480901	691032	3.95%	0.591%
65	13.145	1737	1744	1749	rBV	34661	52869	0.30%	0.045%

66	13.210	1749	1755	1763	rBV	979312	1367263	7.82%	1.169%
67	13.422	1787	1791	1799	rVB	47961	71087	0.41%	0.061%
68	13.545	1805	1812	1816	rBV	143318	237129	1.36%	0.203%
69	13.716	1833	1841	1847	rVV	3484387	4644626	26.57%	3.970%
70	13.757	1847	1848	1852	rVV	31026	40578	0.23%	0.035%

71	13.851	1860	1864	1876	rVB2	74924	118975	0.68%	0.102%
72	13.987	1876	1887	1894	rBV	4200614	6010852	34.38%	5.138%
73	14.157	1909	1916	1922	rBV3	46138	96605	0.55%	0.083%
74	14.298	1932	1940	1952	rBV	2486415	3916653	22.40%	3.348%
75	14.492	1964	1973	1980	rVV	376015	568478	3.25%	0.486%

76	14.563	1980	1985	1991	rVV4	40222	78738	0.45%	0.067%
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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
 Title : SVOA CALIBRATION

77	14.722	2007	2012	2018	rVV7	27578	50039	0.29%	0.043%
78	14.998	2053	2059	2065	rBV3	30414	62246	0.36%	0.053%
79	15.051	2065	2068	2073	rVV	46454	59029	0.34%	0.050%
80	15.134	2075	2082	2087	rVB2	133649	175866	1.01%	0.150%
81	15.292	2102	2109	2117	rVV	4807659	6698025	38.31%	5.725%
82	15.404	2121	2128	2134	rVV	969576	1314397	7.52%	1.124%
83	15.457	2135	2137	2142	rVV3	35475	50273	0.29%	0.043%
84	15.534	2145	2150	2156	rVV2	31869	57112	0.33%	0.049%
85	15.675	2167	2174	2180	rVV2	63947	106554	0.61%	0.091%
86	15.987	2222	2227	2235	rVB	44343	64853	0.37%	0.055%
87	17.039	2397	2406	2415	rBV	2675112	3726164	21.31%	3.185%
88	17.134	2415	2422	2435	rBV2	4601658	6650229	38.04%	5.685%
89	17.351	2454	2459	2464	rBV	28533	37571	0.21%	0.032%
90	17.733	2518	2524	2534	rBV	3377713	4101496	23.46%	3.506%
91	18.051	2571	2578	2583	rBV3	33195	67110	0.38%	0.057%
92	18.169	2593	2598	2601	rBV	57827	79829	0.46%	0.068%
93	18.204	2601	2604	2609	rVB3	35304	49417	0.28%	0.042%
94	18.998	2734	2739	2743	rBV	56837	79393	0.45%	0.068%
95	19.063	2746	2750	2754	rVV	65114	78814	0.45%	0.067%
96	19.428	2803	2812	2819	rBV	4704179	6447450	36.88%	5.511%
97	21.257	3118	3123	3132	rBV2	2056031	2912456	16.66%	2.490%
98	22.657	3357	3361	3367	rBV2	96728	162576	0.93%	0.139%
99	23.951	3572	3581	3598	rVB	2579489	6078613	34.77%	5.196%
100	24.145	3606	3614	3626	rVB	1363195	3227380	18.46%	2.759%

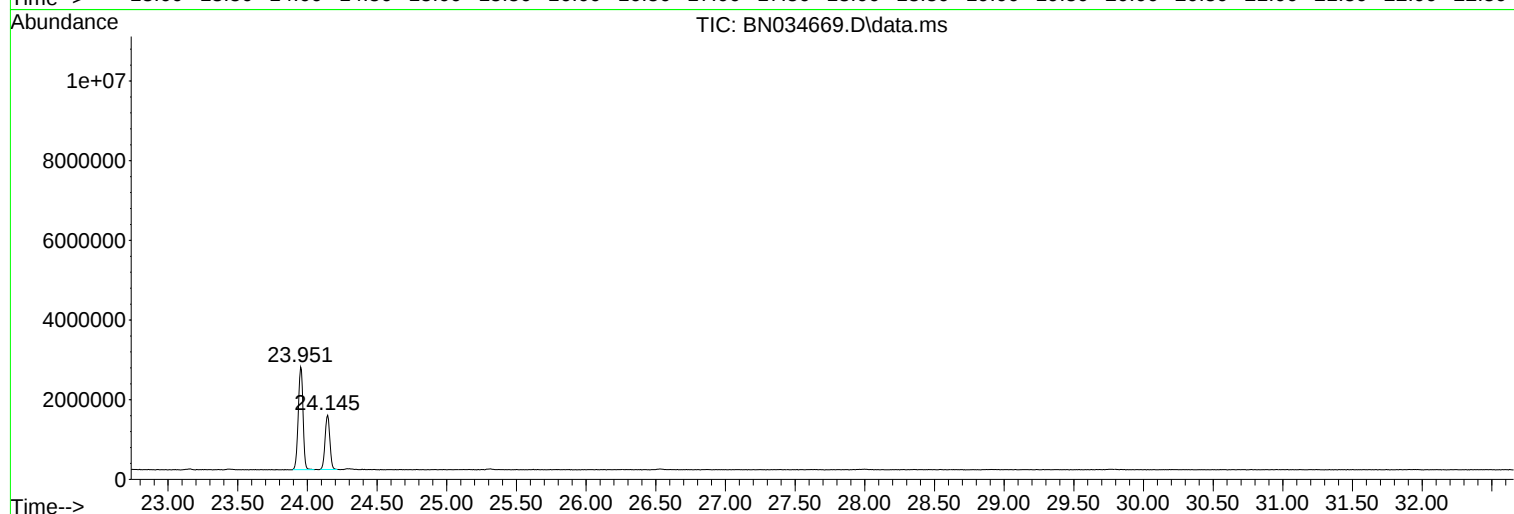
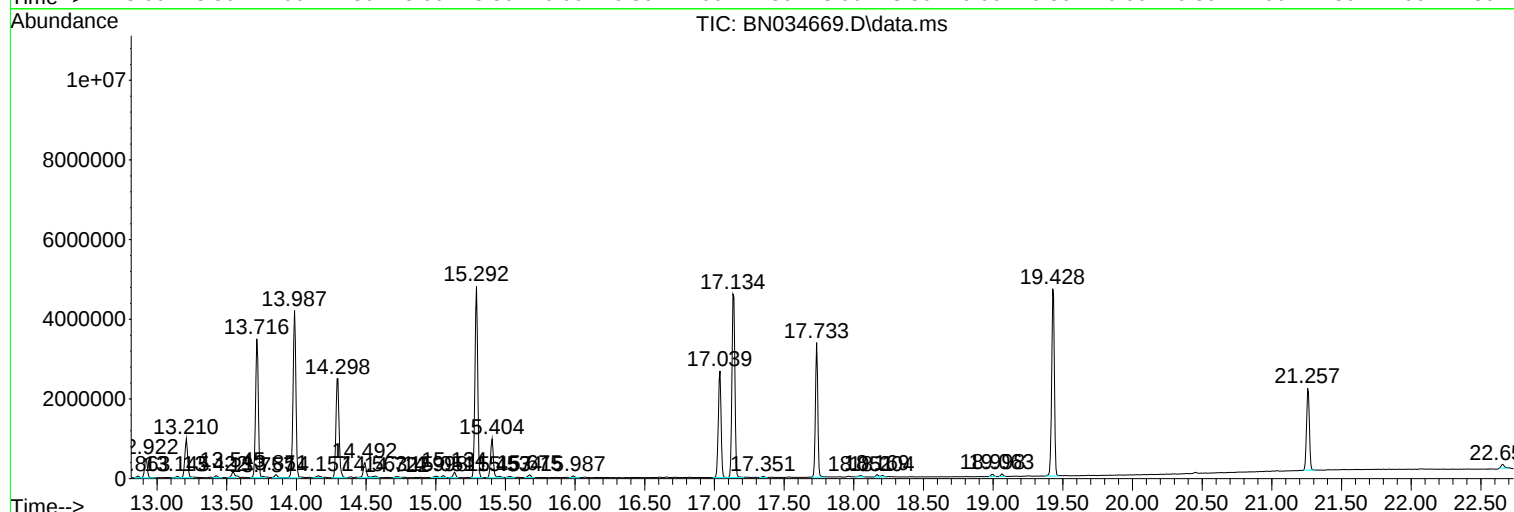
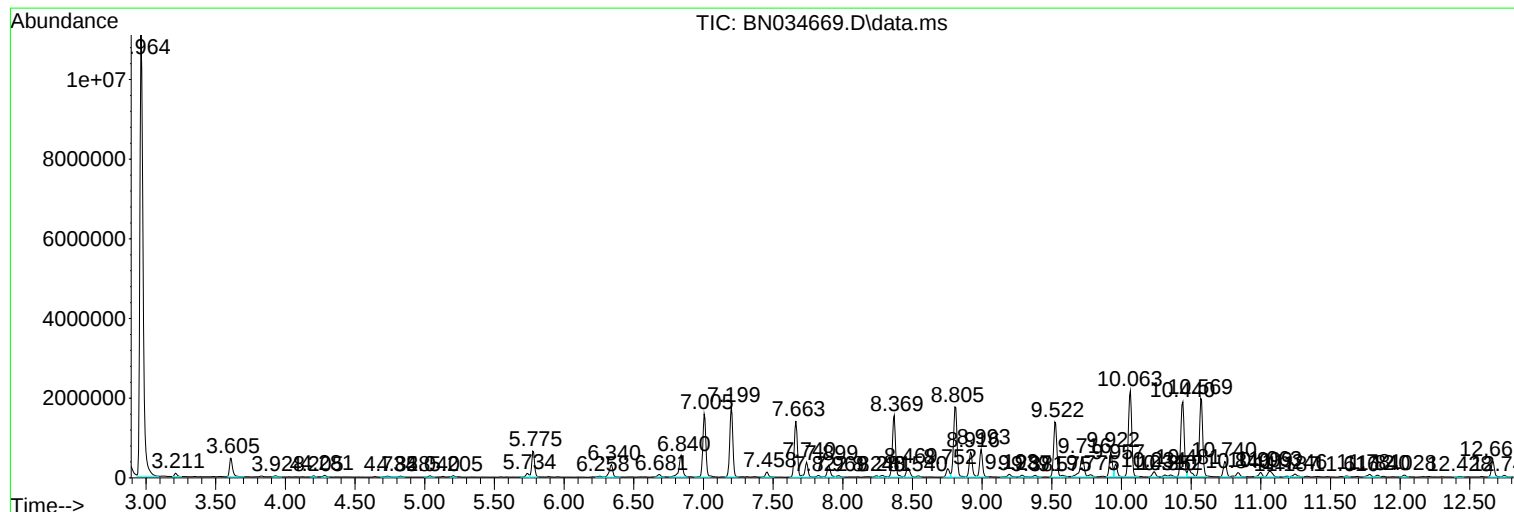
Sum of corrected areas: 116988506

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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



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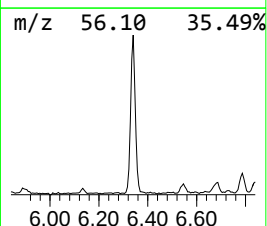
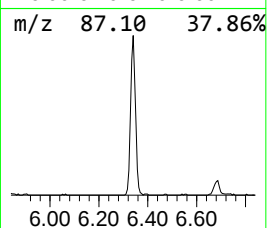
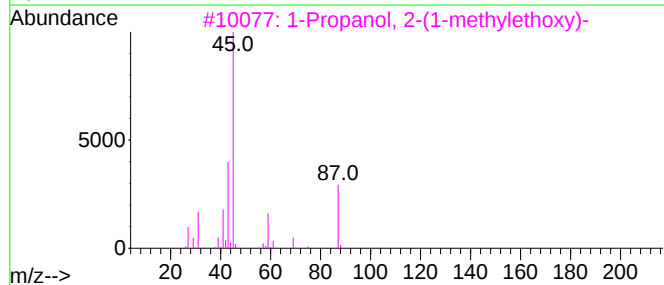
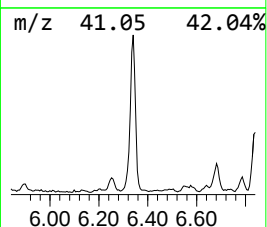
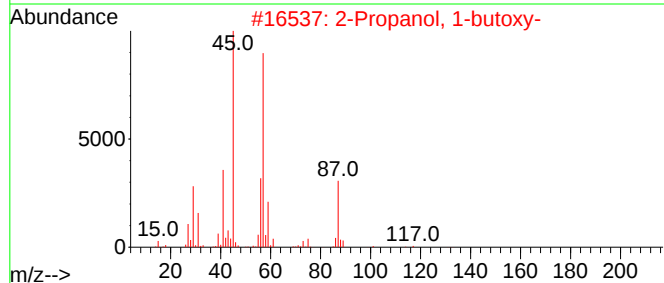
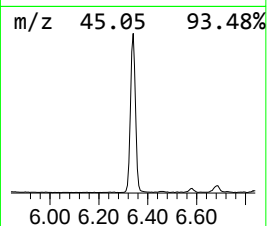
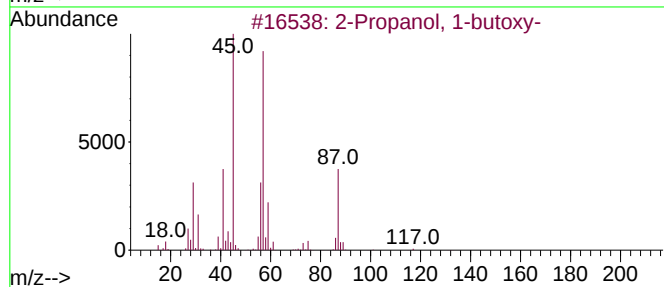
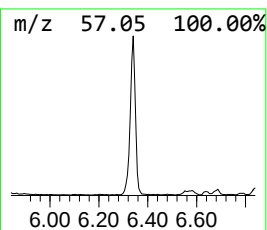
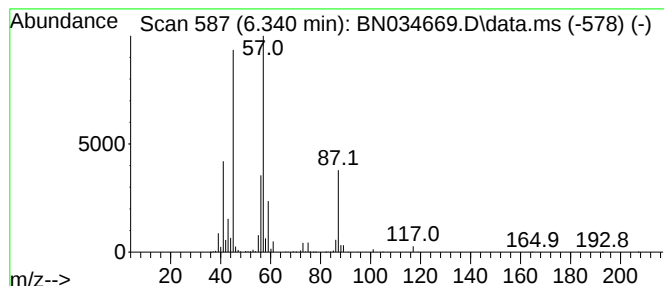
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1-butoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.340	5.02 ng/ul	542703	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-butoxy-			132	C7H16O2	005131-66-8	90
2	2-Propanol, 1-butoxy-			132	C7H16O2	005131-66-8	90
3	1-Propanol, 2-(1-methylethoxy)-			118	C6H14O2	003944-37-4	53
4	1-Propanol, 2-(1-methylethoxy)-			118	C6H14O2	003944-37-4	53
5	3,3-Dimethylbutane-2-ol			102	C6H14O	000464-07-3	53



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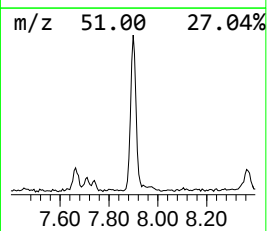
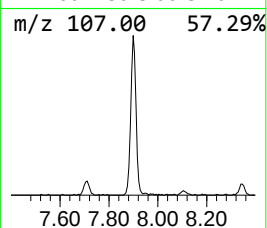
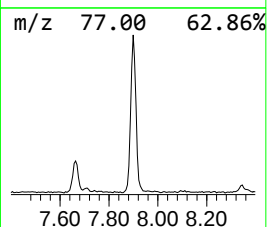
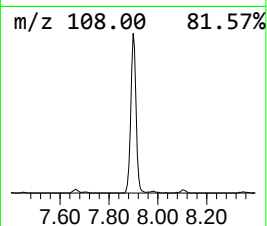
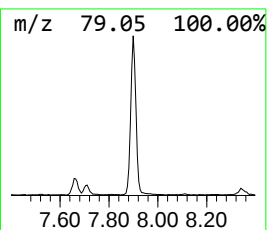
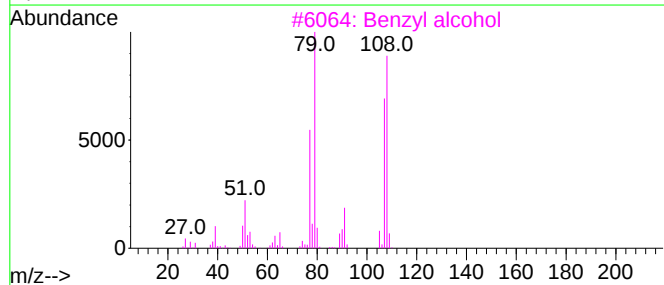
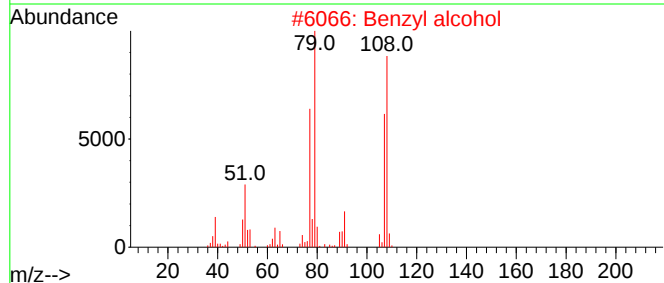
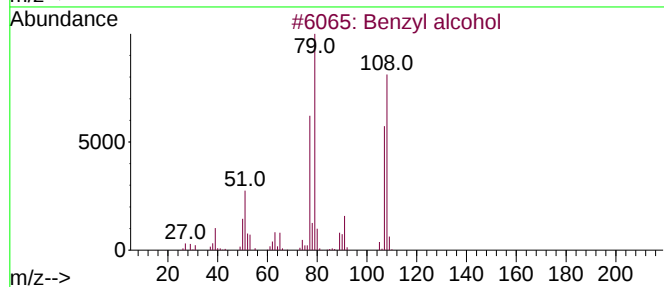
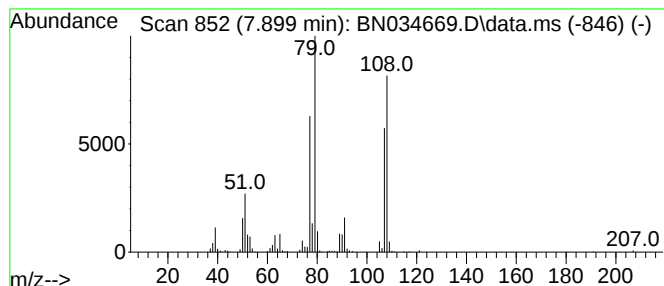
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzyl alcohol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.899	4.39 ng/ul	474614	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl alcohol	108	C7H8O	000100-51-6	98
2			Benzyl alcohol	108	C7H8O	000100-51-6	96
3			Benzyl alcohol	108	C7H8O	000100-51-6	94
4			Benzyl alcohol	108	C7H8O	000100-51-6	94
5			N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91



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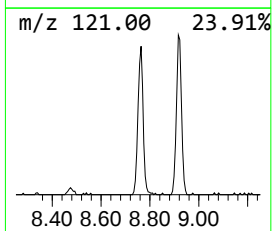
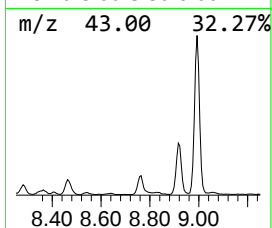
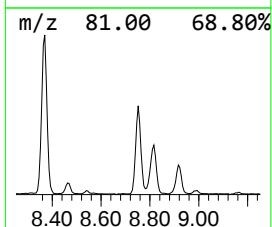
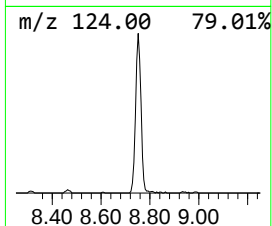
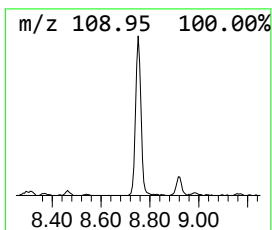
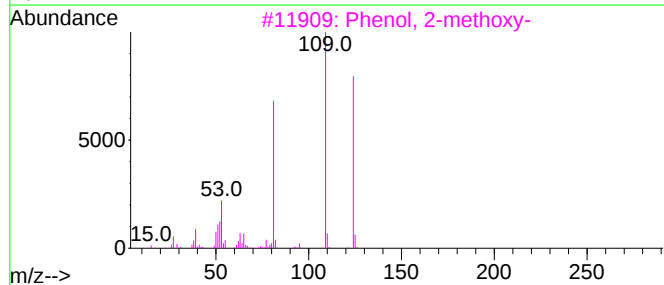
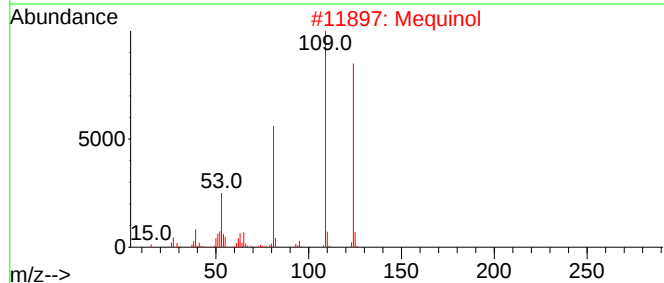
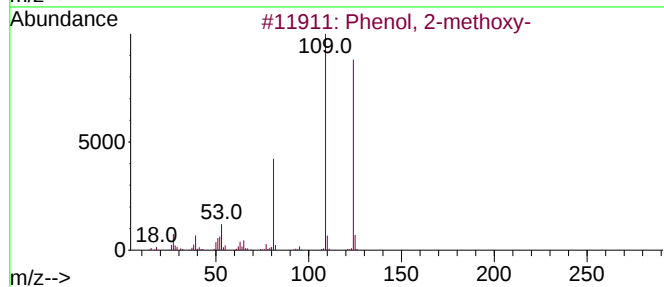
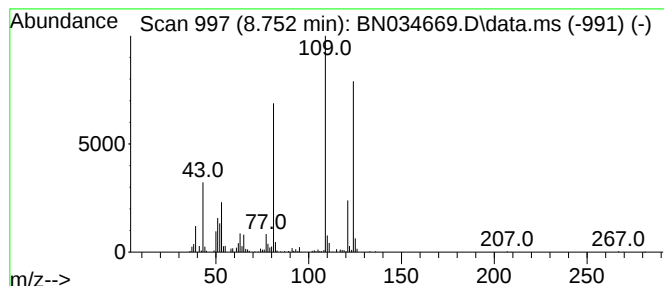
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 2-methoxy- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.752	3.29 ng/ul	355317	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	95
2			Mequinol	124	C7H8O2	000150-76-5	93
3			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	93
4			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
5			Mequinol	124	C7H8O2	000150-76-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034669.D
Acq On : 19 Oct 2024 04:24
Operator : RC/JU
Sample : P4429-14
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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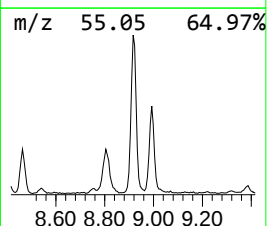
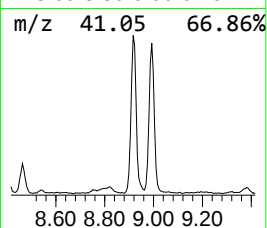
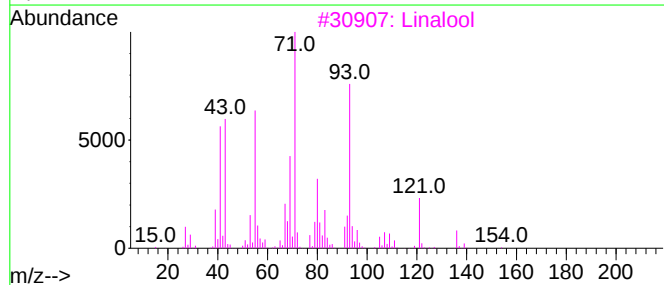
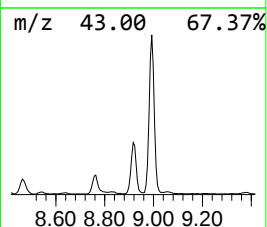
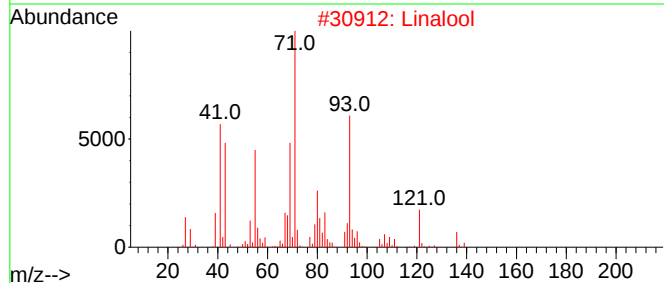
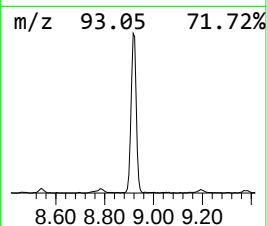
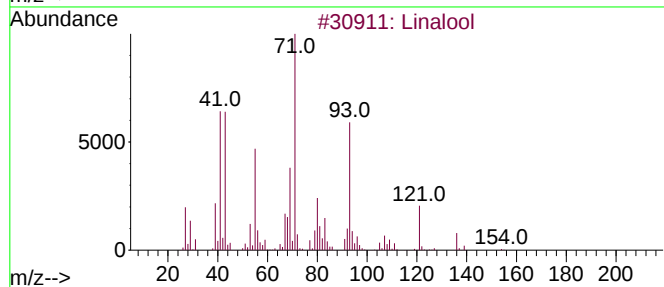
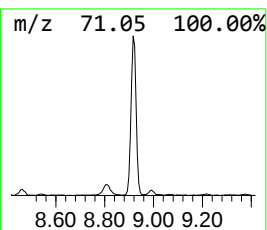
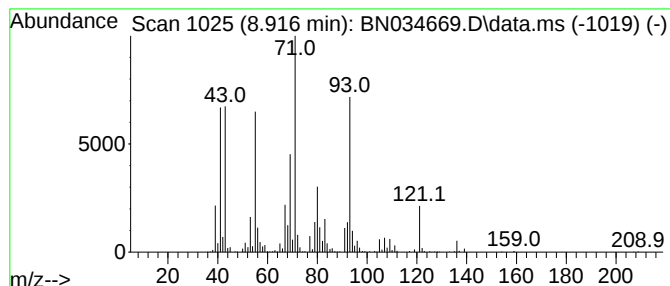
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Linalool Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.916	8.91 ng/ul	962752	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Linalool	154	C10H18O	000078-70-6	87
2			Linalool	154	C10H18O	000078-70-6	80
3			Linalool	154	C10H18O	000078-70-6	58
4			1,5-Dimethyl-1-vinyl-4-hexenyl b...	224	C14H24O2	000078-36-4	46
5			1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034669.D
Acq On : 19 Oct 2024 04:24
Operator : RC/JU
Sample : P4429-14
Misc :
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Instrument :
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ClientSampleId :
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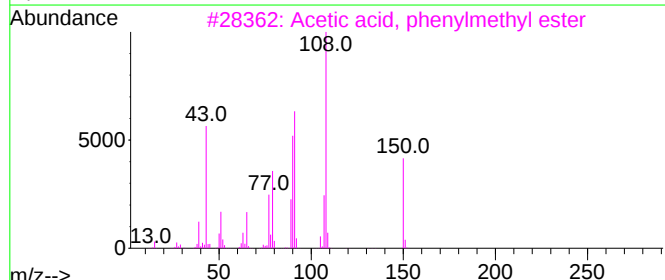
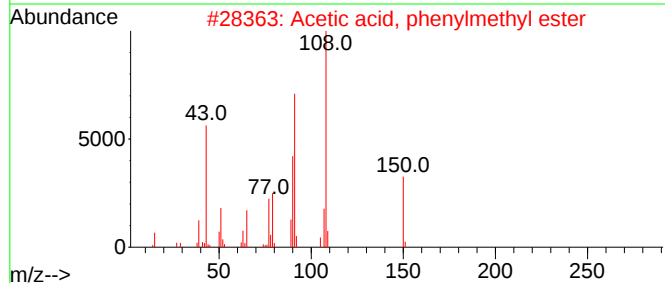
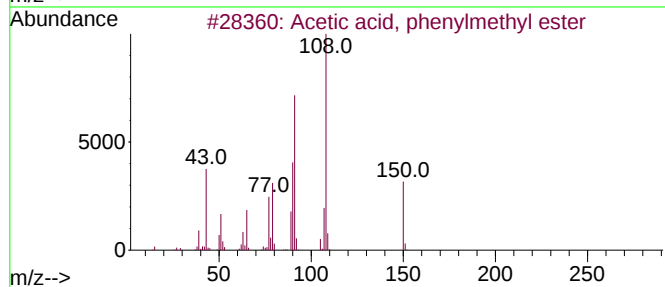
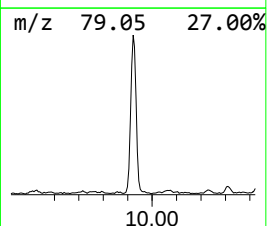
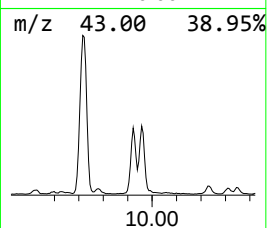
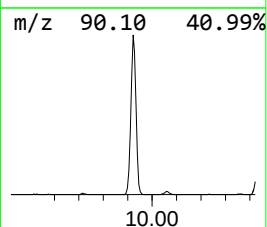
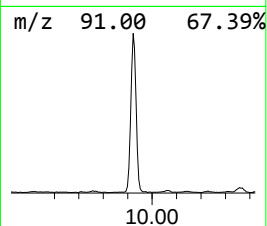
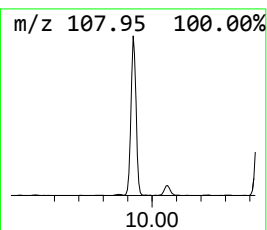
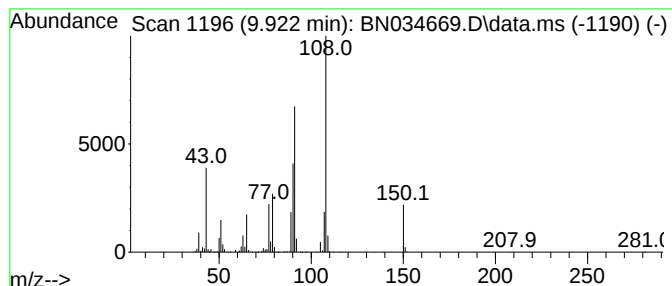
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Acetic acid, phenylmethyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.922	6.05 ng/ul	955074	Naphthalene-d8	10.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	98
2			Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	95
3			Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	95
4			Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	91
5			Acetic acid, phenylmethyl ester	150	C9H10O2	000140-11-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034669.D
Acq On : 19 Oct 2024 04:24
Operator : RC/JU
Sample : P4429-14
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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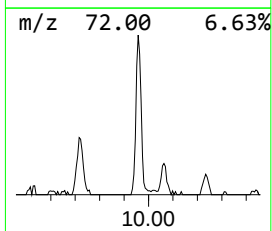
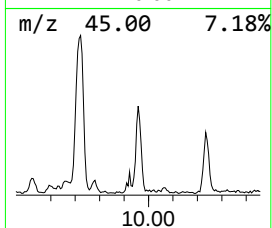
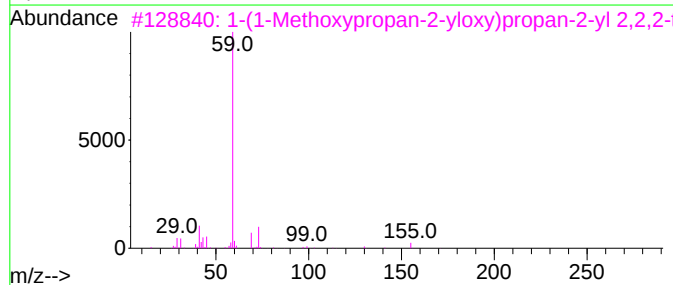
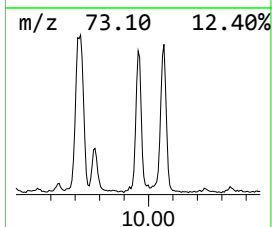
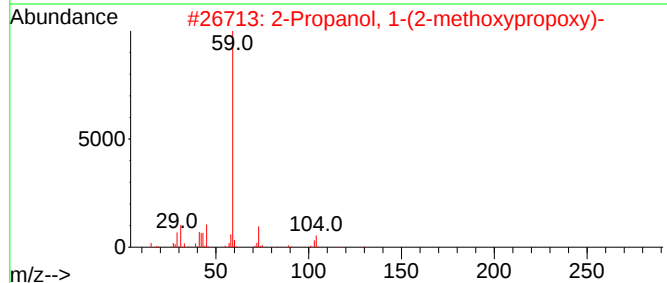
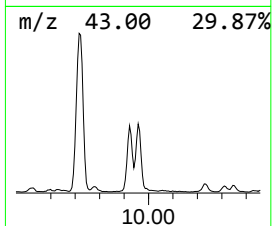
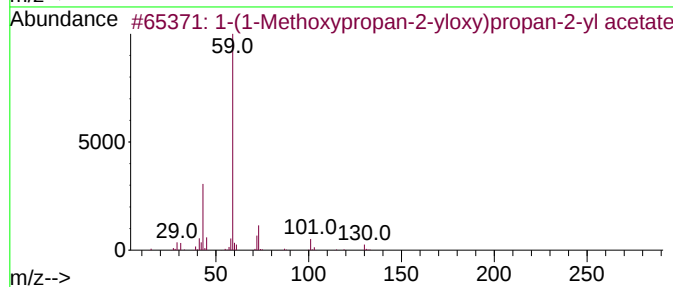
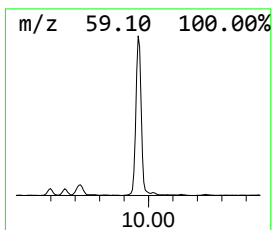
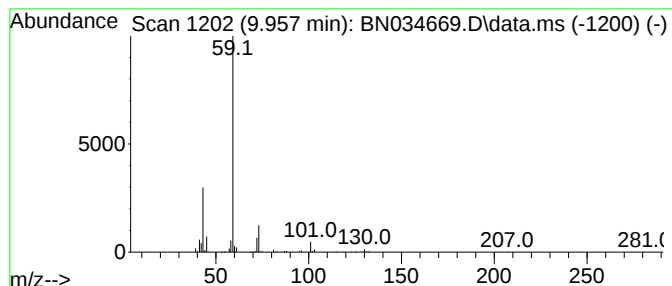
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1-(1-Methoxypropan-2-yloxy)... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.957	2.64 ng/ul	416113	Naphthalene-d8	10.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4	1000367-08-6	90		
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	72		
3	1-(1-Methoxypropan-2-yloxy)propa...	244	C9H15F3O4	1010367-08-7	56		
4	Pentane, 2-methoxy-	102	C6H14O	006795-88-6	53		
5	3-Hexanol	102	C6H14O	000623-37-0	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034669.D
Acq On : 19 Oct 2024 04:24
Operator : RC/JU
Sample : P4429-14
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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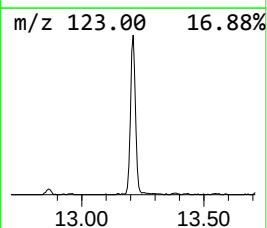
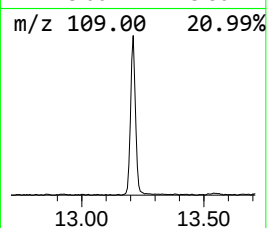
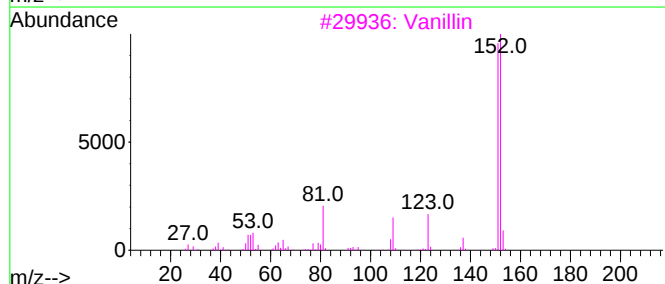
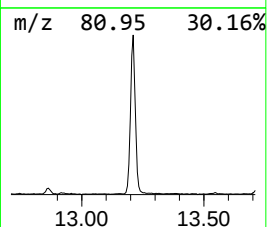
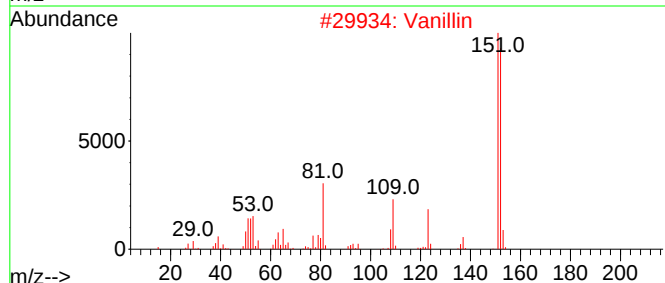
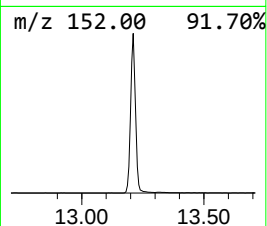
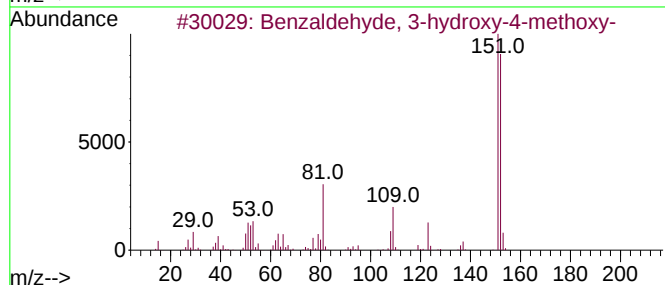
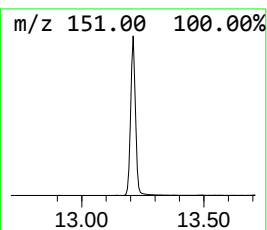
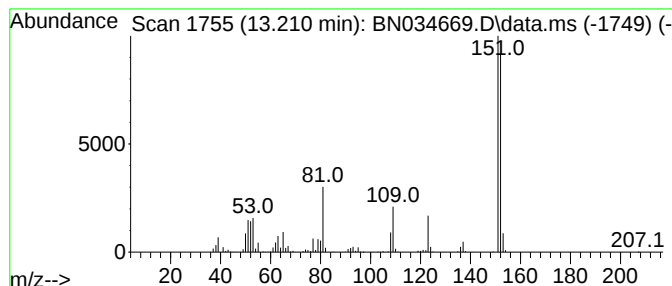
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzaldehyde, 3-hydroxy-4-m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.210	6.98 ng/ul	1367260	Acenaphthene-d10	14.298

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	97
5			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034669.D
Acq On : 19 Oct 2024 04:24
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Misc :
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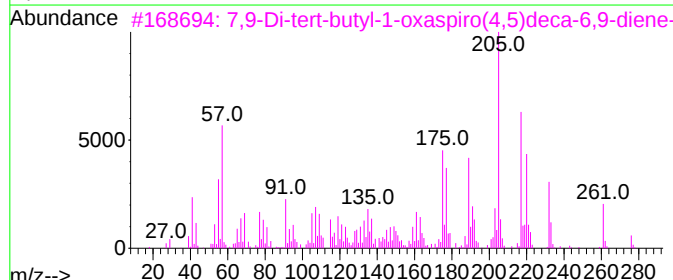
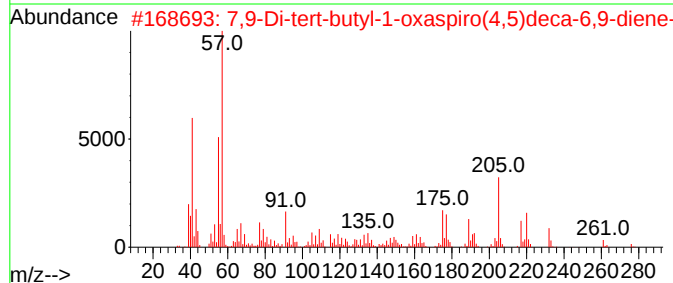
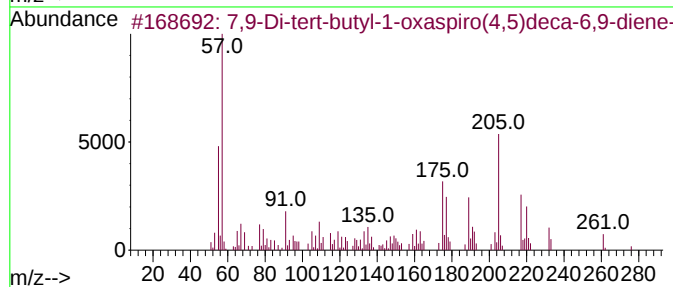
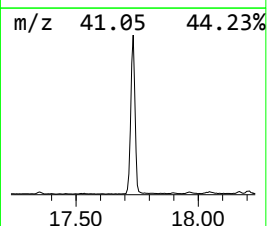
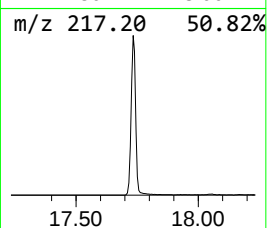
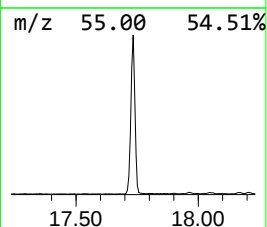
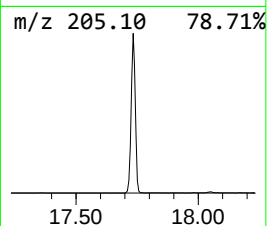
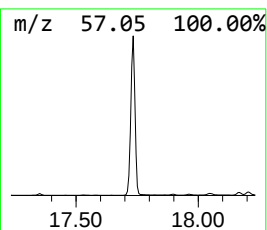
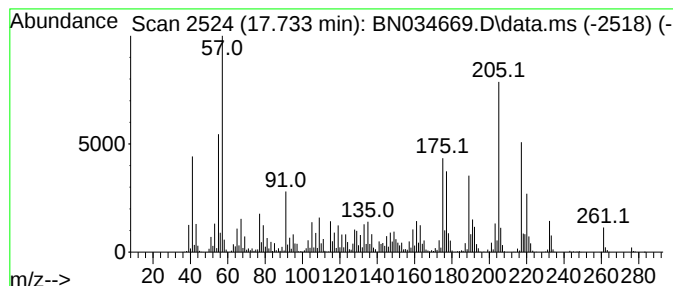
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.733	22.01 ng/ul	4101500	Phenanthrene-d10	17.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	98		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	91		
4	1,3-Pentadiene, 1,1-diphenyl-, (Z)-	220	C17H16	015295-31-5	55		
5	2-Phenazinecarbonitrile	205	C13H7N3	006479-93-2	51		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034669.D
Acq On : 19 Oct 2024 04:24
Operator : RC/JU
Sample : P4429-14
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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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-Propanol, 1-b...	6.340	5.0	ng/ul	542703	1	7.663	2162030	20.0
Benzyl alcohol	7.899	4.4	ng/ul	474614	1	7.663	2162030	20.0
Phenol, 2-methoxy-	8.752	3.3	ng/ul	355317	1	7.663	2162030	20.0
Linalool	8.916	8.9	ng/ul	962752	1	7.663	2162030	20.0
Acetic acid, ph...	9.922	6.0	ng/ul	955074	2	10.440	3156130	20.0
1-(1-Methoxypro...	9.957	2.6	ng/ul	416113	2	10.440	3156130	20.0
Benzaldehyde, 3...	13.210	7.0	ng/ul	1367260	3	14.298	3916650	20.0
7,9-Di-tert-but...	17.733	22.0	ng/ul	4101500	4	17.039	3726160	20.0