

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018224.D
 Acq On : 17 Nov 2023 16:46
 Operator : MA/JU
 Sample : 05369-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDN4

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-BP111023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018224.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	9	13	18	rBV2	10425	13460	0.62%	0.077%
2	3.611	85	89	102	rBV	36029	74724	3.42%	0.427%
3	3.899	134	138	139	rBV4	3261	3894	0.18%	0.022%
4	4.358	214	216	218	rBV3	2453	2291	0.10%	0.013%
5	4.411	220	225	233	rVV	22497	37354	1.71%	0.213%
6	6.246	532	537	540	rBV8	1907	3763	0.17%	0.022%
7	6.969	655	660	672	rVV	32475	78035	3.58%	0.446%
8	7.134	682	688	702	rBV	203754	331334	15.18%	1.893%
9	7.305	711	717	727	rBV	186516	335983	15.39%	1.920%
10	7.487	745	748	752	rVB6	1621	2396	0.11%	0.014%
11	7.652	774	776	782	rVB6	2715	3693	0.17%	0.021%
12	7.781	792	798	815	rBV	142655	268853	12.32%	1.536%
13	8.516	917	923	944	rBV	126628	281624	12.90%	1.609%
14	8.987	997	1003	1018	rBV	274873	525701	24.09%	3.004%
15	9.158	1028	1032	1037	rVB5	4937	8356	0.38%	0.048%
16	9.705	1117	1125	1142	rBV	225284	452384	20.73%	2.585%
17	10.234	1208	1215	1244	rBV	249432	584605	26.78%	3.341%
18	10.599	1269	1277	1289	rBV	234281	405707	18.59%	2.318%
19	10.787	1299	1309	1337	rBV	221172	564062	25.84%	3.223%
20	11.546	1436	1438	1444	rVB6	2152	3231	0.15%	0.018%
21	11.646	1448	1455	1459	rVV7	2339	4107	0.19%	0.023%
22	12.251	1553	1558	1559	rBV5	2871	3903	0.18%	0.022%
23	13.298	1730	1736	1744	rBV2	22419	39379	1.80%	0.225%
24	13.751	1802	1813	1823	rBV3	44276	175510	8.04%	1.003%
25	13.887	1830	1836	1848	rBV	733887	1067948	48.93%	6.103%
26	14.157	1876	1882	1893	rBV	748037	1139921	52.23%	6.514%
27	14.463	1927	1934	1944	rBV	359103	536449	24.58%	3.066%
28	14.828	1989	1996	1997	rBV6	6203	10738	0.49%	0.061%
29	15.357	2082	2086	2091	rVB7	2175	4836	0.22%	0.028%
30	15.463	2098	2104	2113	rBV	1001028	1487122	68.14%	8.498%
31	15.628	2126	2132	2145	rBV	287782	498305	22.83%	2.848%
32	16.181	2224	2226	2231	rVB5	2891	4433	0.20%	0.025%
33	16.269	2237	2241	2245	rVB6	1554	2921	0.13%	0.017%
34	16.334	2250	2252	2257	rBV6	1878	3438	0.16%	0.020%
35	16.492	2275	2279	2280	rBV4	2702	2534	0.12%	0.014%
36	16.551	2285	2289	2292	rVB5	2564	3467	0.16%	0.020%

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37	16.598	2292	2297	2298	rBV5	4251	5240	0.24%	0.030%
38	17.245	2400	2407	2417	rVV	446233	626541	28.71%	3.580%
39	17.345	2417	2424	2441	rBV	1064883	1622277	74.33%	9.271%
40	17.492	2447	2449	2455	rVB7	4653	6027	0.28%	0.034%
41	17.875	2511	2514	2520	rBV7	5787	8921	0.41%	0.051%
42	18.098	2547	2552	2560	rBV4	20997	38839	1.78%	0.222%
43	18.310	2583	2588	2597	rBV	180311	266193	12.20%	1.521%
44	19.045	2710	2713	2716	rBV5	2836	3826	0.18%	0.022%
45	19.175	2731	2735	2740	rBV3	14183	21102	0.97%	0.121%
46	19.257	2744	2749	2754	rVB	12299	21068	0.97%	0.120%
47	19.345	2761	2764	2773	rBV7	13785	24421	1.12%	0.140%
48	19.475	2783	2786	2790	rBV6	7688	8983	0.41%	0.051%
49	19.604	2802	2808	2820	rBV	1590602	2135642	97.85%	12.204%
50	21.380	3105	3110	3121	rBV	525918	767628	35.17%	4.387%
51	23.686	3494	3502	3514	rBV2	1092601	2182610	100.00%	12.473%
52	23.851	3523	3530	3540	rBV	372272	793057	36.34%	4.532%

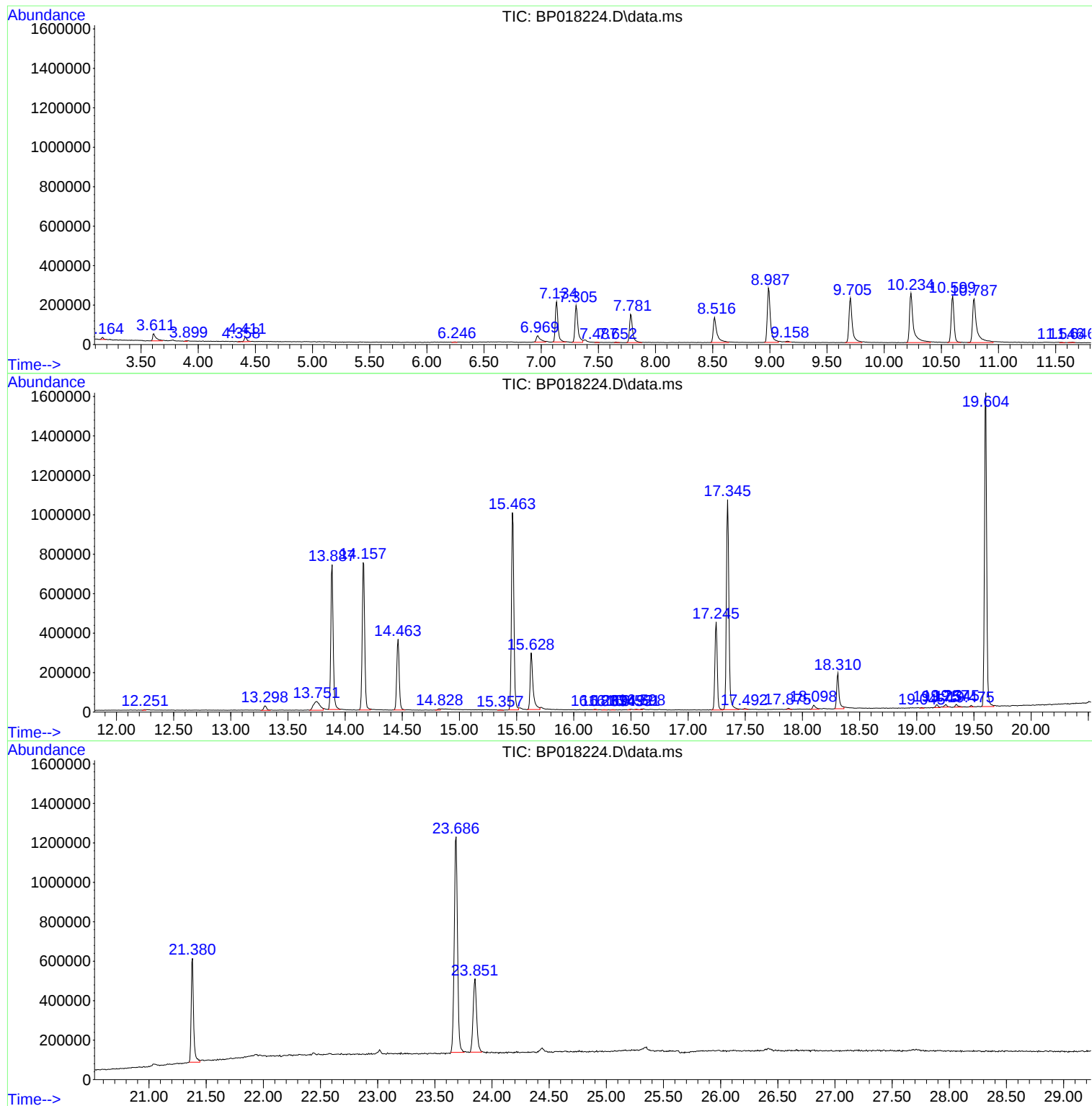
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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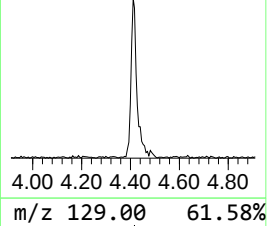
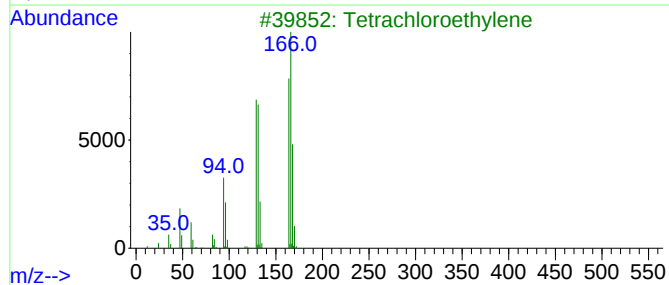
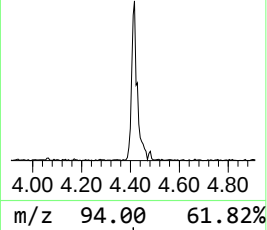
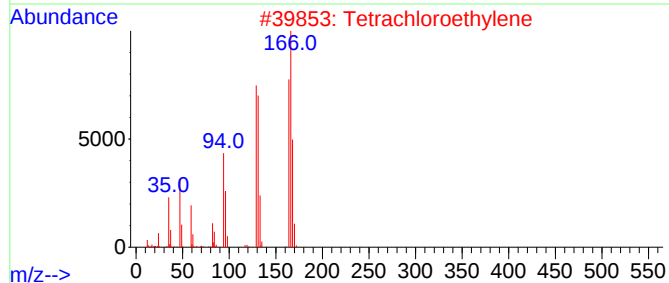
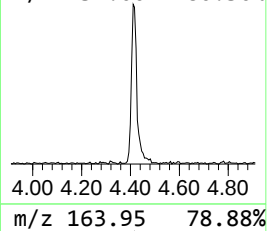
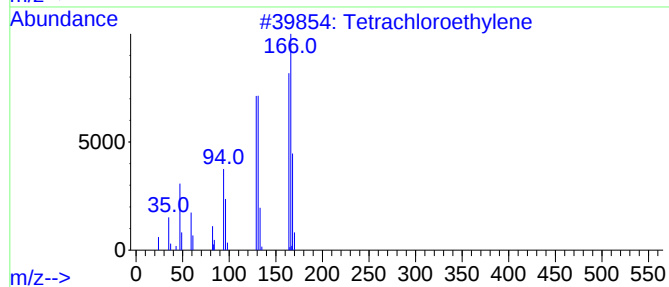
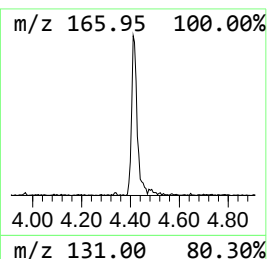
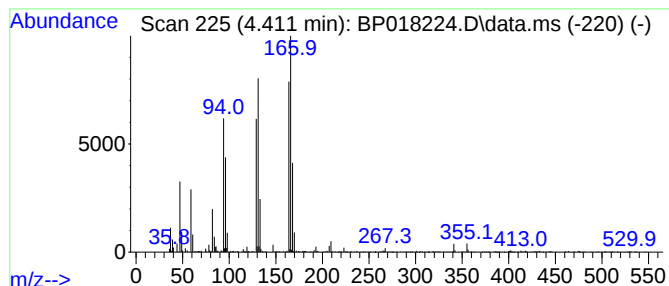
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.411	2.78 ng/ul	37354	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	93
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	91
5			1,3-Dichloro-2-fluorobenzene	164	C6H3Cl2F	002268-05-5	35



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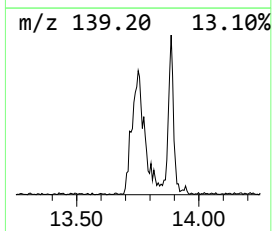
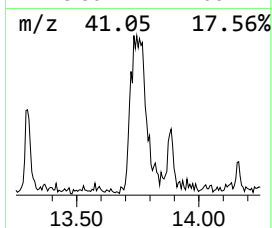
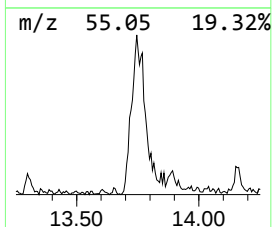
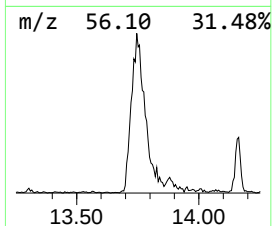
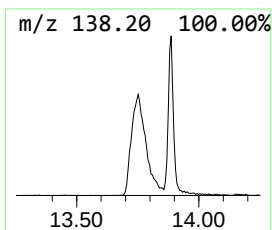
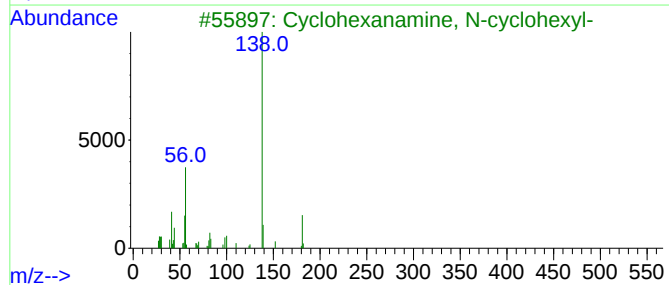
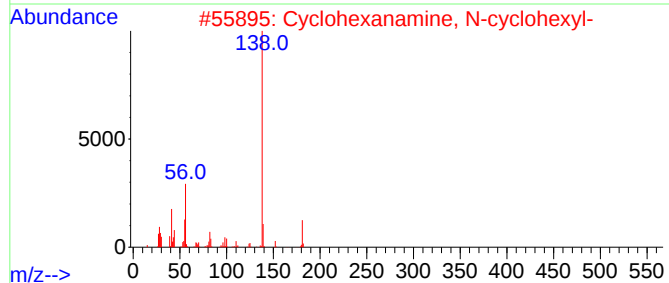
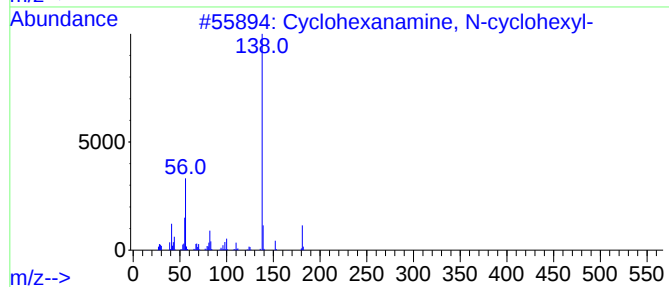
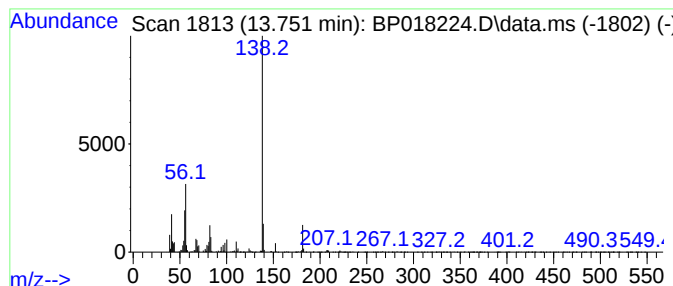
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclohexanamine, N-cyclohexyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	6.54 ng/ul	175510	Acenaphthene-d10	14.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	97
2			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
3			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	91
4			(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	91
5			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	91



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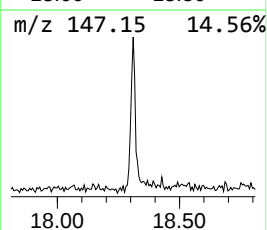
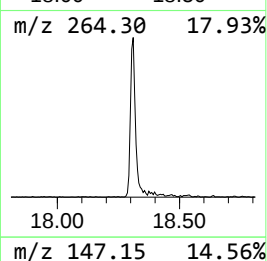
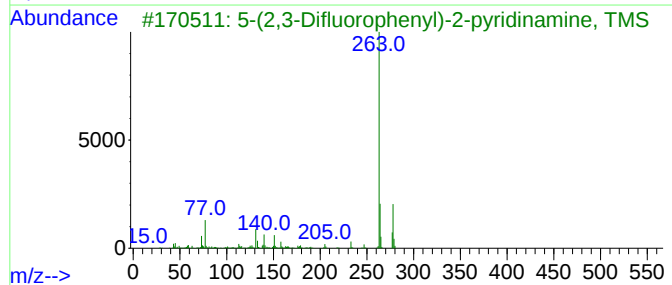
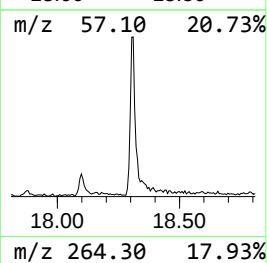
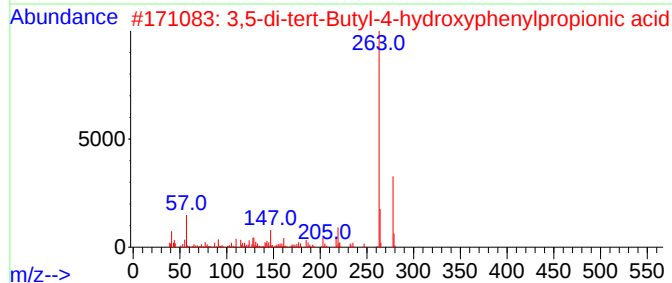
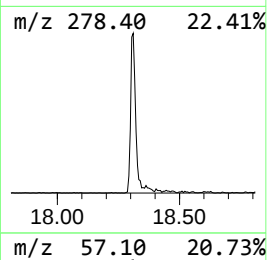
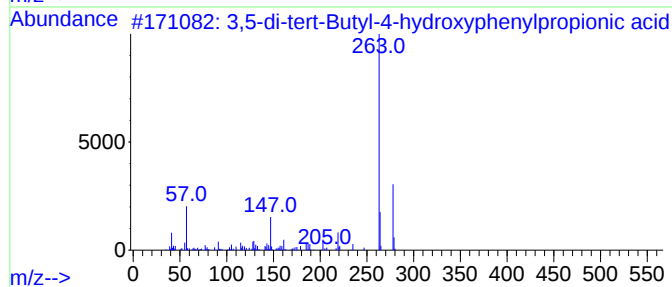
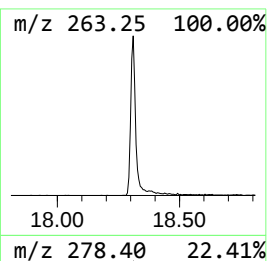
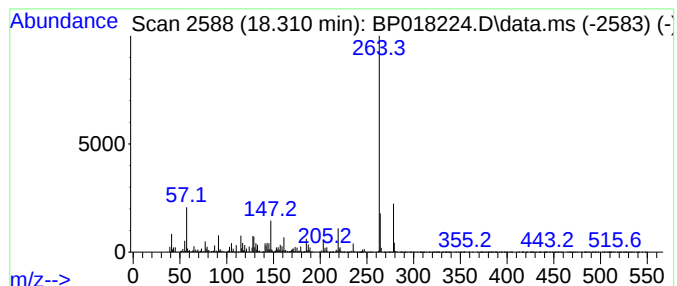
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.310	8.50 ng/ul	266193	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	96
2			3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3			5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	74
4			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	64
5			Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	59



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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
Tetrachloroethy...	4.411	2.8	ng/ul	37354	1	7.781	268853	20.0
Cyclohexanamine...	13.751	6.5	ng/ul	175510	3	14.463	536449	20.0
3,5-di-tert-But...	18.310	8.5	ng/ul	266193	4	17.245	626541	20.0