

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
 Data File : BN029776.D
 Acq On : 14 Feb 2024 19:04
 Operator : MA/JU
 Sample : P1289-16
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AG5

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-BN020724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN029776.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.322	36	40	45	rBV	36452	43966	0.93%	0.082%
2	3.734	107	110	126	rBV	220700	383406	8.15%	0.717%
3	3.958	145	148	154	rVB5	9317	13646	0.29%	0.026%
4	4.052	161	164	170	rVB	21635	29866	0.64%	0.056%
5	4.993	321	324	331	rVB	15969	25315	0.54%	0.047%
6	5.493	405	409	414	rBV	13972	20853	0.44%	0.039%
7	5.928	477	483	484	rBV5	7087	8692	0.18%	0.016%
8	5.958	484	488	501	rVB	21673	43360	0.92%	0.081%
9	6.440	566	570	575	rVB5	7919	13303	0.28%	0.025%
10	6.522	579	584	589	rVB6	7816	12635	0.27%	0.024%
11	6.869	639	643	648	rBV5	7608	14787	0.31%	0.028%
12	6.975	657	661	666	rBV7	7701	13749	0.29%	0.026%
13	7.040	666	672	682	rVB	179872	304923	6.48%	0.570%
14	7.210	695	701	710	rBV	747394	1207428	25.67%	2.258%
15	7.399	726	733	745	rBV	740896	1170535	24.89%	2.189%
16	7.869	805	813	819	rBV	709108	1107264	23.54%	2.071%
17	7.934	819	824	832	rVB	198633	302429	6.43%	0.566%
18	8.122	850	856	867	rVB2	22788	43192	0.92%	0.081%
19	8.434	905	909	914	rBV7	7048	14558	0.31%	0.027%
20	8.499	916	920	925	rVB2	9609	16487	0.35%	0.031%
21	8.581	927	934	944	rVV	544483	878559	18.68%	1.643%
22	8.663	944	948	949	rVV3	8106	11854	0.25%	0.022%
23	8.699	949	954	960	rVB	43612	68090	1.45%	0.127%
24	8.969	993	1000	1005	rBV	89895	147970	3.15%	0.277%
25	9.034	1005	1011	1019	rVV	879339	1455783	30.95%	2.723%
26	9.146	1027	1030	1035	rVB5	8801	13875	0.30%	0.026%
27	9.204	1035	1040	1047	rVB2	28247	48410	1.03%	0.091%
28	9.428	1073	1078	1084	rVB2	25940	41162	0.88%	0.077%
29	9.522	1089	1094	1097	rBV6	9831	19054	0.41%	0.036%
30	9.587	1101	1105	1110	rVB5	15125	24912	0.53%	0.047%
31	9.751	1125	1133	1145	rBV	672849	1151964	24.49%	2.155%
32	9.910	1155	1160	1162	rBV4	10318	15959	0.34%	0.030%
33	9.993	1169	1174	1182	rVB3	22193	38444	0.82%	0.072%
34	10.151	1197	1201	1204	rBV5	6703	10734	0.23%	0.020%
35	10.193	1204	1208	1214	rVB7	7658	14887	0.32%	0.028%
36	10.287	1216	1224	1235	rBV	949034	1623377	34.52%	3.036%

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37	10.451	1246	1252	1261	rVB8	17732	37895	0.81%	0.071%
38	10.587	1268	1275	1281	rBV5	24037	63050	1.34%	0.118%
39	10.663	1281	1288	1295	rBV	877935	1512735	32.17%	2.829%
40	10.728	1295	1299	1305	rVB5	27406	46407	0.99%	0.087%
41	10.810	1305	1313	1324	rBV	935944	1656515	35.22%	3.098%
42	10.969	1337	1340	1348	rVB8	7904	12654	0.27%	0.024%
43	11.057	1348	1355	1363	rVV8	10487	26105	0.56%	0.049%
44	11.210	1374	1381	1386	rBV	1308380	1962223	41.72%	3.670%
45	11.269	1386	1391	1397	rBV	1470659	2275844	48.39%	4.257%
46	11.475	1418	1426	1433	rVV3	61440	138794	2.95%	0.260%
47	11.545	1433	1438	1455	rVB2	45350	102406	2.18%	0.192%
48	11.845	1485	1489	1492	rBV3	8140	11163	0.24%	0.021%
49	11.928	1498	1503	1509	rVB2	12287	21903	0.47%	0.041%
50	12.004	1509	1516	1522	rBV8	10530	25277	0.54%	0.047%
51	12.175	1540	1545	1550	rBV6	8247	12261	0.26%	0.023%
52	12.257	1550	1559	1564	rVV2	17981	36332	0.77%	0.068%
53	12.428	1583	1588	1593	rVB5	9883	17208	0.37%	0.032%
54	12.722	1635	1638	1647	rVB9	9968	15915	0.34%	0.030%
55	12.792	1647	1650	1657	rBV6	10401	19382	0.41%	0.036%
56	12.869	1657	1663	1670	rBV	84004	139102	2.96%	0.260%
57	12.981	1676	1682	1687	rBV	17239	26071	0.55%	0.049%
58	13.122	1700	1706	1717	rBV	169331	259566	5.52%	0.485%
59	13.263	1720	1730	1735	rVV10	10872	20940	0.45%	0.039%
60	13.357	1742	1746	1750	rBV	12216	17347	0.37%	0.032%
61	13.416	1750	1756	1764	rBV	1360106	2126280	45.21%	3.977%
62	13.745	1807	1812	1817	rBV8	7273	13120	0.28%	0.025%
63	13.839	1825	1828	1832	rVB5	7049	10213	0.22%	0.019%
64	13.934	1834	1844	1850	rVV	1771266	2426609	51.60%	4.539%
65	13.981	1850	1852	1859	rVV	31838	42831	0.91%	0.080%
66	14.045	1859	1863	1868	rVV7	10321	16975	0.36%	0.032%
67	14.204	1881	1890	1898	rBV2	1982025	2939017	62.49%	5.497%
68	14.386	1916	1921	1930	rVB6	10213	24444	0.52%	0.046%
69	14.510	1935	1942	1953	rBV	1199236	1789297	38.05%	3.347%
70	14.598	1953	1957	1963	rBV9	5341	10462	0.22%	0.020%
71	14.722	1971	1978	1993	rBV	103409	218589	4.65%	0.409%
72	14.939	2009	2015	2021	rVB10	10314	20833	0.44%	0.039%
73	15.116	2039	2045	2049	rVB6	11043	19701	0.42%	0.037%
74	15.198	2055	2059	2062	rBV3	8447	12011	0.26%	0.022%
75	15.263	2066	2070	2075	rVB	53001	67276	1.43%	0.126%
76	15.345	2080	2084	2097	rVB	114828	172052	3.66%	0.322%

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 Title : SVOA CALIBRATION

77	15.510	2105	2112	2126	rVB	2393912	3473396	73.86%	6.497%
78	15.633	2126	2133	2144	rBV	697890	1000473	21.27%	1.871%
79	15.751	2149	2153	2158	rVB5	12589	17163	0.36%	0.032%
80	15.886	2169	2176	2181	rBV	60317	87098	1.85%	0.163%
81	16.298	2240	2246	2249	rVB4	9652	14811	0.31%	0.028%
82	16.869	2337	2343	2348	rBV7	7507	12938	0.28%	0.024%
83	17.263	2401	2410	2418	rBV	1413409	2005769	42.65%	3.752%
84	17.363	2420	2427	2440	rBV2	2642868	3756396	79.87%	7.026%
85	17.463	2442	2444	2449	rVB4	8173	12188	0.26%	0.023%
86	17.563	2457	2461	2465	rVB	19881	26140	0.56%	0.049%
87	17.657	2474	2477	2486	rVB8	10371	19465	0.41%	0.036%
88	17.945	2519	2526	2531	rBV	425092	529950	11.27%	0.991%
89	18.175	2560	2565	2570	rVV	31438	45619	0.97%	0.085%
90	18.251	2574	2578	2586	rVB	36982	61296	1.30%	0.115%
91	18.380	2597	2600	2604	rBV4	9072	12010	0.26%	0.022%
92	18.610	2635	2639	2643	rBV4	18950	24880	0.53%	0.047%
93	19.227	2739	2744	2750	rBV3	32829	48719	1.04%	0.091%
94	19.298	2750	2756	2760	rBV	29487	43498	0.92%	0.081%
95	19.463	2780	2784	2789	rBV3	18839	24461	0.52%	0.046%
96	19.663	2812	2818	2826	rVV	3283275	4369490	92.91%	8.173%
97	20.668	2986	2989	2994	rBV7	10569	16619	0.35%	0.031%
98	21.463	3119	3124	3139	rBV	1618879	2129107	45.27%	3.982%
99	23.680	3493	3501	3512	rBV2	2377063	4702988	100.00%	8.797%
100	23.833	3520	3527	3538	rVB2	1134008	2307437	49.06%	4.316%

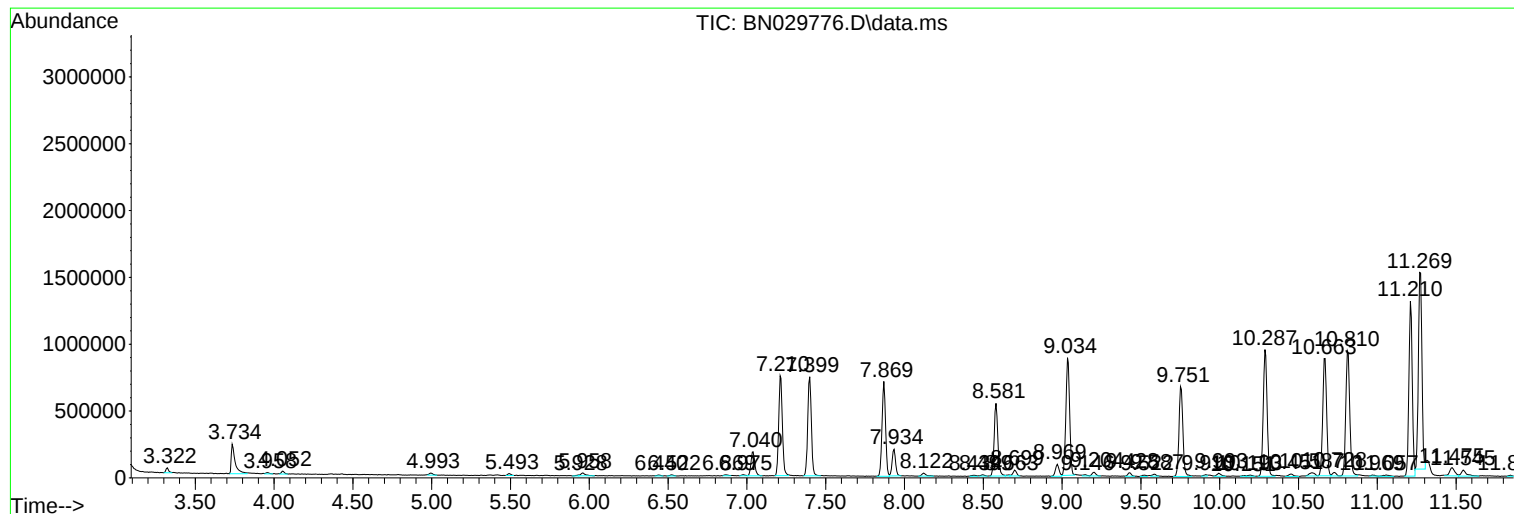
Sum of corrected areas: 53464144

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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



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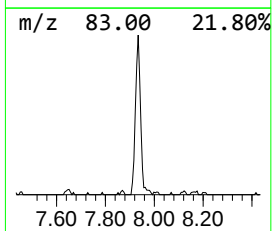
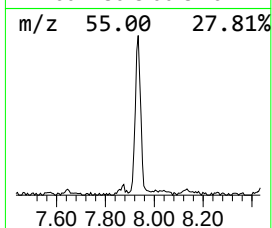
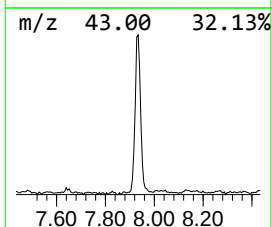
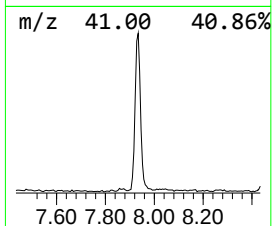
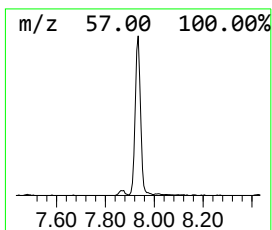
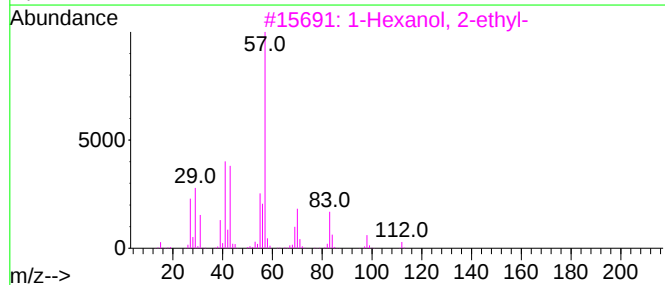
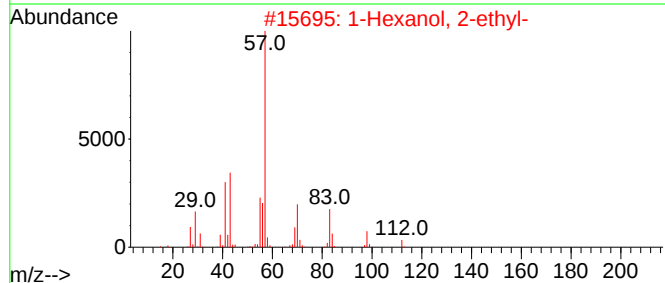
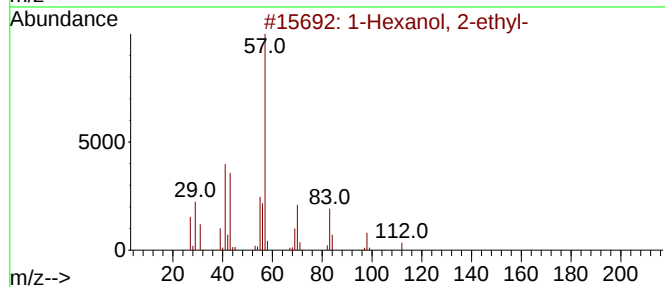
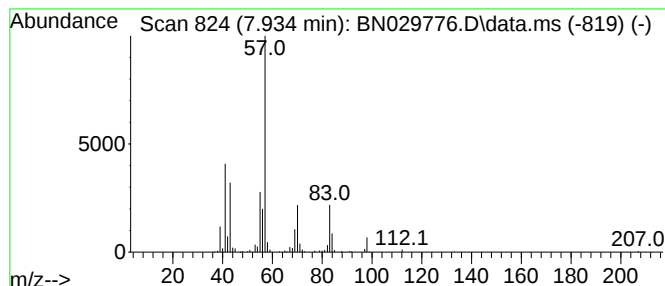
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.934	5.46 ng/ul	302429	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	59



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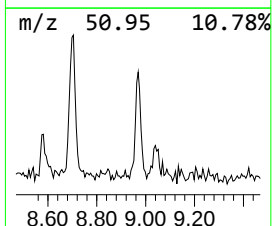
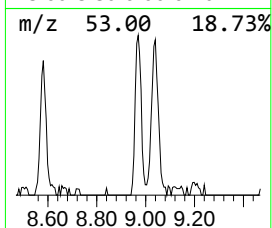
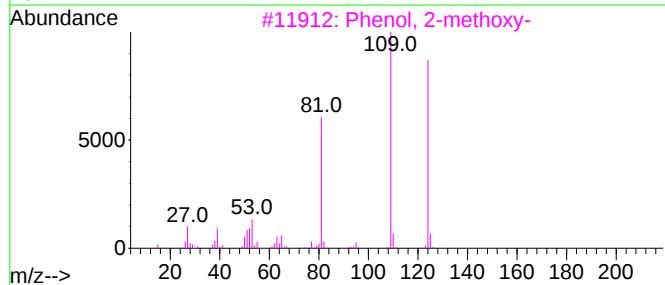
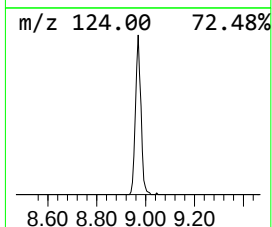
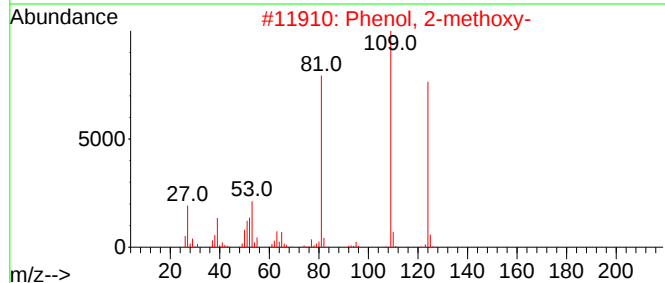
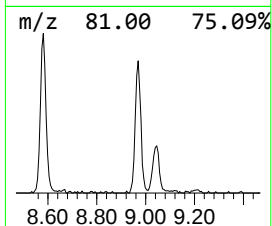
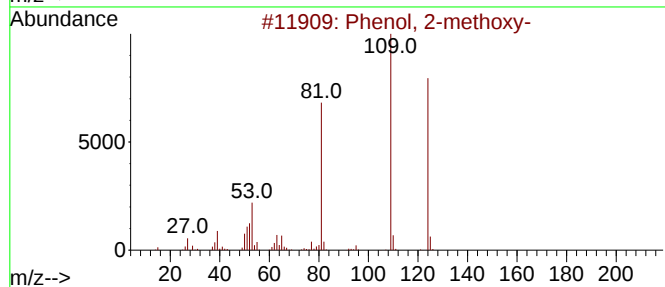
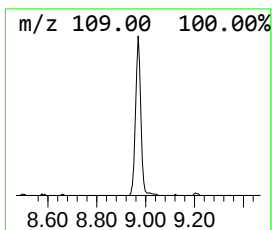
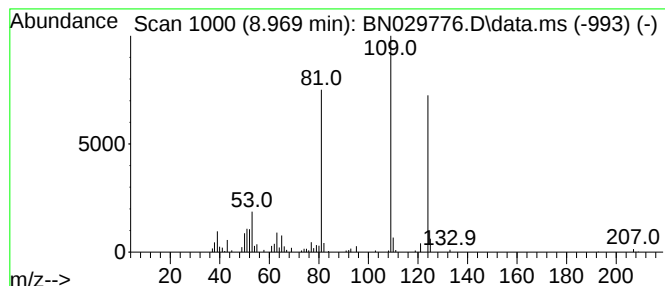
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 2-methoxy- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.969	2.67 ng/ul	147970	1,4-Dichlorobenzene-d4	7.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	97
2			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	97
3			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
4			Mequinol	124	C7H8O2	000150-76-5	90
5			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	87



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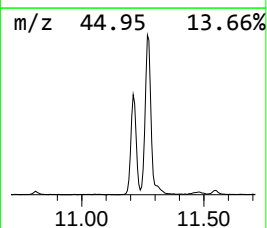
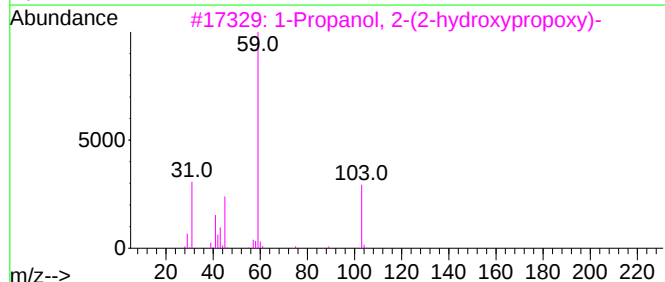
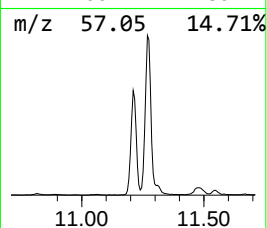
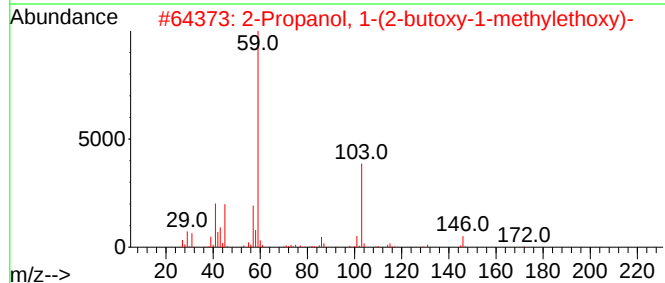
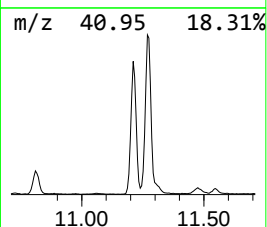
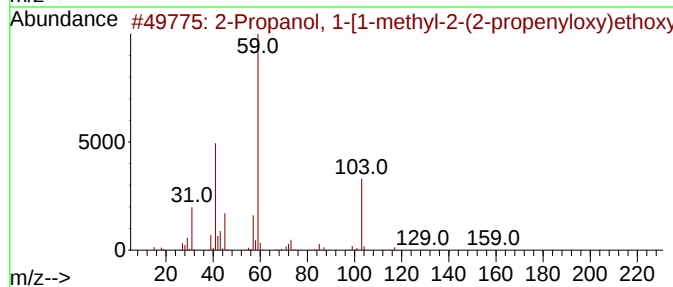
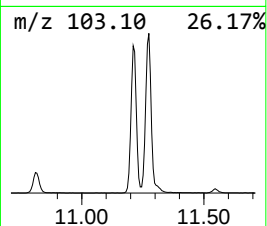
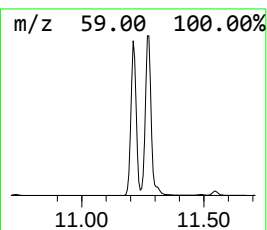
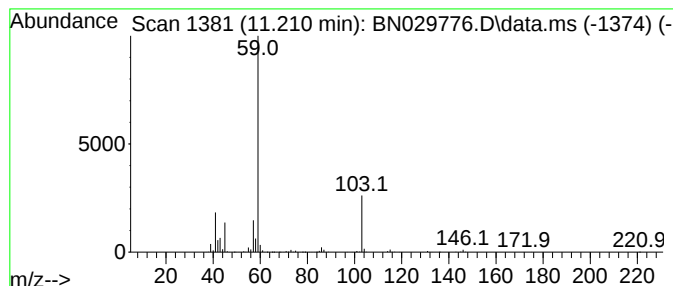
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1-[1-methyl-2-(... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.210	25.94 ng/ul	1962220	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	78		
2	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	74		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		
4	1-(1-tert-Butoxypropan-2-yloxy)p...	190	C10H22O3	132739-31-2	64		
5	2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029776.D
Acq On : 14 Feb 2024 19:04
Operator : MA/JU
Sample : P1289-16
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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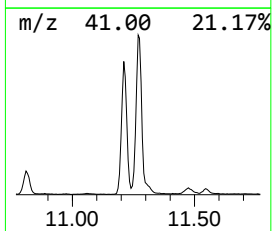
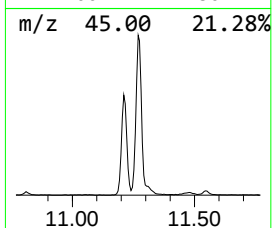
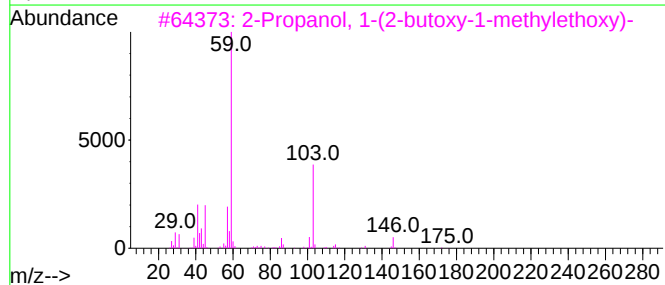
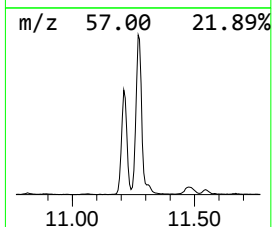
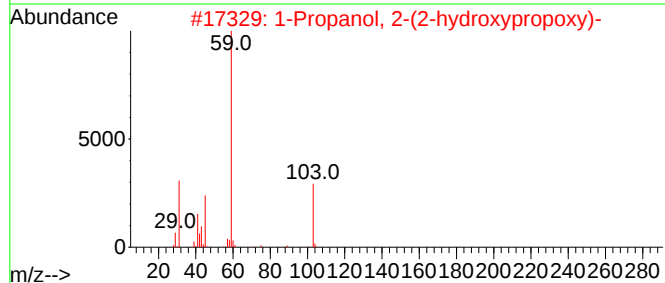
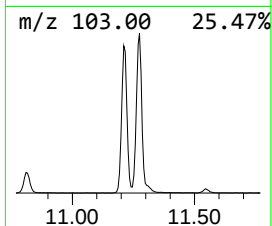
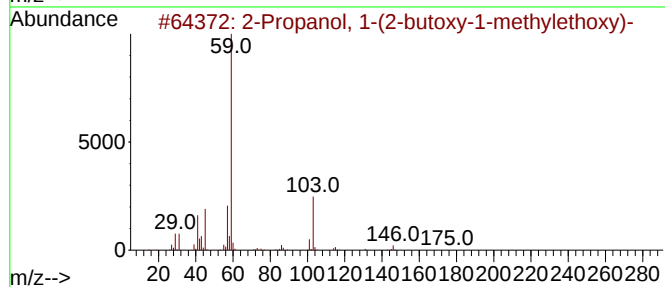
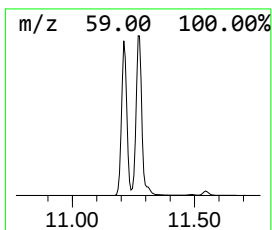
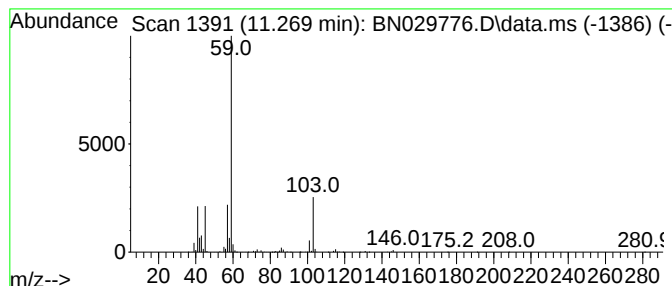
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.269	30.09 ng/ul	2275840	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	90		
2	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
3	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	74		
4	1,4,7-Trimethyl-3,6-dioxaoctane-...	192	C9H20O4	1000423-63-2	72		
5	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	72		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029776.D
Acq On : 14 Feb 2024 19:04
Operator : MA/JU
Sample : P1289-16
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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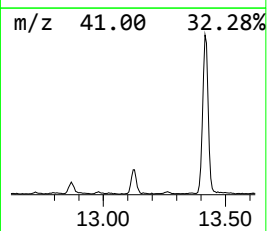
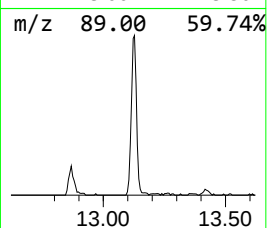
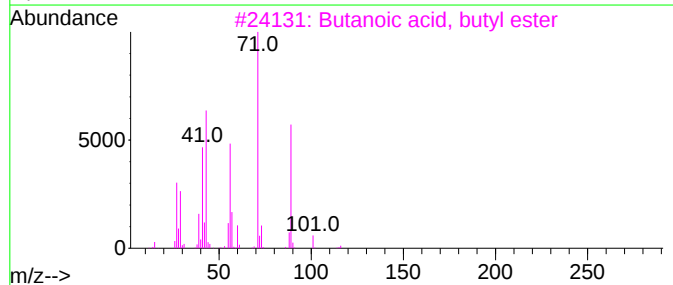
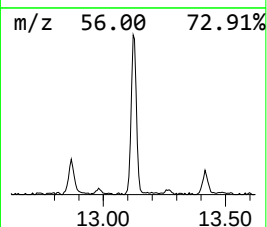
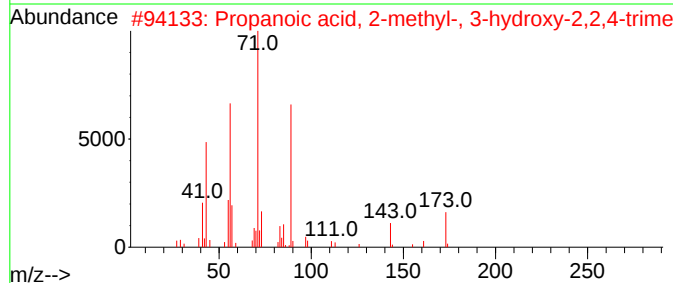
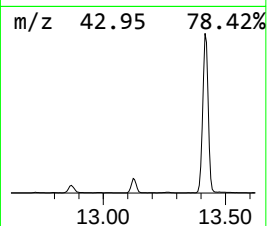
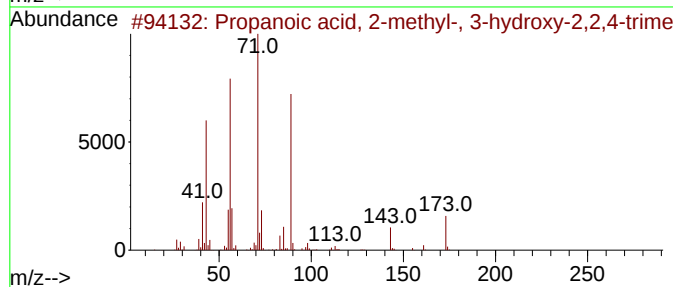
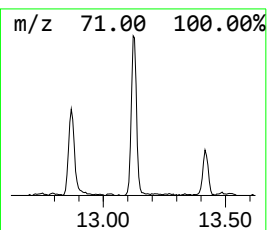
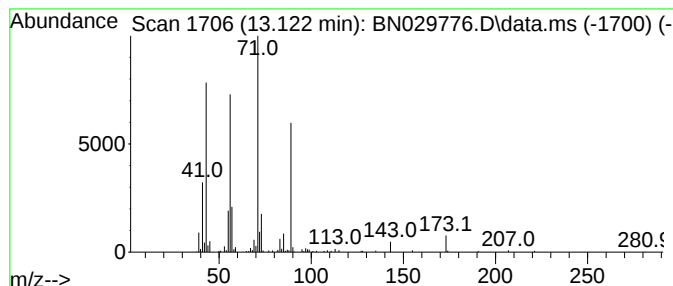
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Propanoic acid, 2-methyl-, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.122	2.90 ng/ul	259566	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	86
2			Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	000077-68-9	64
3			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64
4			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64
5			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029776.D
Acq On : 14 Feb 2024 19:04
Operator : MA/JU
Sample : P1289-16
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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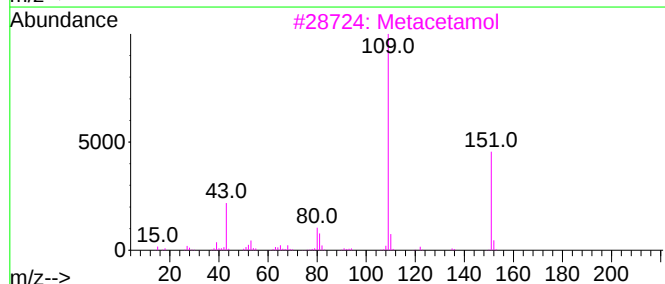
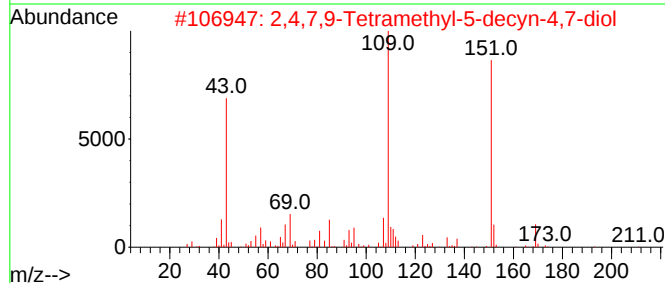
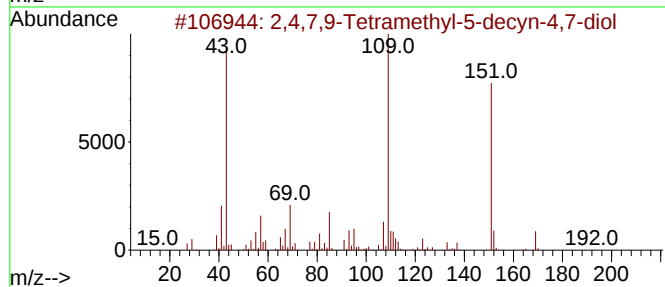
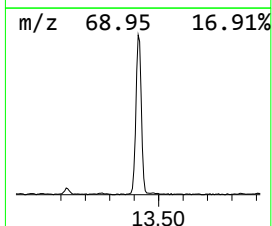
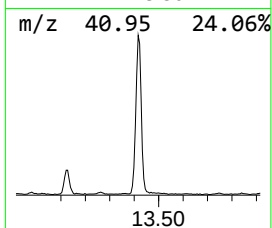
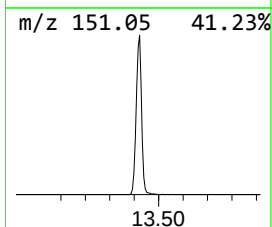
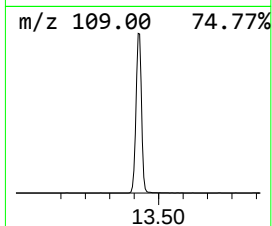
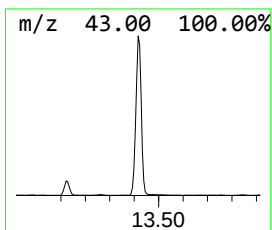
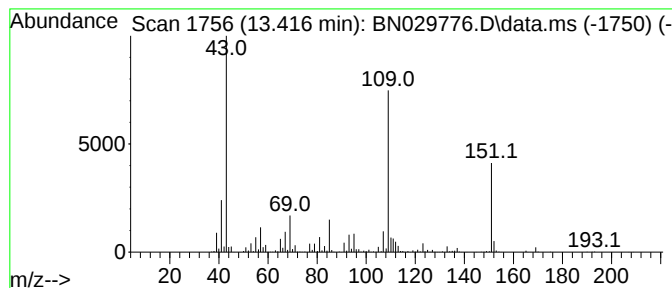
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.416	23.77 ng/ul	2126280	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	83		
3	Metacetamol	151	C8H9NO2	000621-42-1	64		
4	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		
5	Benzenamine, 4-propoxy-	151	C9H13NO	004469-80-1	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029776.D
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Sample : P1289-16
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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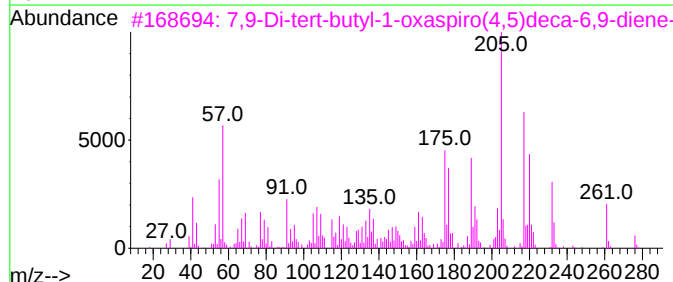
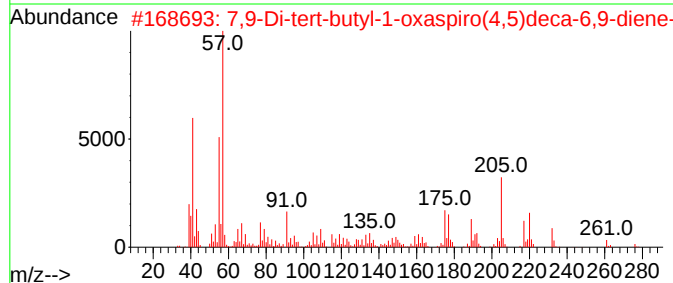
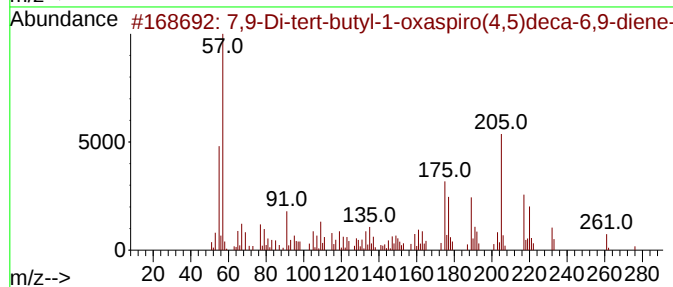
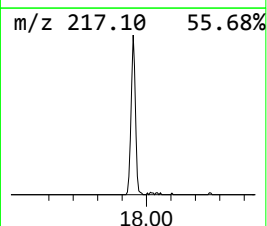
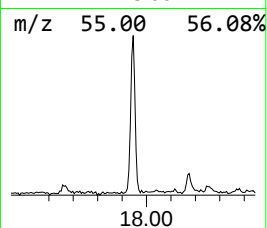
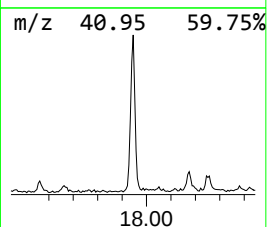
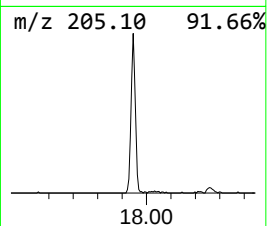
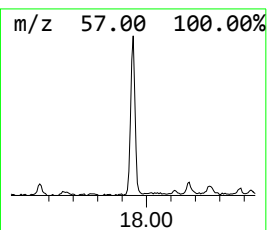
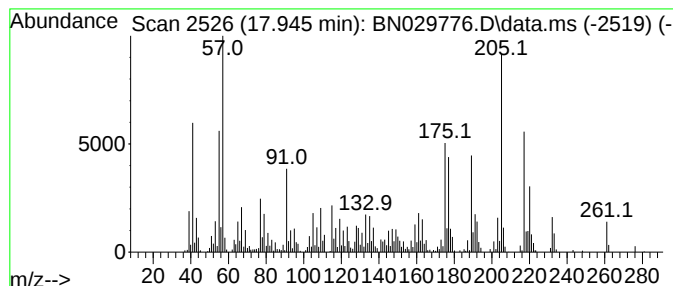
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.945	5.28 ng/ul	529950	Phenanthrene-d10	17.263

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	99
3			7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	91
4			(1R,3aR,5aR,9aS)-1,4,4,7-Tetrame...	220	C15H24O	104188-25-2	52
5			Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021424\
Data File : BN029776.D
Acq On : 14 Feb 2024 19:04
Operator : MA/JU
Sample : P1289-16
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.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
1-Hexanol, 2-et...	7.934	5.5	ng/ul	302429	1	7.869	1107260	20.0
Phenol, 2-methoxy-	8.969	2.7	ng/ul	147970	1	7.869	1107260	20.0
2-Propanol, 1-[...	11.210	25.9	ng/ul	1962220	2	10.663	1512740	20.0
2-Propanol, 1-(...	11.269	30.1	ng/ul	2275840	2	10.663	1512740	20.0
Propanoic acid,...	13.122	2.9	ng/ul	259566	3	14.510	1789300	20.0
2,4,7,9-Tetrame...	13.416	23.8	ng/ul	2126280	3	14.510	1789300	20.0
7,9-Di-tert-but...	17.945	5.3	ng/ul	529950	4	17.263	2005770	20.0