

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
 Data File : BP014662.D
 Acq On : 11 Apr 2023 18:35
 Operator : CG/JU
 Sample : 02251-11
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMC9

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-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014662.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.364	27	30	34	rBV4	10567	13978	0.41%	0.046%
2	3.817	105	107	124	rBV	52874	123988	3.66%	0.407%
3	4.017	137	141	147	rVB7	6923	12547	0.37%	0.041%
4	4.111	154	157	162	rVB	10013	11155	0.33%	0.037%
5	4.499	219	223	225	rVV4	3223	4304	0.13%	0.014%
6	4.528	227	228	235	rVB7	2828	5655	0.17%	0.019%
7	5.258	349	352	360	rVB5	9555	17884	0.53%	0.059%
8	5.493	388	392	397	rBV	7347	12030	0.36%	0.039%
9	6.011	476	480	489	rVB2	13582	21708	0.64%	0.071%
10	7.069	656	660	662	rBV5	2578	3634	0.11%	0.012%
11	7.152	668	674	689	rBV2	68818	164673	4.86%	0.540%
12	7.305	694	700	713	rBV	375378	640257	18.91%	2.100%
13	7.505	726	734	755	rBV	354643	649424	19.18%	2.130%
14	7.975	807	814	821	rBV	426450	728009	21.50%	2.388%
15	8.034	821	824	835	rVB	44261	79629	2.35%	0.261%
16	8.269	859	864	868	rBV6	4498	8889	0.26%	0.029%
17	8.434	889	892	895	rBV5	2278	3607	0.11%	0.012%
18	8.558	907	913	915	rBV7	2809	4354	0.13%	0.014%
19	8.616	918	923	930	rBV3	15368	34728	1.03%	0.114%
20	8.699	931	937	954	rVV	243964	478637	14.14%	1.570%
21	8.828	954	959	965	rVB7	6599	14381	0.42%	0.047%
22	9.081	997	1002	1008	rBV	152963	267709	7.91%	0.878%
23	9.152	1008	1014	1024	rVV	453658	788293	23.28%	2.586%
24	9.305	1036	1040	1047	rVB8	8332	14081	0.42%	0.046%
25	9.516	1071	1076	1082	rVB4	7573	13054	0.39%	0.043%
26	9.875	1129	1137	1153	rBV	357698	668585	19.75%	2.193%
27	10.040	1158	1165	1168	rBV7	5498	12157	0.36%	0.040%
28	10.275	1201	1205	1207	rBV5	3652	4254	0.13%	0.014%
29	10.310	1207	1211	1215	rVV6	5673	8234	0.24%	0.027%
30	10.422	1221	1230	1248	rBV	400232	883434	26.10%	2.898%
31	10.710	1272	1279	1286	rBV5	31517	63099	1.86%	0.207%
32	10.793	1286	1293	1300	rVV	611035	1025542	30.29%	3.364%
33	10.846	1300	1302	1308	rVB4	17509	26395	0.78%	0.087%
34	10.946	1313	1319	1343	rBV	317997	752265	22.22%	2.468%
35	11.416	1394	1399	1402	rBV6	2998	5776	0.17%	0.019%
36	11.487	1404	1411	1420	rVB10	10005	27694	0.82%	0.091%

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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 >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
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37	11.746	1451	1455	1460	rVB8	2725	4945	0.15%	0.016%
38	11.810	1465	1466	1471	rBV5	2646	3674	0.11%	0.012%
39	11.987	1489	1496	1499	rBV5	9938	20015	0.59%	0.066%
40	12.016	1499	1501	1506	rVB3	10101	13276	0.39%	0.044%
41	12.128	1514	1520	1527	rBV6	17016	43347	1.28%	0.142%
42	12.399	1560	1566	1573	rVB7	8574	16896	0.50%	0.055%
43	12.716	1615	1620	1623	rVB7	2890	4690	0.14%	0.015%
44	12.863	1641	1645	1648	rBV5	3500	4439	0.13%	0.015%
45	12.928	1651	1656	1659	rVV7	3856	6034	0.18%	0.020%
46	12.975	1659	1664	1672	rVB10	7085	13937	0.41%	0.046%
47	13.110	1681	1687	1688	rBV5	5033	7439	0.22%	0.024%
48	13.228	1701	1707	1714	rVB6	8576	17285	0.51%	0.057%
49	13.475	1744	1749	1752	rBV3	12132	20104	0.59%	0.066%
50	13.504	1752	1754	1765	rVB3	10851	19784	0.58%	0.065%
51	14.034	1834	1844	1857	rBV	1006591	1558718	46.04%	5.113%
52	14.140	1860	1862	1869	rVB8	5103	8416	0.25%	0.028%
53	14.322	1886	1893	1902	rBV	1169083	1670699	49.35%	5.480%
54	14.398	1902	1906	1911	rVB8	9140	17890	0.53%	0.059%
55	14.516	1920	1926	1933	rBV2	33177	64602	1.91%	0.212%
56	14.628	1939	1945	1954	rVB	898501	1301914	38.46%	4.271%
57	14.934	1991	1997	2000	rBV8	11219	23202	0.69%	0.076%
58	15.228	2044	2047	2049	rVV4	3321	3822	0.11%	0.013%
59	15.304	2055	2060	2065	rBV4	10866	22225	0.66%	0.073%
60	15.375	2068	2072	2078	rVV2	16699	25674	0.76%	0.084%
61	15.445	2080	2084	2093	rVB4	13824	26520	0.78%	0.087%
62	15.622	2108	2114	2128	rBV	1439235	2191188	64.72%	7.188%
63	15.763	2131	2138	2153	rVV	299052	571086	16.87%	1.873%
64	15.863	2153	2155	2162	rVB7	10815	15844	0.47%	0.052%
65	15.998	2173	2178	2186	rVB5	10072	18411	0.54%	0.060%
66	16.210	2208	2214	2217	rBV7	4881	7869	0.23%	0.026%
67	16.316	2230	2232	2239	rVB8	3353	4563	0.13%	0.015%
68	16.422	2245	2250	2253	rBV7	3942	5940	0.18%	0.019%
69	16.816	2311	2317	2318	rBV6	2865	4903	0.14%	0.016%
70	16.975	2340	2344	2349	rBV8	6755	10965	0.32%	0.036%
71	17.187	2375	2380	2385	rVB7	3997	7182	0.21%	0.024%
72	17.392	2408	2415	2425	rBV	1091279	1592934	47.05%	5.225%
73	17.492	2425	2432	2446	rVV	1659140	2512976	74.23%	8.243%
74	17.669	2458	2462	2466	rVB	16937	22163	0.65%	0.073%
75	17.863	2491	2495	2498	rBV6	5156	8694	0.26%	0.029%
76	18.045	2520	2526	2536	rBV	411782	506313	14.96%	1.661%

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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 >
 Peak separation: 5

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77	18.204	2550	2553	2558	rVB7	8539	13647	0.40%	0.045%
78	18.275	2561	2565	2569	rBV6	9365	17756	0.52%	0.058%
79	18.334	2572	2575	2578	rBV	13761	15858	0.47%	0.052%
80	18.475	2595	2599	2602	rBV4	16882	27081	0.80%	0.089%
81	18.510	2602	2605	2611	rVB8	12940	19736	0.58%	0.065%
82	18.745	2643	2645	2648	rBV3	4719	4743	0.14%	0.016%
83	19.304	2736	2740	2746	rVB4	25297	35209	1.04%	0.115%
84	19.369	2749	2751	2755	rVB2	16473	20460	0.60%	0.067%
85	19.722	2806	2811	2828	rBV	2419094	3184490	94.06%	10.446%
86	21.480	3106	3110	3121	rBV	1156963	1832400	54.13%	6.011%
87	23.821	3501	3508	3519	rVB2	1695930	3385441	100.00%	11.105%
88	23.980	3529	3535	3548	rVB2	857454	1933937	57.13%	6.344%

Sum of corrected areas: 30485379

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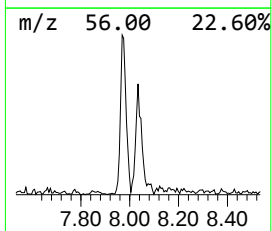
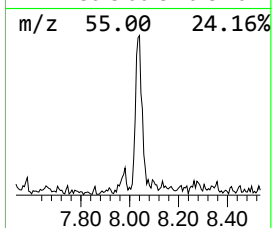
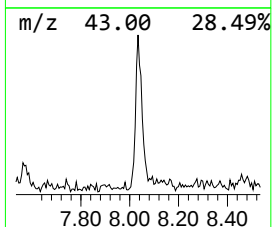
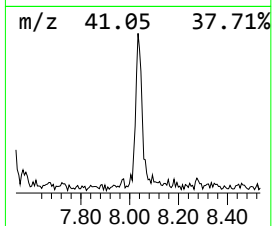
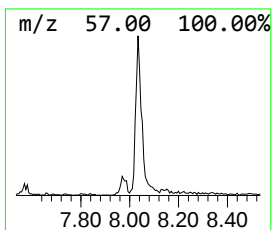
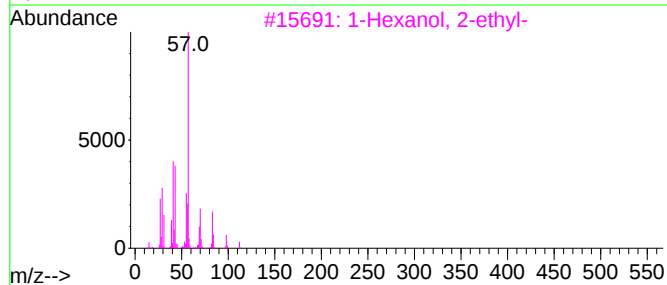
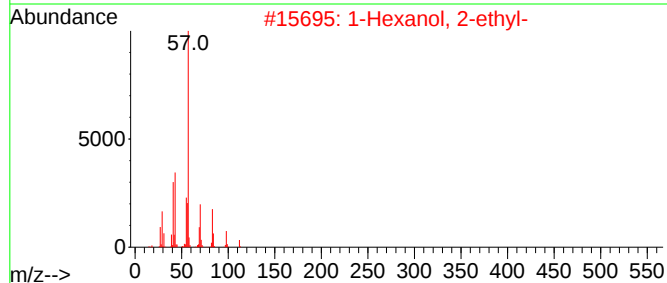
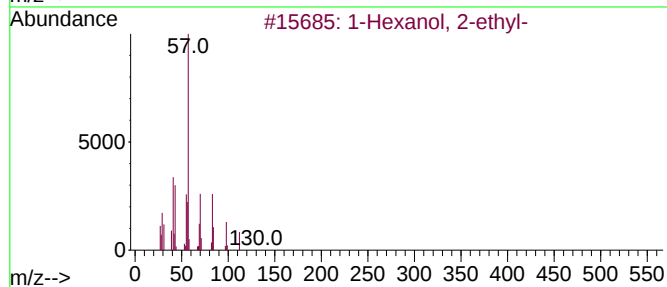
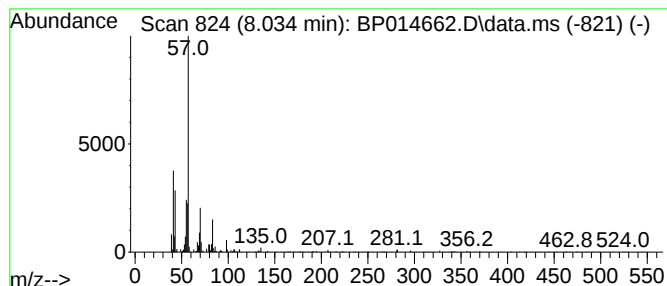
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.034	2.19 ng/ul	79629	1,4-Dichlorobenzene-d4	7.975

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	56
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	56
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	56
5	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	45



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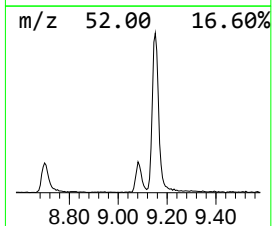
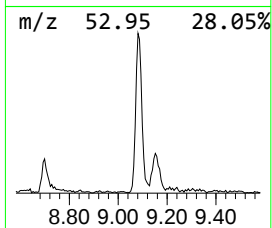
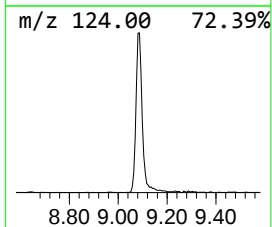
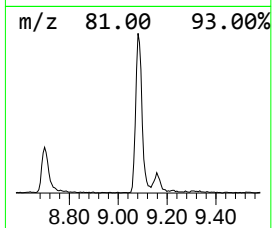
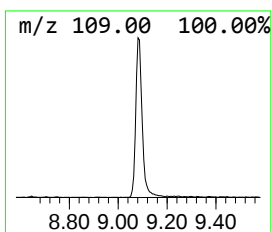
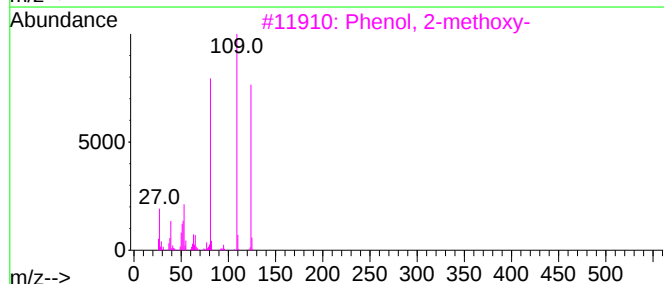
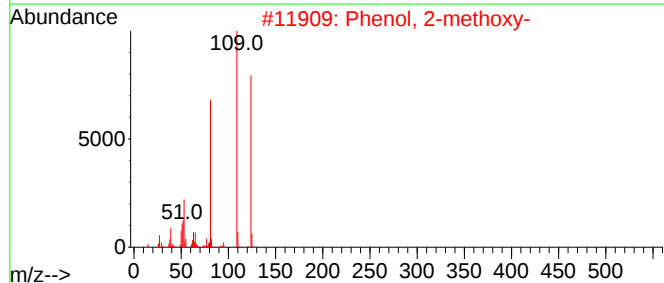
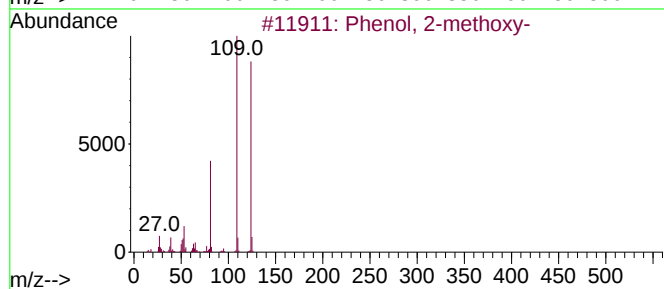
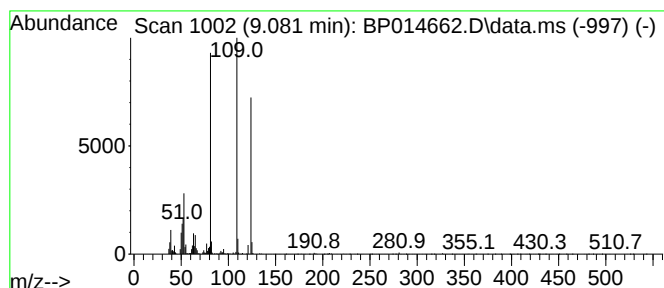
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 2-methoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.081	7.35 ng/ul	267709	1,4-Dichlorobenzene-d4	7.975

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	95
2			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	95
3			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	95
4			Mequinol	124	C7H8O2	000150-76-5	91
5			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	91



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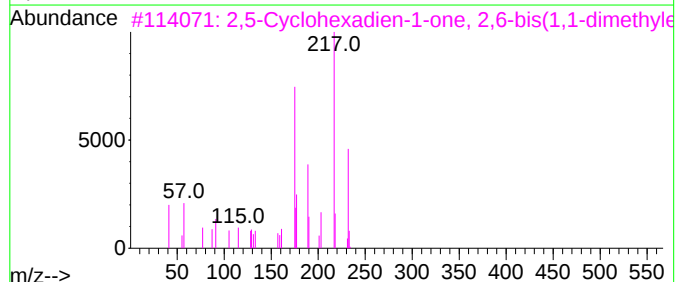
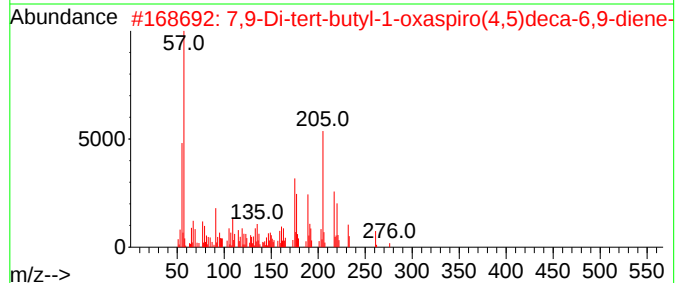
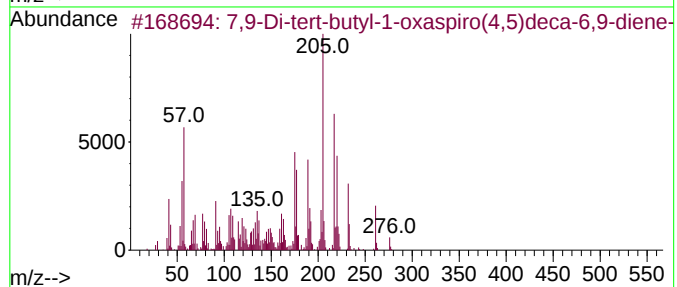
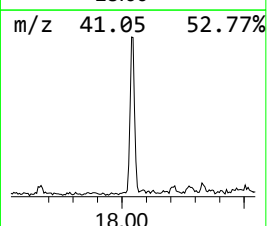
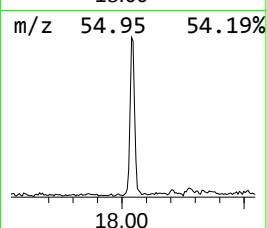
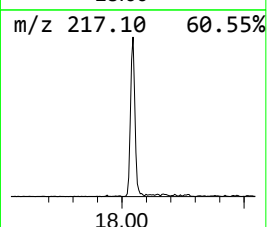
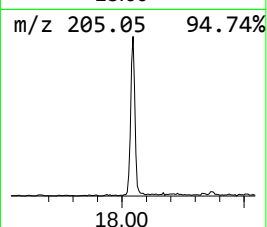
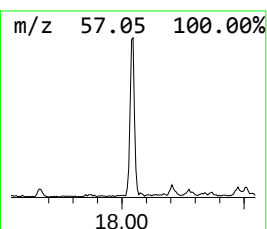
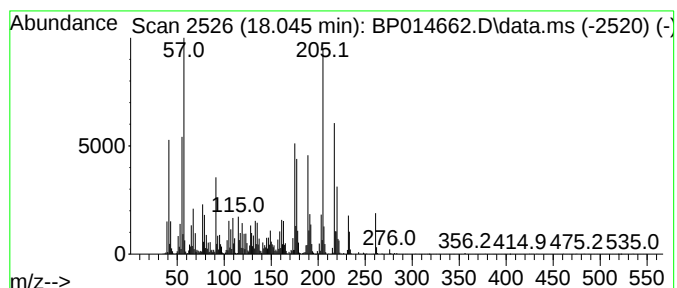
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.045	6.36 ng/ul	506313	Phenanthrene-d10	17.392	
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	95
3	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	86
4	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	64
5	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	50



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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
1-Hexanol, 2-et...	8.034	2.2	ng/ul	79629	1	7.975	728009	20.0
Phenol, 2-methoxy-	9.081	7.3	ng/ul	267709	1	7.975	728009	20.0
7,9-Di-tert-but...	18.045	6.4	ng/ul	506313	4	17.392	1592930	20.0