

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
 Data File : BP019964.D
 Acq On : 22 Mar 2024 20:41
 Operator : MA/JU
 Sample : P1813-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0LX9

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-BP032224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019964.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	23	27	37	rVB	34575	52542	1.42%	0.146%
2	3.499	68	70	74	rVB4	5969	6195	0.17%	0.017%
3	3.646	91	95	111	rBV	158502	279281	7.56%	0.778%
4	3.940	142	145	150	rVB2	8871	12975	0.35%	0.036%
5	4.058	161	165	171	rVB8	3702	6106	0.17%	0.017%
6	4.287	202	204	213	rVV10	4621	8798	0.24%	0.025%
7	4.370	214	218	226	rVB3	15131	25376	0.69%	0.071%
8	5.017	323	328	332	rVB4	6425	9509	0.26%	0.027%
9	5.128	343	347	353	rVB7	4078	7819	0.21%	0.022%
10	5.981	485	492	495	rBV3	7917	15547	0.42%	0.043%
11	6.622	595	601	604	rBV7	3289	6930	0.19%	0.019%
12	6.811	624	633	645	rBV	127213	246854	6.68%	0.688%
13	6.893	645	647	653	rVB6	2612	4685	0.13%	0.013%
14	6.969	653	660	674	rBV	446300	791579	21.42%	2.206%
15	7.163	683	693	709	rBV	453466	776447	21.01%	2.164%
16	7.369	725	728	734	rVB8	4144	7204	0.19%	0.020%
17	7.522	751	754	760	rVB7	2739	5187	0.14%	0.014%
18	7.616	763	770	779	rVV	406596	654616	17.72%	1.825%
19	7.716	783	787	794	rVB5	11591	22266	0.60%	0.062%
20	7.981	822	832	839	rBV3	20152	52339	1.42%	0.146%
21	8.146	852	860	865	rBV	17899	33924	0.92%	0.095%
22	8.269	876	881	884	rBV7	6863	14106	0.38%	0.039%
23	8.322	884	890	904	rVB	367024	634112	17.16%	1.767%
24	8.516	915	923	929	rVB7	10198	26664	0.72%	0.074%
25	8.622	932	941	947	rBV9	10322	23965	0.65%	0.067%
26	8.769	957	966	978	rBV	575198	1026662	27.78%	2.862%
27	8.969	991	1000	1009	rVB	40547	77762	2.10%	0.217%
28	9.110	1017	1024	1030	rBV4	34803	68972	1.87%	0.192%
29	9.187	1032	1037	1047	rVB	14633	27654	0.75%	0.077%
30	9.328	1054	1061	1065	rBV8	8004	17090	0.46%	0.048%
31	9.387	1067	1071	1072	rBV4	5845	5294	0.14%	0.015%
32	9.410	1072	1075	1079	rVB6	6944	9057	0.25%	0.025%
33	9.475	1079	1086	1095	rBV	439142	806782	21.83%	2.249%
34	9.763	1130	1135	1146	rVB7	20160	47856	1.30%	0.133%
35	9.910	1157	1160	1165	rVB7	2967	4310	0.12%	0.012%
36	10.010	1169	1177	1187	rBV	554118	1167998	31.61%	3.256%

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Integrator: RTE

Smoothing : OFF

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
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37	10.134	1196	1198	1203	rVB6	3938	6228	0.17%	0.017%
38	10.210	1205	1211	1218	rVB4	24175	50941	1.38%	0.142%
39	10.369	1231	1238	1251	rBV	526780	1015158	27.47%	2.830%
40	10.528	1256	1265	1276	rBV	333840	697493	18.88%	1.944%
41	10.693	1290	1293	1296	rBV4	4760	6311	0.17%	0.018%
42	10.728	1296	1299	1301	rBV3	7616	10245	0.28%	0.029%
43	10.857	1317	1321	1323	rBV4	4193	5804	0.16%	0.016%
44	10.934	1329	1334	1337	rBV3	14119	23695	0.64%	0.066%
45	10.975	1337	1341	1346	rVB6	9925	17944	0.49%	0.050%
46	11.093	1355	1361	1367	rBV	66318	107074	2.90%	0.298%
47	11.146	1367	1370	1374	rBV4	8790	14156	0.38%	0.039%
48	11.199	1375	1379	1380	rBV4	6101	8490	0.23%	0.024%
49	11.363	1402	1407	1410	rBV6	8272	14645	0.40%	0.041%
50	11.404	1410	1414	1415	rVV3	12337	17343	0.47%	0.048%
51	11.434	1416	1419	1427	rVB5	21152	45765	1.24%	0.128%
52	11.581	1439	1444	1451	rBV7	20661	52215	1.41%	0.146%
53	11.640	1451	1454	1460	rVB3	16347	24356	0.66%	0.068%
54	11.722	1460	1468	1473	rBV5	49484	110093	2.98%	0.307%
55	11.775	1473	1477	1482	rBV6	13765	19327	0.52%	0.054%
56	11.828	1482	1486	1494	rVB10	15259	40336	1.09%	0.112%
57	11.916	1494	1501	1506	rBV2	46696	104566	2.83%	0.291%
58	11.969	1506	1510	1512	rVV4	10950	13857	0.38%	0.039%
59	12.010	1512	1517	1522	rVV2	22333	39761	1.08%	0.111%
60	12.069	1523	1527	1531	rBV7	19516	29166	0.79%	0.081%
61	12.163	1540	1543	1549	rVB6	25341	38929	1.05%	0.109%
62	12.263	1555	1560	1566	rBV8	13297	32311	0.87%	0.090%
63	12.328	1567	1571	1580	rVB9	22327	45295	1.23%	0.126%
64	12.428	1580	1588	1594	rBV7	28723	66994	1.81%	0.187%
65	12.487	1594	1598	1601	rBV5	18817	36944	1.00%	0.103%
66	12.728	1636	1639	1640	rBV3	9872	10676	0.29%	0.030%
67	12.846	1657	1659	1665	rVB6	11198	15897	0.43%	0.044%
68	12.946	1671	1676	1677	rBV4	14760	20323	0.55%	0.057%
69	13.034	1686	1691	1695	rBV3	46028	87775	2.38%	0.245%
70	13.269	1726	1731	1734	rBV4	25498	45295	1.23%	0.126%
71	13.428	1754	1758	1760	rBV5	13114	17412	0.47%	0.049%
72	13.469	1761	1765	1769	rVV4	32269	48587	1.31%	0.135%
73	13.516	1771	1773	1779	rVB	26177	40488	1.10%	0.113%
74	13.675	1794	1800	1807	rVB	1218494	1828475	49.48%	5.096%
75	13.875	1831	1834	1838	rBV5	21723	32267	0.87%	0.090%
76	13.940	1839	1845	1854	rVB	1477201	2172137	58.78%	6.054%

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 Title : SVOA CALIBRATION

77	14.081	1862	1869	1877	rBV	75743	153228	4.15%	0.427%
78	14.181	1883	1886	1891	rVB6	33747	40498	1.10%	0.113%
79	14.245	1892	1897	1907	rVB	880940	1268385	34.33%	3.535%
80	14.345	1912	1914	1915	rBV2	8931	6830	0.18%	0.019%
81	14.457	1928	1933	1938	rBV4	53718	97014	2.63%	0.270%
82	14.516	1938	1943	1952	rVB2	80887	179995	4.87%	0.502%
83	14.657	1964	1967	1978	rVB2	31391	74127	2.01%	0.207%
84	14.757	1982	1984	1991	rVB4	21049	34004	0.92%	0.095%
85	15.187	2053	2057	2061	rBV2	44238	73865	2.00%	0.206%
86	15.275	2067	2072	2085	rVB	1879846	2706569	73.25%	7.544%
87	15.410	2088	2095	2108	rBV	504422	932313	25.23%	2.599%
88	15.540	2113	2117	2126	rVB2	59246	124572	3.37%	0.347%
89	15.816	2159	2164	2175	rVB2	146097	287139	7.77%	0.800%
90	16.363	2252	2257	2266	rVB	78861	133670	3.62%	0.373%
91	17.081	2372	2379	2388	rVB	793363	1499613	40.58%	4.180%
92	17.192	2392	2398	2410	rBV2	1887583	2909191	78.73%	8.109%
93	18.039	2538	2542	2551	rBV	209554	349493	9.46%	0.974%
94	19.386	2768	2771	2777	rVB3	65214	87115	2.36%	0.243%
95	19.586	2799	2805	2812	rVB	2271322	3437252	93.02%	9.581%
96	19.657	2814	2817	2822	rVB	82341	122284	3.31%	0.341%
97	21.233	3081	3085	3092	rVB2	156719	235578	6.38%	0.657%
98	21.551	3134	3139	3145	rVB2	1028346	1608887	43.54%	4.484%
99	24.610	3649	3659	3669	rVB2	1352762	3695197	100.00%	10.300%
100	24.839	3690	3698	3711	rVB	618063	1821205	49.29%	5.076%

Sum of corrected areas: 35877261

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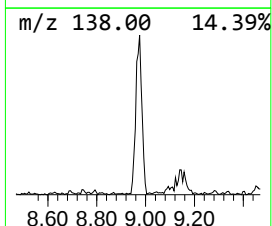
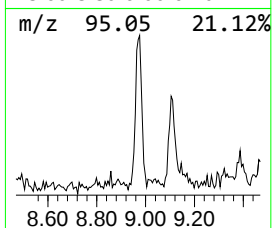
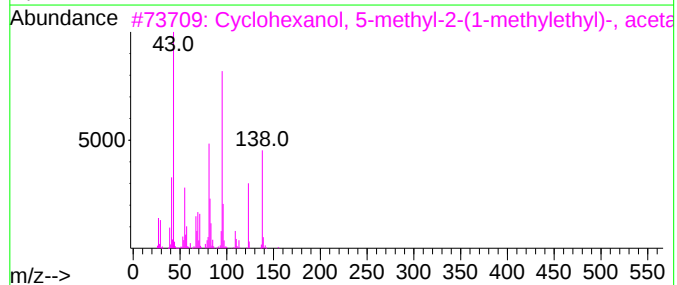
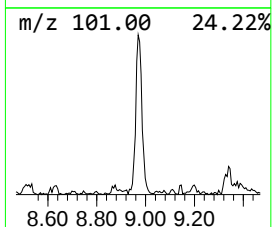
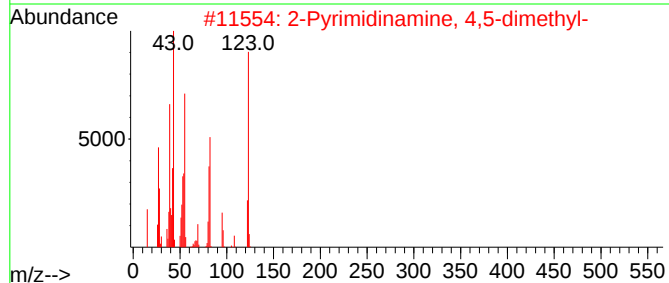
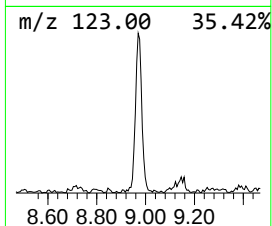
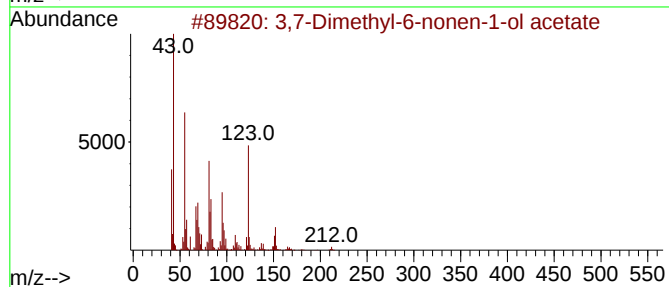
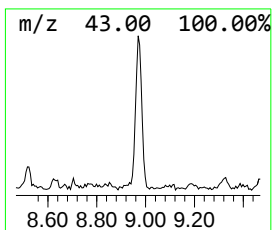
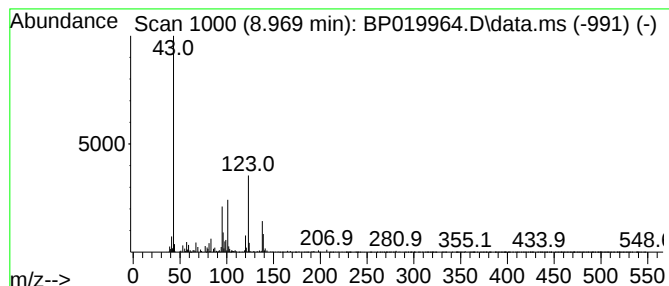
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.969	2.38 ng/ul	77762	1,4-Dichlorobenzene-d4	7.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,7-Dimethyl-6-nonen-1-ol acetate	212	C13H24O2	1010131-34-9	37
2			2-Pyrimidinamine, 4,5-dimethyl-	123	C6H9N3	001193-74-4	30
3			Cyclohexanol, 5-methyl-2-(1-meth...	198	C12H22O2	016409-45-3	25
4			Cyclohexanol, 5-methyl-2-(1-meth...	198	C12H22O2	020777-45-1	17
5			Menthyl acetate	198	C12H22O2	000089-48-5	17



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ALS Vial : 6 Sample Multiplier: 1

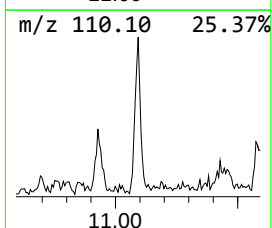
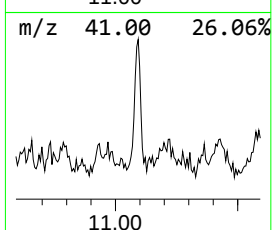
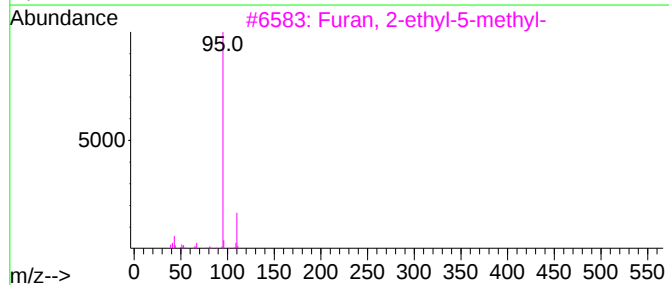
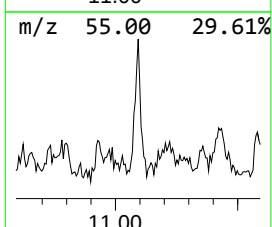
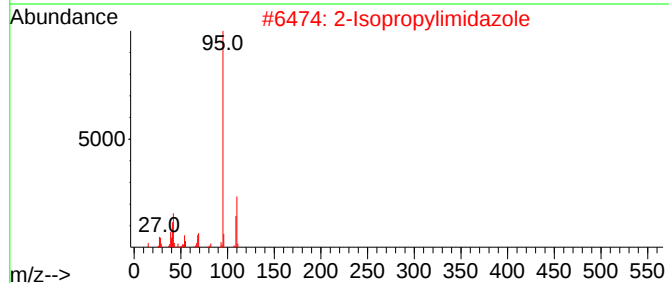
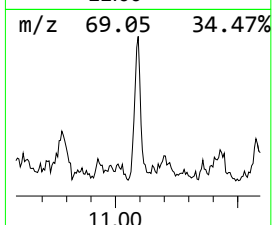
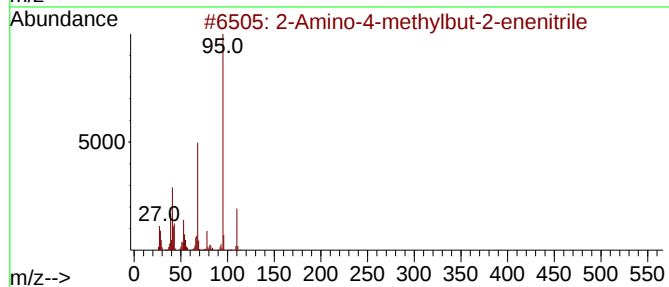
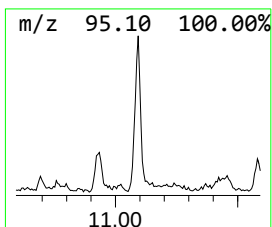
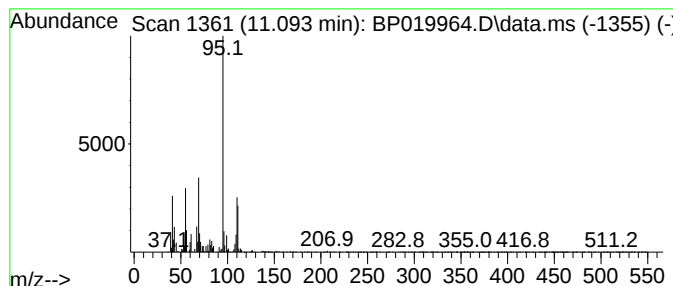
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Amino-4-methylbut-2-eneni... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.093	2.11 ng/ul	107074	Naphthalene-d8	10.369	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-4-methylbut-2-enenitrile	110	C6H10N2	1000197-39-9	59
2	2-Isopropylimidazole	110	C6H10N2	036947-68-9	58
3	Furan, 2-ethyl-5-methyl-	110	C7H10O	001703-52-2	58
4	1H-Imidazole, 2-ethyl-4-methyl-	110	C6H10N2	000931-36-2	53
5	endo-Borneol	154	C10H18O	000507-70-0	50



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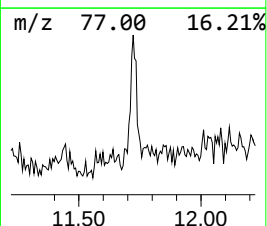
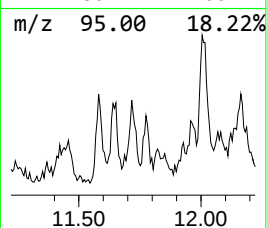
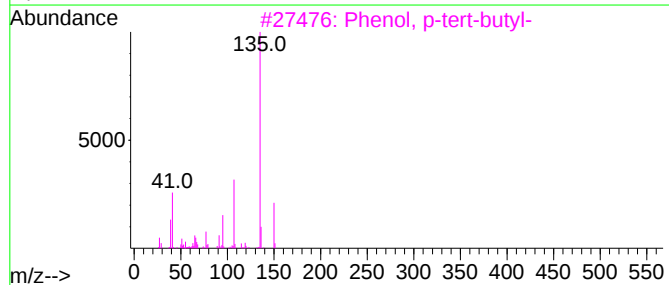
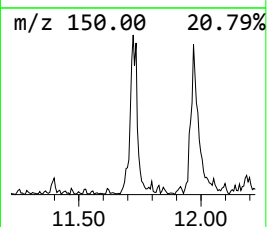
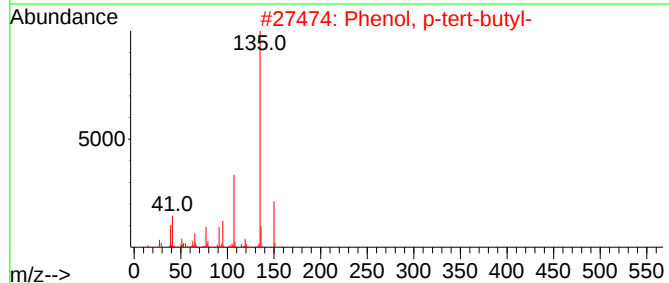
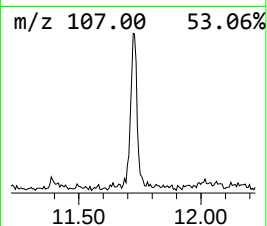
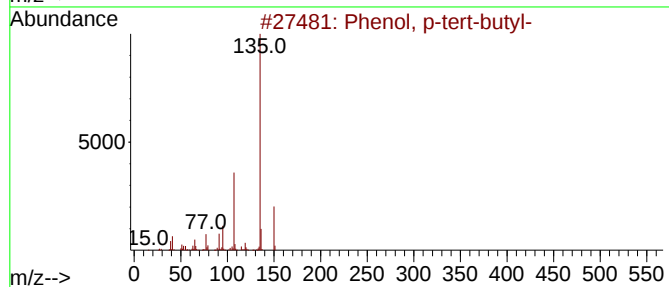
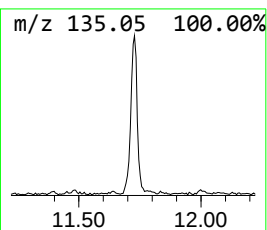
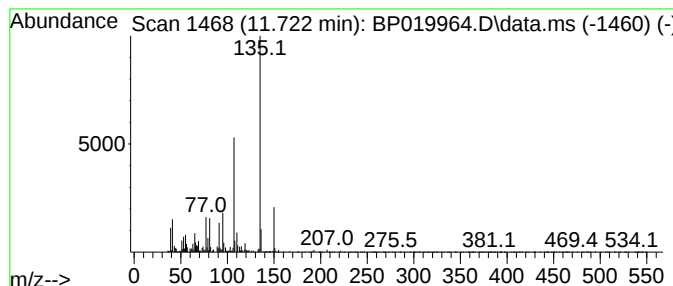
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Phenol, p-tert-butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.722	2.17 ng/ul	110093	Naphthalene-d8	10.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90
4			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	90
5			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	90



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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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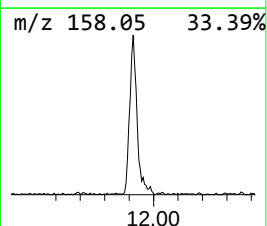
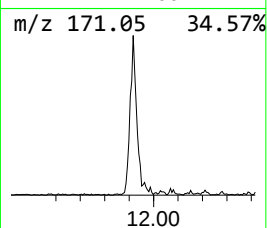
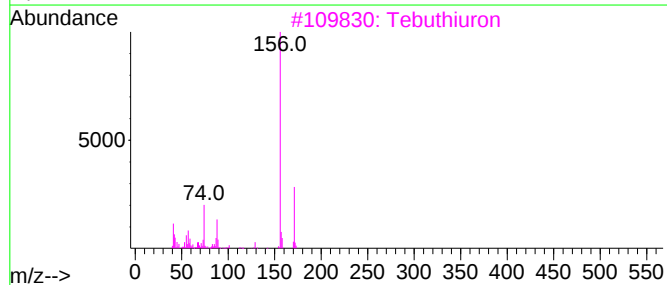
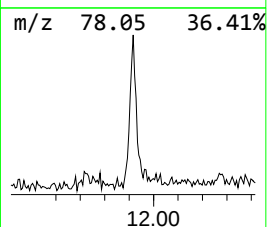
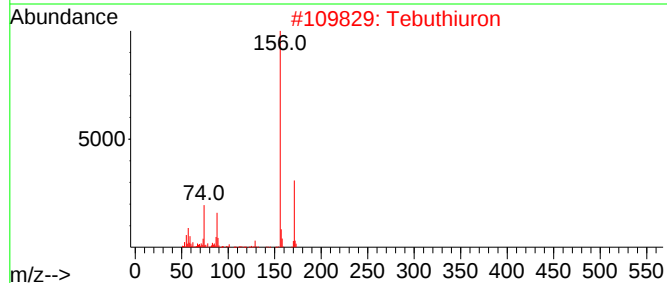
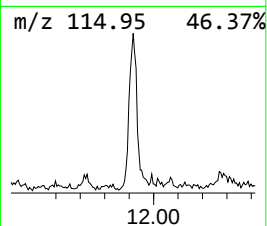
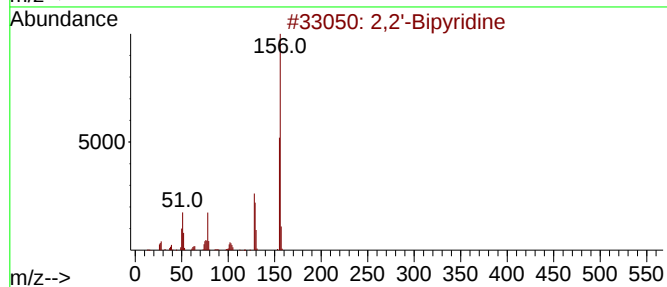
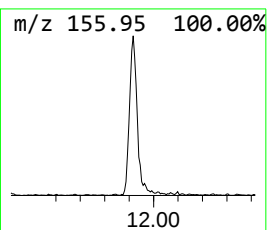
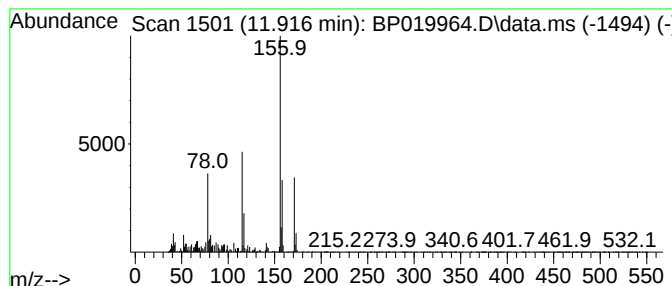
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	2.06 ng/ul	104566	Naphthalene-d8	10.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2'-Bipyridine	156	C10H8N2	000366-18-7	38
2			Tebuthiuron	228	C9H16N4OS	034014-18-1	37
3			Tebuthiuron	228	C9H16N4OS	034014-18-1	37
4			5,6-Dimethyl-2-thiouracil	156	C6H8N2OS	028456-54-4	35
5			Triacetyl-tetrahydro-veratramine...	553	C33H47NO6	1000256-71-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019964.D
Acq On : 22 Mar 2024 20:41
Operator : MA/JU
Sample : P1813-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0LX9

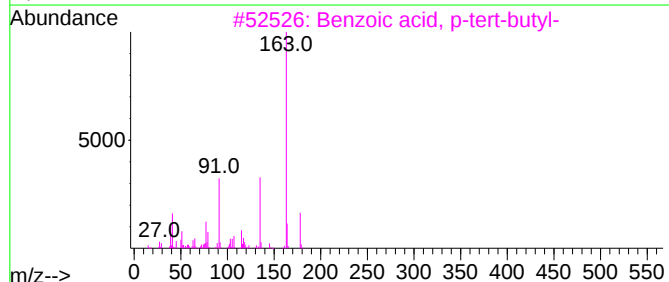
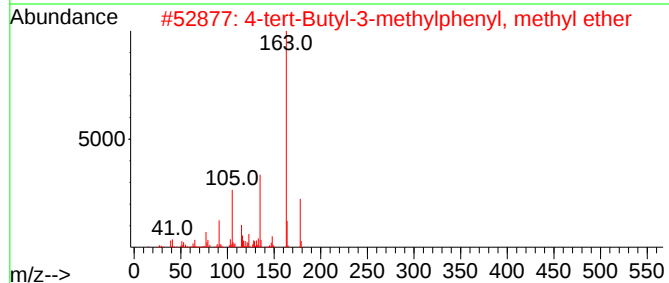
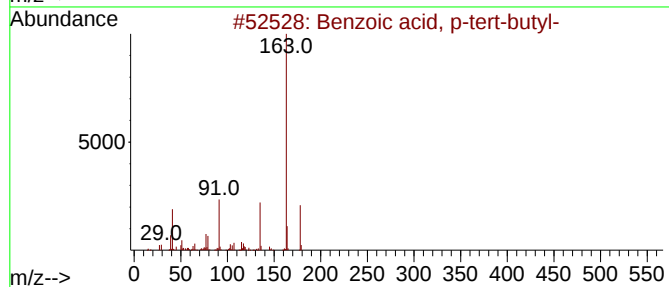
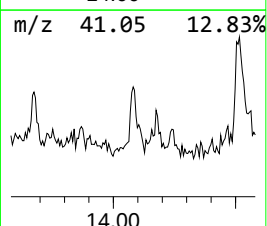
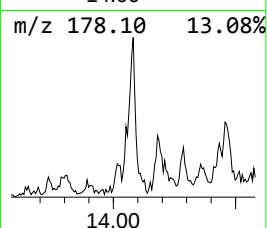
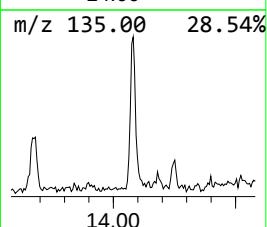
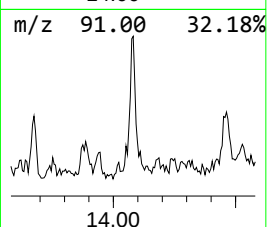
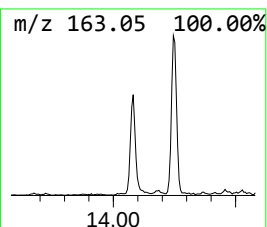
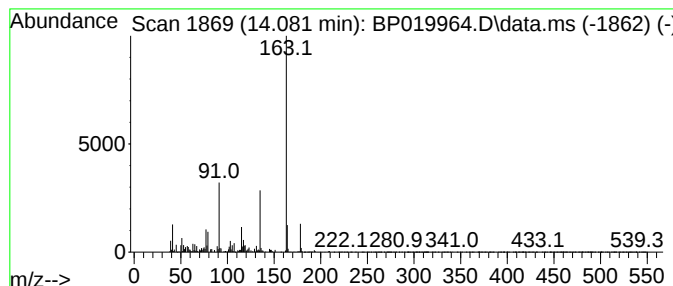
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzoic acid, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	2.42 ng/ul	153228	Acenaphthene-d10	14.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	91
2			4-tert-Butyl-3-methylphenyl, met...	178	C12H18O	031268-79-8	64
3			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	60
4			1,3-Benzenedicarboxylic acid, di...	194	C10H10O4	001459-93-4	59
5			6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019964.D
Acq On : 22 Mar 2024 20:41
Operator : MA/JU
Sample : P1813-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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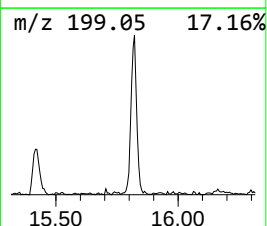
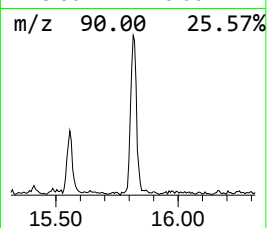
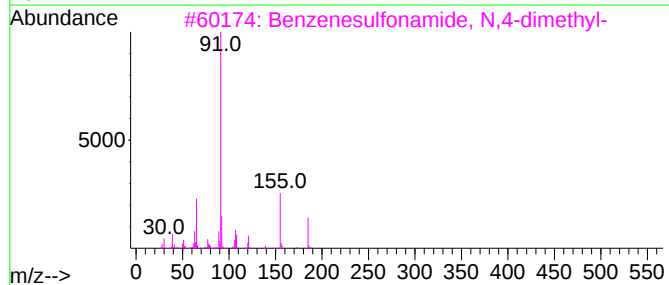
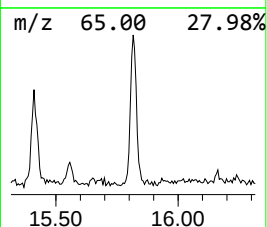
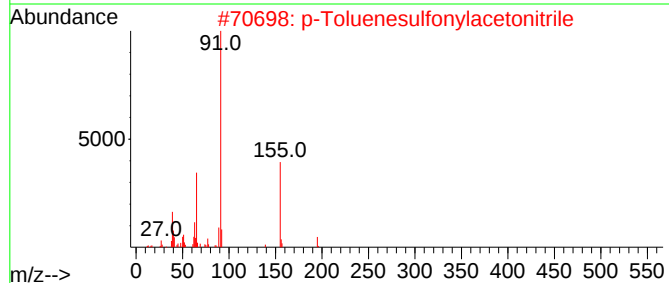
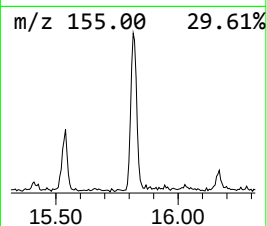
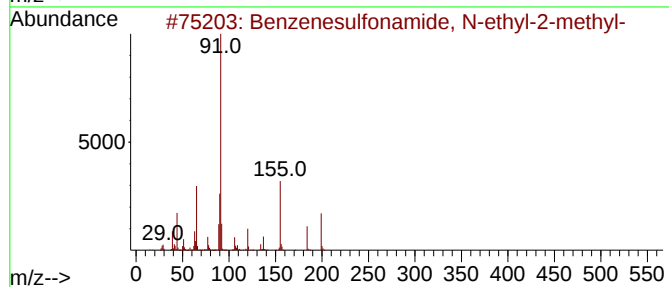
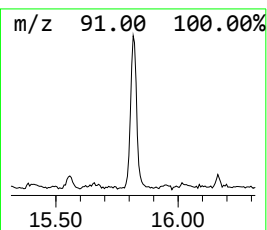
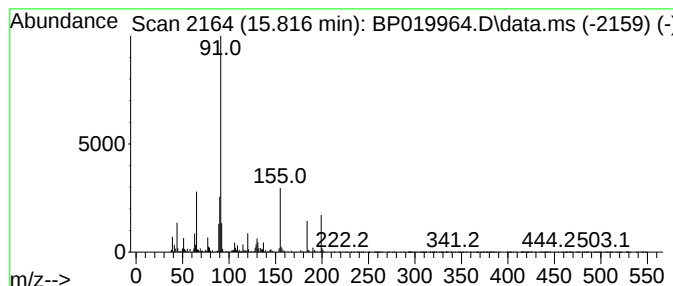
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.816	3.83 ng/ul	287139	Phenanthrene-d10	17.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2			p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	58
3			Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	50
4			Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	49
5			Benzenesulfonic acid, 3-methyl-,...	200	C9H12O3S	126865-08-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019964.D
Acq On : 22 Mar 2024 20:41
Operator : MA/JU
Sample : P1813-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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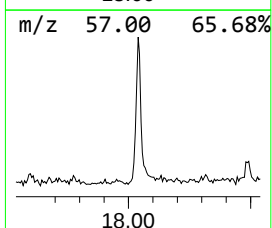
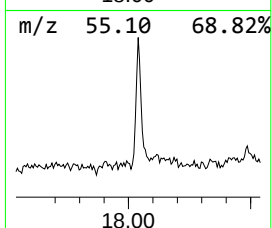
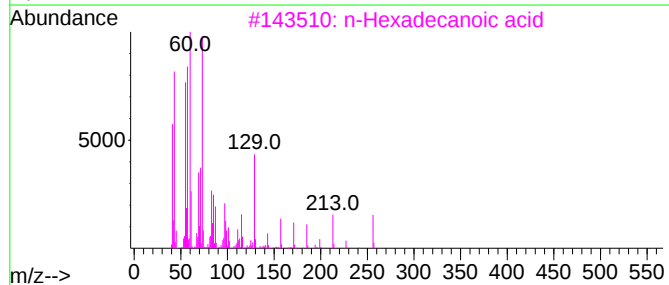
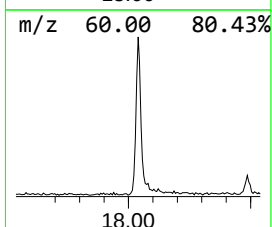
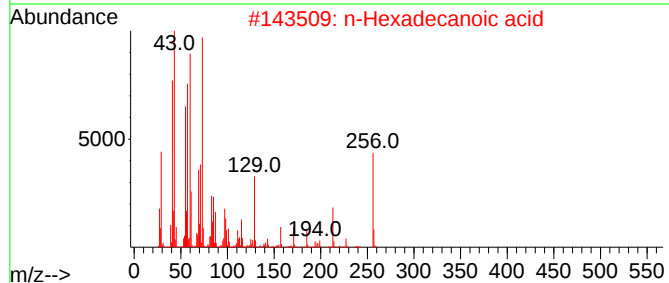
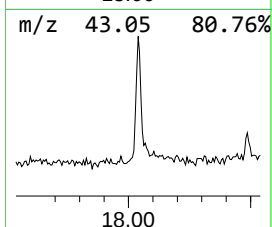
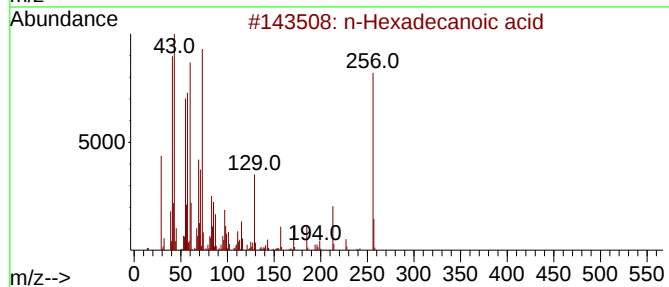
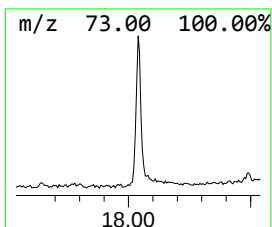
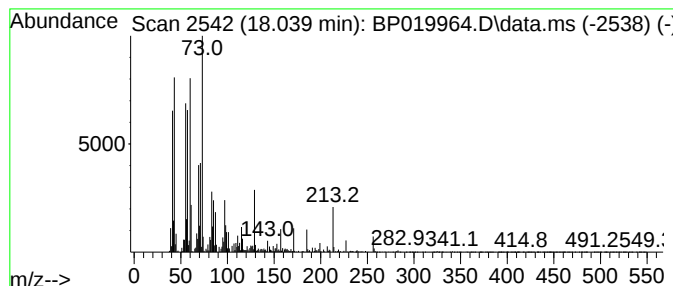
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.039	4.66 ng/ul	349493	Phenanthrene-d10	17.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			Tridecanoic acid	214	C13H26O2	000638-53-9	91
5			Octadecanoic acid	284	C18H36O2	000057-11-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019964.D
Acq On : 22 Mar 2024 20:41
Operator : MA/JU
Sample : P1813-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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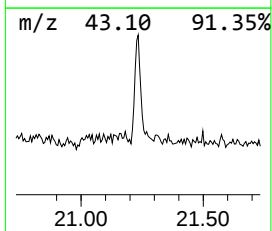
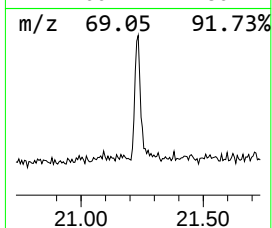
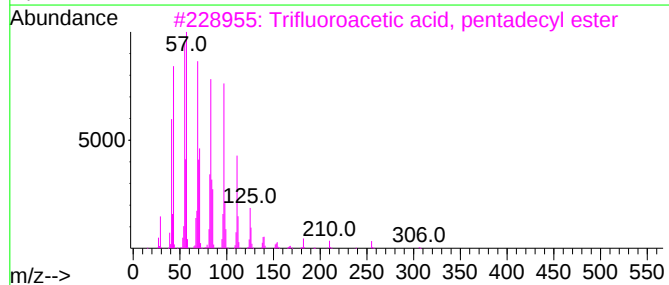
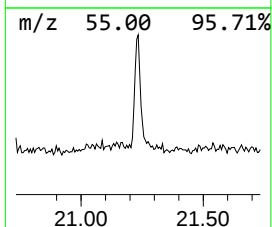
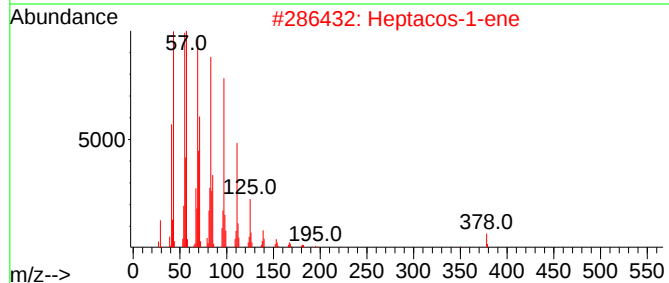
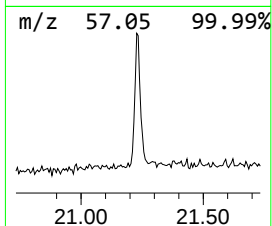
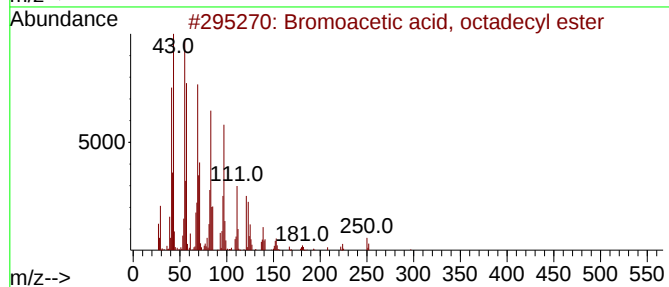
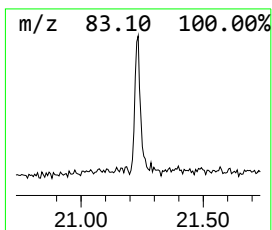
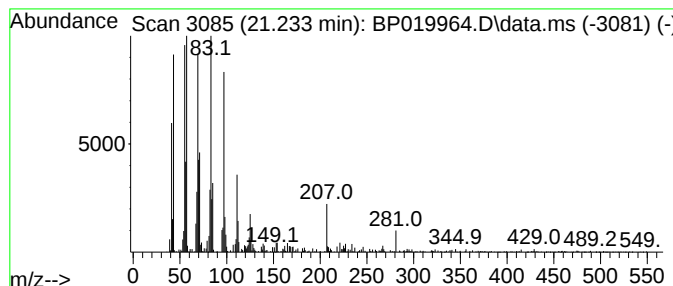
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Bromoacetic acid, octadecyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.233	2.93 ng/ul	235578	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	87
2			Heptacos-1-ene	378	C27H54	015306-27-1	81
3			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	74
4			1-Tricosene	322	C23H46	018835-32-0	70
5			Pentacos-1-ene	350	C25H50	016980-85-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019964.D
Acq On : 22 Mar 2024 20:41
Operator : MA/JU
Sample : P1813-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.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
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2-Amino-4-methy...	11.093	2.1	ng/ul	107074	2	10.369	1015160	20.0
Phenol, p-tert-...	11.722	2.2	ng/ul	110093	2	10.369	1015160	20.0
unknown-02	11.916	2.1	ng/ul	104566	2	10.369	1015160	20.0
Benzoic acid, p...	14.081	2.4	ng/ul	153228	3	14.246	1268390	20.0
Benzenesulfonam...	15.816	3.8	ng/ul	287139	4	17.081	1499610	20.0
n-Hexadecanoic ...	18.039	4.7	ng/ul	349493	4	17.081	1499610	20.0
Bromoacetic aci...	21.233	2.9	ng/ul	235578	5	21.551	1608890	20.0