

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
 Data File : BP006350.D
 Acq On : 26 Jul 2021 22:00
 Operator : CG/JU
 Sample : M3091-20
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0058

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-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006350.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.305	33	37	50	rVB	71105	106183	1.76%	0.150%
2	3.711	102	106	122	rVB	382589	622719	10.33%	0.879%
3	4.028	155	160	166	rVB	24751	39553	0.66%	0.056%
4	4.940	311	315	321	rVB	25650	36608	0.61%	0.052%
5	5.858	466	471	475	rBV	88900	132586	2.20%	0.187%
6	5.899	475	478	484	rVB	24063	34171	0.57%	0.048%
7	6.452	567	572	579	rBV	31535	50213	0.83%	0.071%
8	6.958	652	658	669	rVB	308059	502182	8.33%	0.709%
9	7.128	681	687	700	rVB	929711	1485977	24.65%	2.098%
10	7.322	711	720	729	rBV	963190	1529664	25.38%	2.159%
11	7.416	731	736	742	rVV8	9704	20042	0.33%	0.028%
12	7.587	761	765	771	rBV2	10317	19873	0.33%	0.028%
13	7.787	790	799	806	rBV	982402	1525499	25.31%	2.154%
14	7.852	806	810	822	rVB5	31740	62141	1.03%	0.088%
15	8.028	833	840	845	rBV5	10643	23303	0.39%	0.033%
16	8.493	909	919	926	rVV	802262	1245024	20.65%	1.758%
17	8.581	929	934	936	rVV	42100	63643	1.06%	0.090%
18	8.611	936	939	945	rVB	41400	61822	1.03%	0.087%
19	8.881	978	985	990	rBV	282512	458214	7.60%	0.647%
20	8.940	990	995	1002	rVB	1103673	1731525	28.73%	2.444%
21	9.040	1007	1012	1021	rVV3	89880	174275	2.89%	0.246%
22	9.116	1021	1025	1033	rVB3	23090	37010	0.61%	0.052%
23	9.658	1108	1117	1129	rBV	827507	1322137	21.93%	1.866%
24	9.752	1129	1133	1137	rVV	17054	26145	0.43%	0.037%
25	9.840	1137	1148	1155	rVV2	445846	1085896	18.01%	1.533%
26	10.081	1179	1189	1197	rBV	634505	1001835	16.62%	1.414%
27	10.193	1201	1208	1219	rVB	1243693	1968428	32.66%	2.779%
28	10.363	1230	1237	1245	rBV	437235	676908	11.23%	0.956%
29	10.440	1245	1250	1253	rVV4	23026	37065	0.61%	0.052%
30	10.493	1253	1259	1265	rVV3	40058	75069	1.25%	0.106%
31	10.569	1265	1272	1279	rVV	1301692	2151309	35.69%	3.037%
32	10.634	1279	1283	1289	rVV4	46966	84191	1.40%	0.119%
33	10.710	1289	1296	1311	rVB	1248711	2059757	34.17%	2.908%
34	10.875	1318	1324	1329	rBV7	11972	20596	0.34%	0.029%
35	11.122	1355	1366	1371	rBV	153283	246175	4.08%	0.348%
36	11.181	1371	1376	1381	rVV	168729	272155	4.52%	0.384%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: 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-BP072421.M
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37	11.222	1381	1383	1389	rVB	37767	52012	0.86%	0.073%
38	11.752	1461	1473	1481	rBV4	14392	35705	0.59%	0.050%
39	12.004	1511	1516	1521	rVB5	21035	37527	0.62%	0.053%
40	12.104	1523	1533	1538	rBV2	77744	145841	2.42%	0.206%
41	12.157	1538	1542	1546	rVV	25102	37374	0.62%	0.053%
42	12.204	1546	1550	1555	rVB	18211	27614	0.46%	0.039%
43	12.775	1642	1647	1654	rVB	83786	128900	2.14%	0.182%
44	13.034	1685	1691	1697	rBV	131980	192804	3.20%	0.272%
45	13.263	1725	1730	1737	rVV	38943	59916	0.99%	0.085%
46	13.334	1737	1742	1748	rVV	68733	112559	1.87%	0.159%
47	13.657	1789	1797	1808	rBV2	51996	94938	1.58%	0.134%
48	13.834	1815	1827	1837	rBV	2344731	3221097	53.44%	4.547%
49	13.951	1843	1847	1852	rVB6	20839	31657	0.53%	0.045%
50	14.104	1862	1873	1886	rBV	2647149	3750784	62.23%	5.295%
51	14.275	1899	1902	1907	rVV3	14140	24678	0.41%	0.035%
52	14.322	1907	1910	1916	rVB4	14992	24472	0.41%	0.035%
53	14.410	1916	1925	1933	rBV	1796779	2735205	45.38%	3.861%
54	14.504	1936	1941	1945	rVB5	19573	28799	0.48%	0.041%
55	14.610	1951	1959	1967	rBV	294493	432756	7.18%	0.611%
56	14.840	1992	1998	2002	rBV5	17592	25980	0.43%	0.037%
57	15.016	2024	2028	2032	rVV3	21236	27197	0.45%	0.038%
58	15.104	2039	2043	2048	rVV3	13101	24034	0.40%	0.034%
59	15.169	2048	2054	2059	rVV	17541	32607	0.54%	0.046%
60	15.228	2059	2064	2070	rVV2	31772	59258	0.98%	0.084%
61	15.281	2070	2073	2077	rVB2	19539	25350	0.42%	0.036%
62	15.410	2088	2095	2102	rBV	3331745	4731934	78.50%	6.680%
63	15.522	2107	2114	2129	rBV	995531	1350889	22.41%	1.907%
64	15.645	2130	2135	2140	rBV	39370	58662	0.97%	0.083%
65	15.781	2153	2158	2163	rVB3	31507	45428	0.75%	0.064%
66	16.045	2199	2203	2206	rBV5	15252	21934	0.36%	0.031%
67	16.098	2206	2212	2217	rVV6	26017	60040	1.00%	0.085%
68	16.204	2226	2230	2236	rVB7	19617	35728	0.59%	0.050%
69	16.316	2244	2249	2253	rBV6	15439	29226	0.48%	0.041%
70	16.363	2253	2257	2261	rVB5	16757	25183	0.42%	0.036%
71	16.404	2261	2264	2270	rBV2	30240	45579	0.76%	0.064%
72	16.563	2287	2291	2297	rVB	79285	107712	1.79%	0.152%
73	16.634	2298	2303	2306	rVV	25576	37327	0.62%	0.053%
74	16.669	2306	2309	2313	rVV4	27442	40884	0.68%	0.058%
75	16.722	2314	2318	2322	rVV4	16302	26151	0.43%	0.037%
76	16.769	2324	2326	2336	rVB7	12751	20372	0.34%	0.029%

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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	16.904	2345	2349	2353	rBV7	13712	28095	0.47%	0.040%
78	16.945	2353	2356	2359	rVV4	25140	31682	0.53%	0.045%
79	16.992	2360	2364	2371	rVB4	16434	29095	0.48%	0.041%
80	17.163	2386	2393	2402	rBV	2161244	3068983	50.91%	4.332%
81	17.263	2402	2410	2417	rBV	3901552	5515306	91.50%	7.786%
82	17.316	2417	2419	2424	rVB	43176	49917	0.83%	0.070%
83	17.792	2496	2500	2503	rBV3	23775	29659	0.49%	0.042%
84	17.845	2503	2509	2516	rBV	1722737	2067917	34.31%	2.919%
85	17.963	2525	2529	2536	rVB	52611	74479	1.24%	0.105%
86	18.069	2542	2547	2552	rBV	451196	581650	9.65%	0.821%
87	18.157	2559	2562	2570	rVB7	34294	58812	0.98%	0.083%
88	18.269	2578	2581	2585	rBV2	27476	33955	0.56%	0.048%
89	18.804	2668	2672	2675	rBV	28568	37524	0.62%	0.053%
90	18.892	2683	2687	2690	rBV	39856	51347	0.85%	0.072%
91	19.081	2715	2719	2724	rBV2	48058	65767	1.09%	0.093%
92	19.157	2727	2732	2736	rBV3	44928	88168	1.46%	0.124%
93	19.304	2753	2757	2764	rBV	300584	393035	6.52%	0.555%
94	19.445	2778	2781	2785	rVB	49114	53166	0.88%	0.075%
95	19.504	2785	2791	2798	rBV	4684356	6027740	100.00%	8.509%
96	20.986	3039	3043	3050	rBV	442235	562198	9.33%	0.794%
97	21.257	3084	3089	3097	rBV	2780483	3355015	55.66%	4.736%
98	23.451	3455	3462	3470	rBV2	3025812	5948779	98.69%	8.398%
99	23.539	3473	3477	3481	rVV2	201309	332857	5.52%	0.470%
100	23.604	3481	3488	3499	rVB	1734327	3356123	55.68%	4.738%

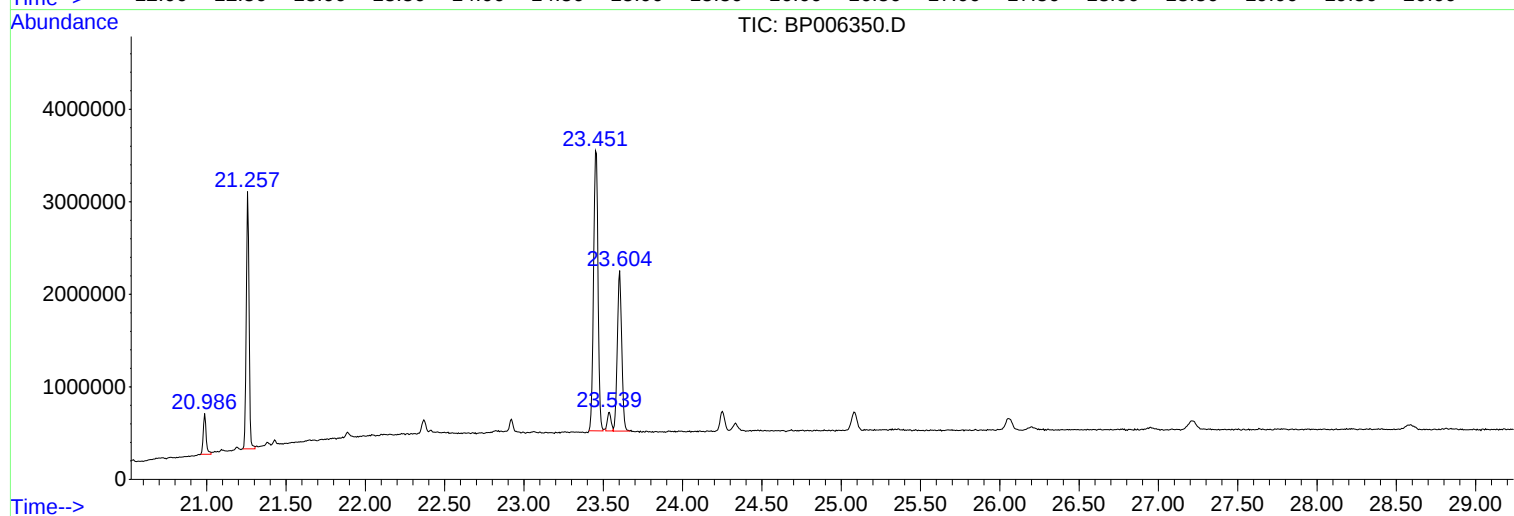
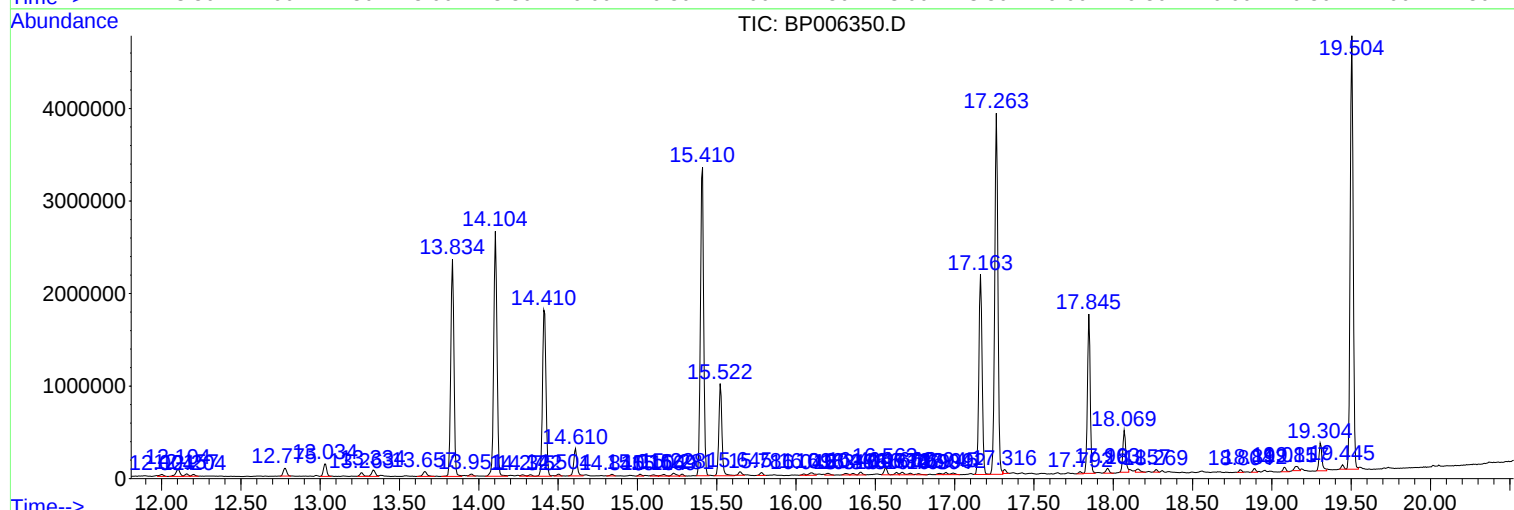
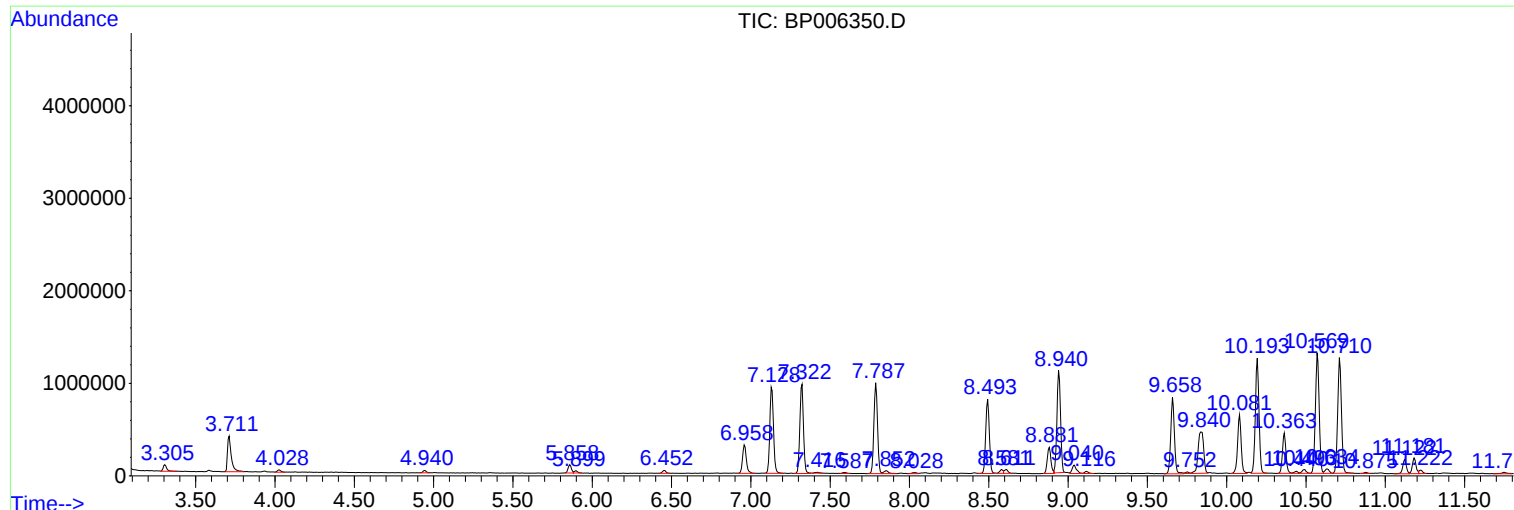
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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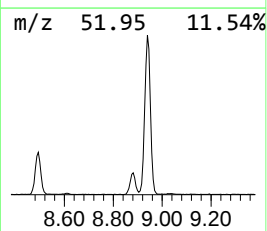
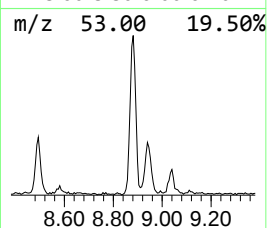
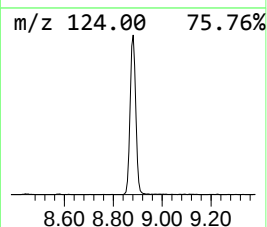
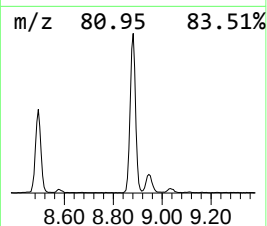
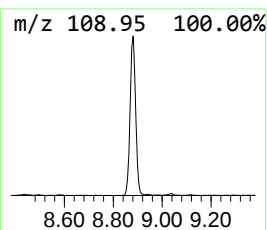
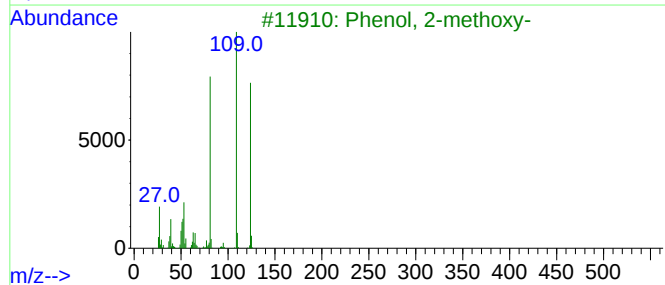
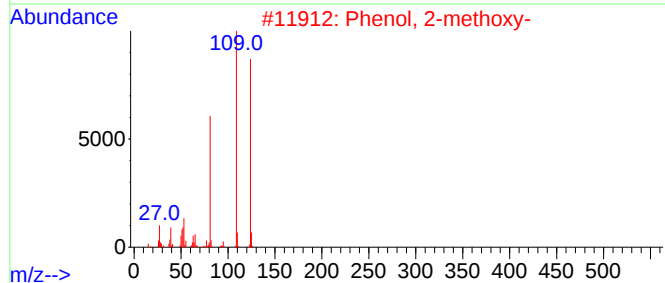
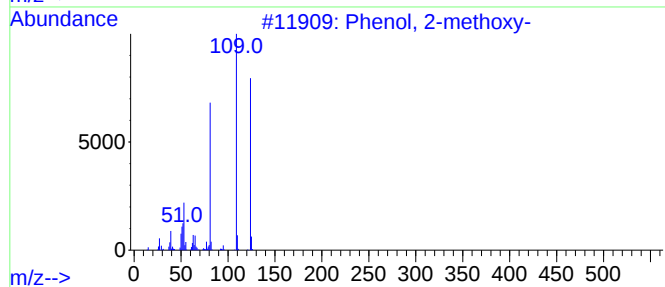
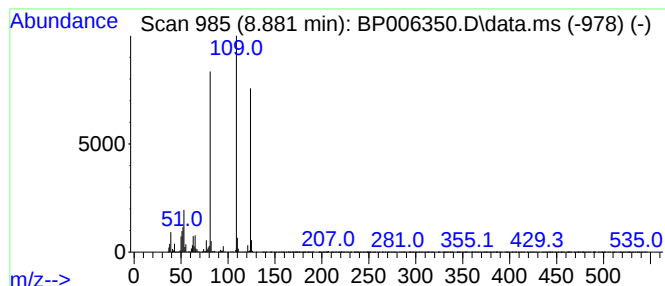
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 2-methoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.880	6.01 ng/ul	458214	1,4-Dichlorobenzene-d4	7.787

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	91
3			Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	91
4			Mequinol	124	C7H8O2	000150-76-5	91
5			Mequinol	124	C7H8O2	000150-76-5	64



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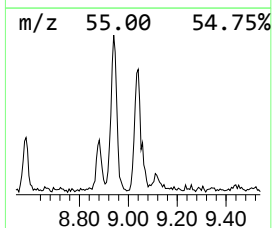
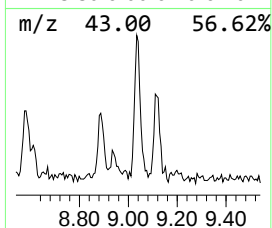
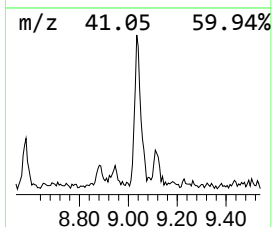
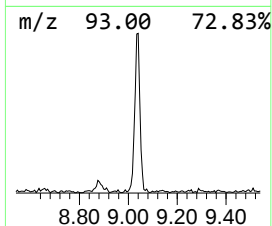
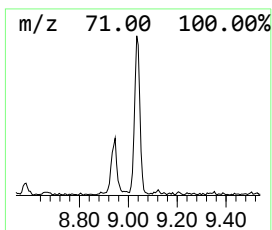
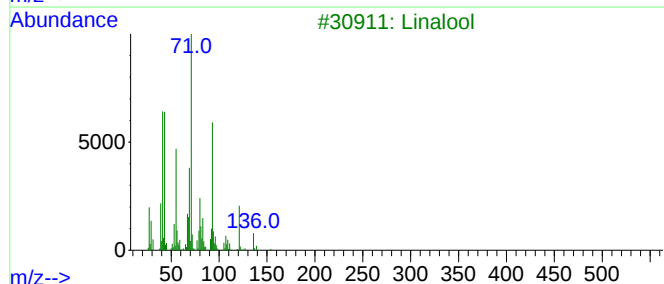
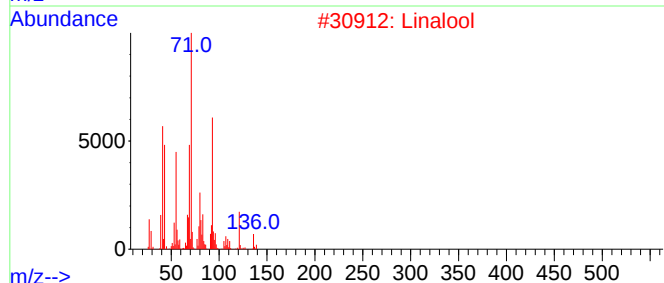
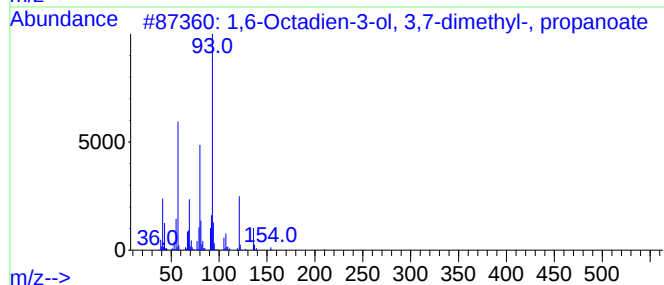
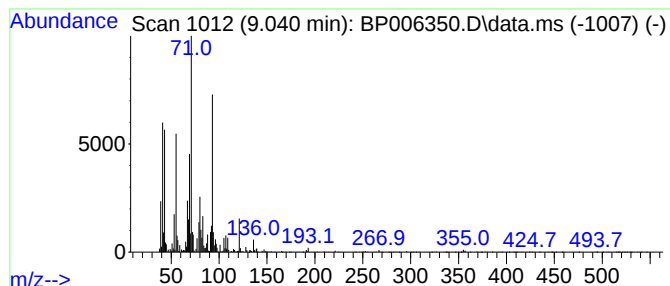
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,6-Octadien-3-ol, 3,7-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.040	2.28 ng/ul	174275	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6-Octadien-3-ol, 3,7-dimethyl-...	210	C13H22O2	000144-39-8	83		
2	Linalool	154	C10H18O	000078-70-6	80		
3	Linalool	154	C10H18O	000078-70-6	72		
4	1,5-Dimethyl-1-vinyl-4-hexenyl b...	224	C14H24O2	000078-36-4	43		
5	1,6-Octadien-3-ol, 3,7-dimethyl-...	182	C11H18O2	000115-99-1	43		



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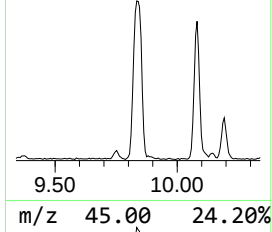
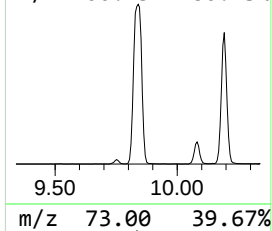
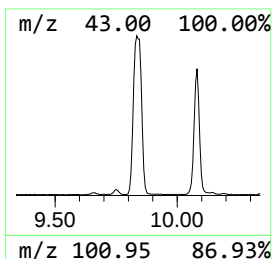
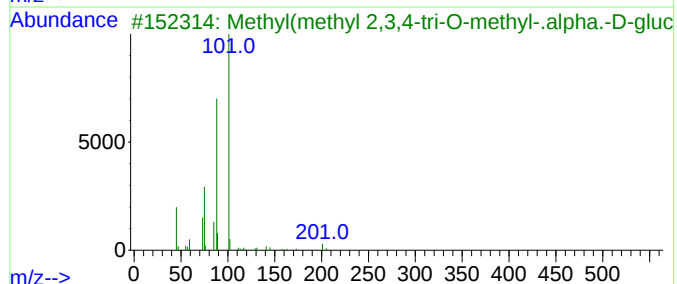
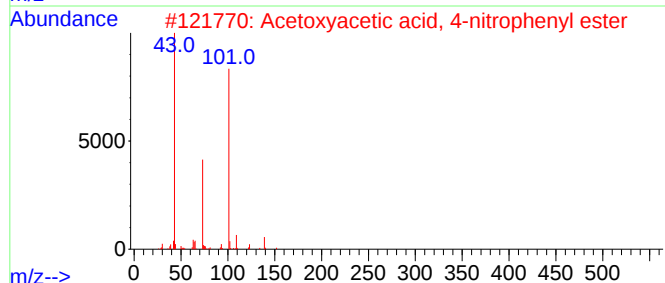
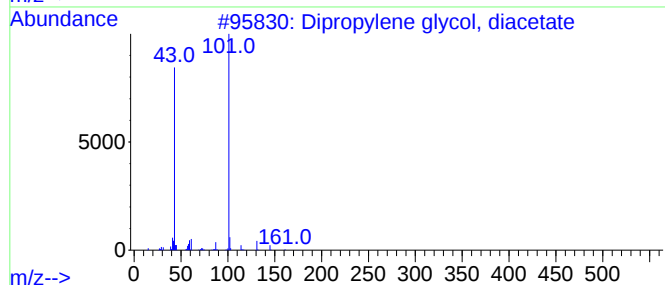
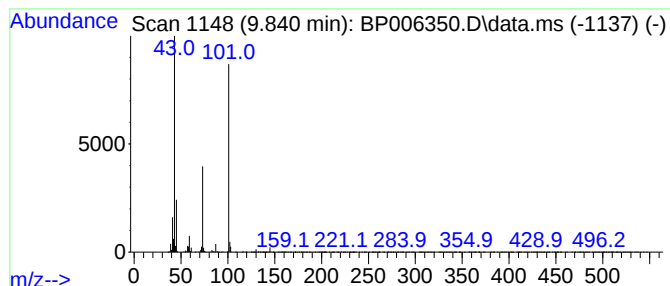
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Dipropylene glycol, diacetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.840	10.10 ng/ul	1085900	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dipropylene glycol, diacetate	218	C10H18O5	1000506-25-8	59
2			Acetoxyacetic acid, 4-nitropheny...	239	C10H9NO6	1000307-54-5	45
3			Methyl(methyl 2,3,4-tri-O-methyl...	264	C11H20O7	052729-97-2	43
4			2,3,4-Trimethyl-d-xylose	192	C8H16O5	007434-17-5	39
5			1,3-Dioxolane-2-acetic acid, 2,4...	188	C9H16O4	090293-83-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
Operator : CG/JU
Sample : M3091-20
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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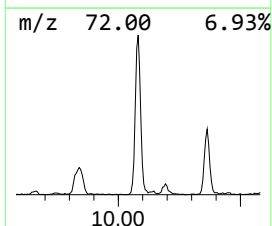
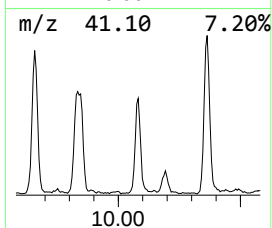
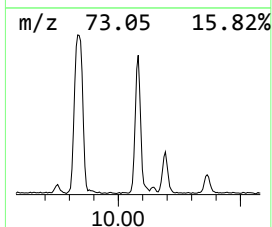
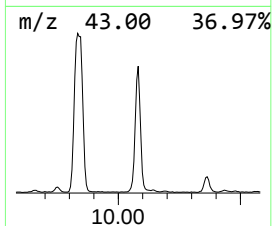
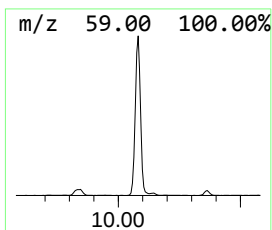
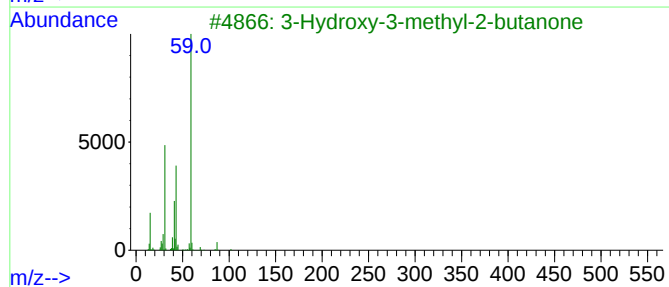
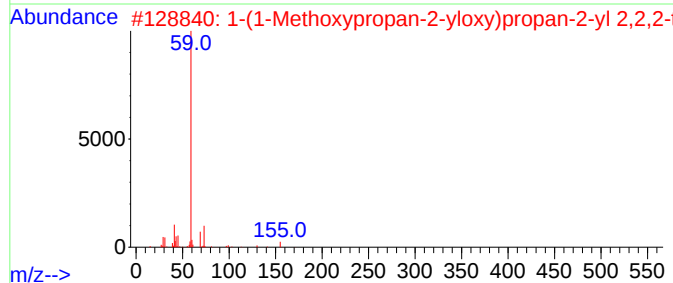
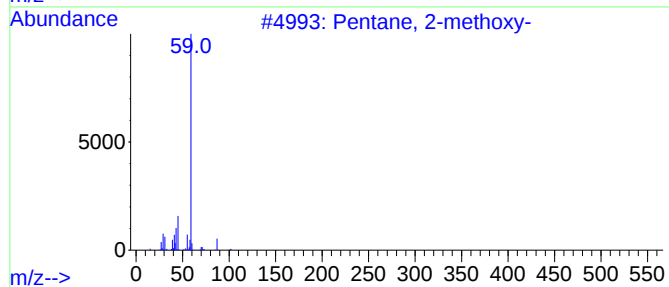
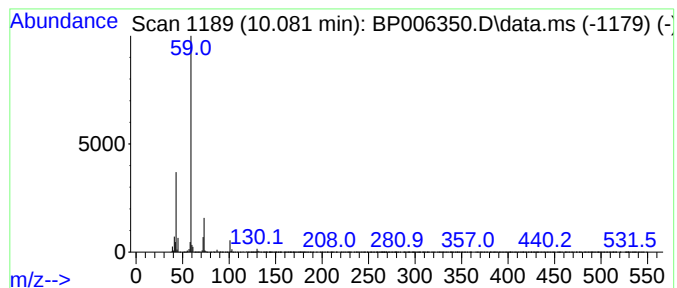
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Pentane, 2-methoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.080	9.31 ng/ul	1001840	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methoxy-	102	C6H14O	006795-88-6	59
2			1-(1-Methoxypropan-2-yloxy)propa...	244	C9H15F3O4	1010367-08-7	56
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	53
4			1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4	1000367-08-6	47
5			2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
Operator : CG/JU
Sample : M3091-20
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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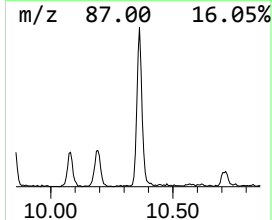
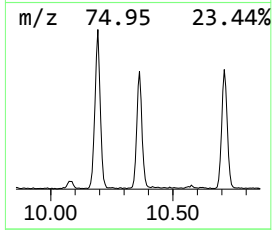
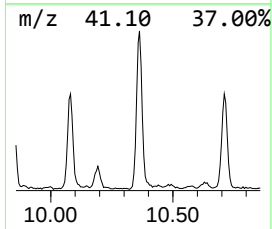
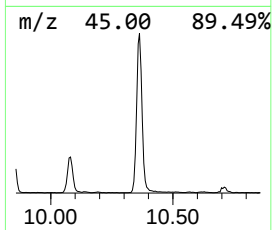
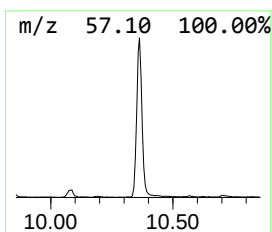
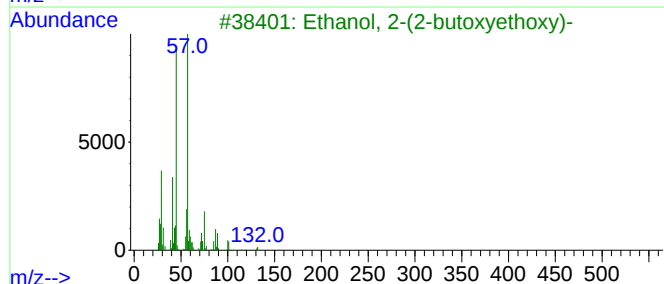
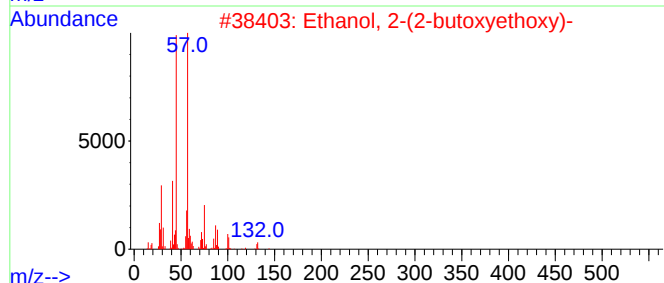
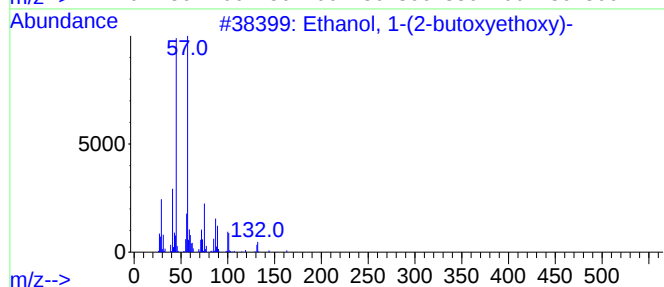
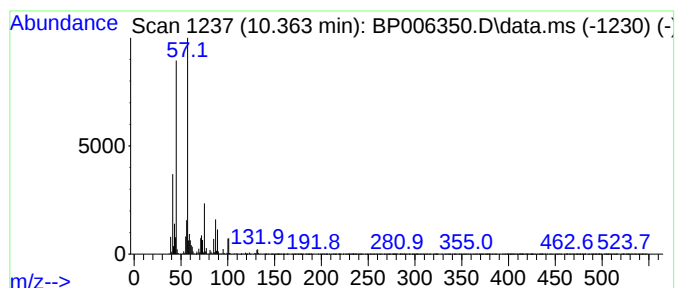
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	6.29 ng/ul	676908	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
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ClientSampleId :
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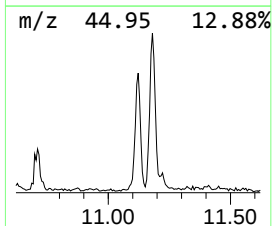
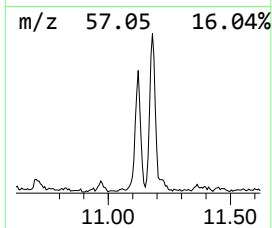
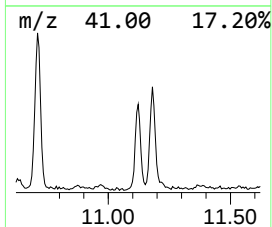
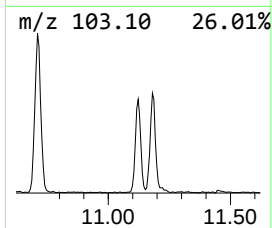
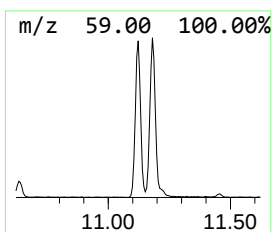
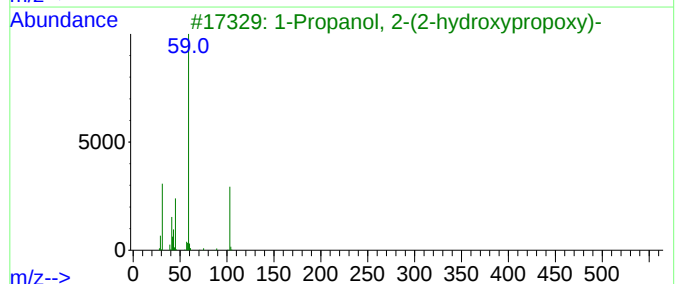
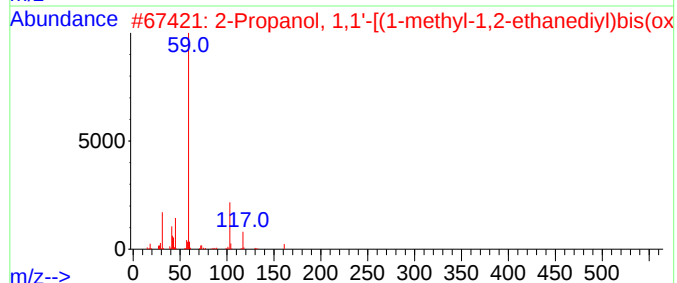
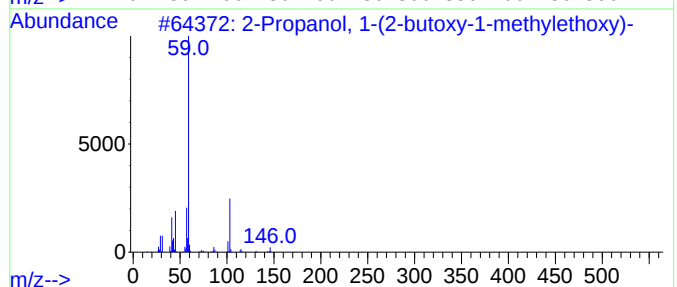
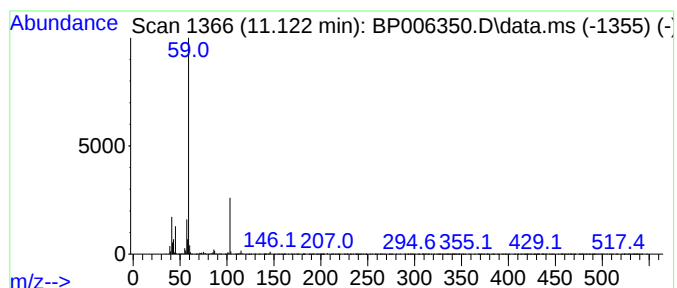
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.120	2.29 ng/ul	246175	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83		
2	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	72		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	72		
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		
5	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
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Sample : M3091-20
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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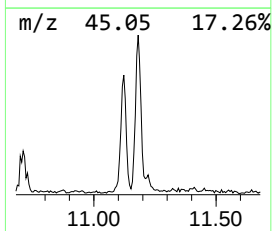
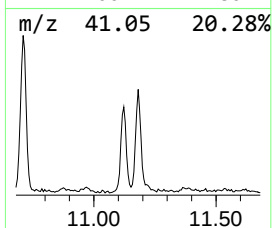
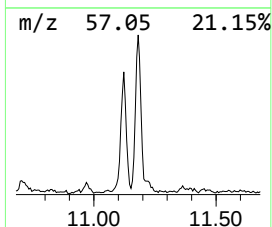
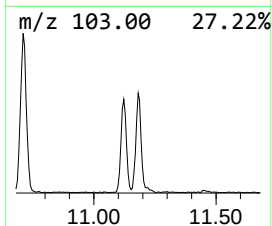
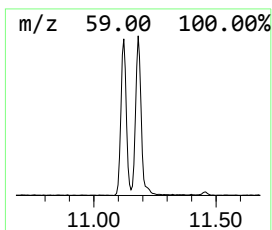
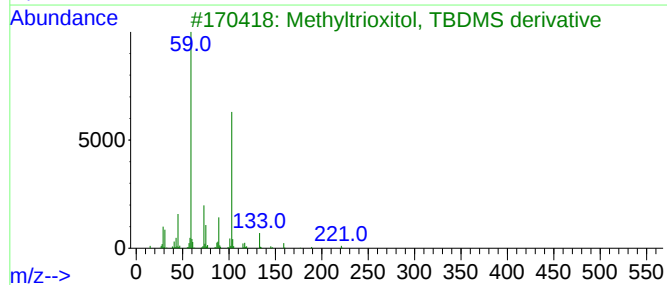
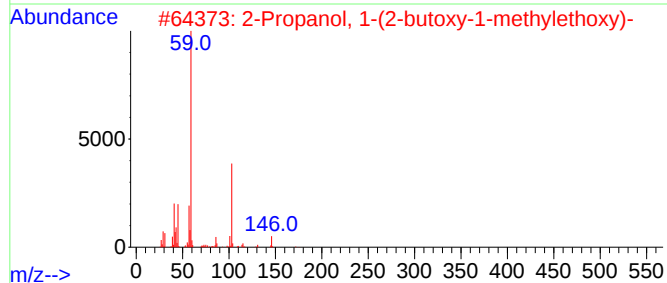
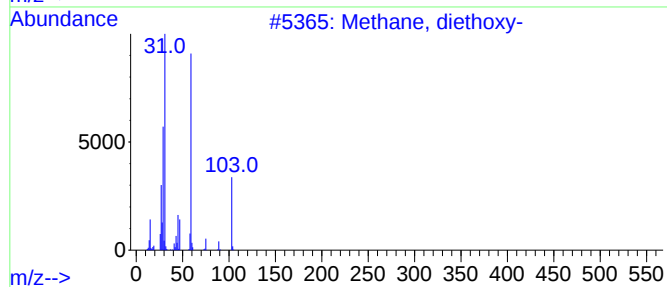
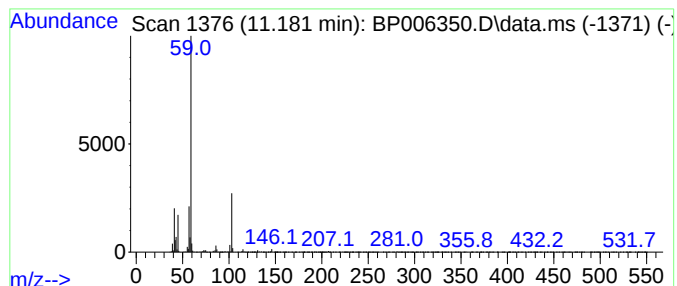
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methane, diethoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.180	2.53 ng/ul	272155	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methane, diethoxy-	104	C5H12O2	000462-95-3	86
2			2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83
3			Methyltrioxitol, TBDMS derivative	278	C13H30O4Si	1000352-09-5	83
4			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
5			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
Operator : CG/JU
Sample : M3091-20
Misc :
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Instrument :
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ClientSampleId :
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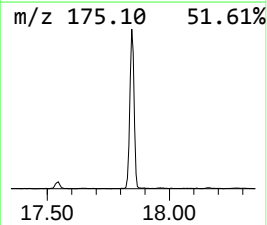
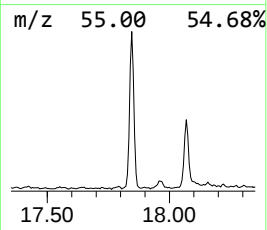
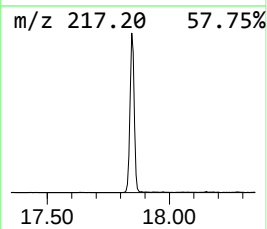
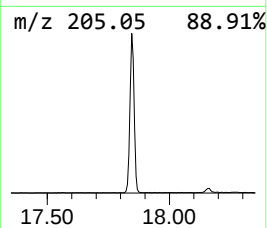
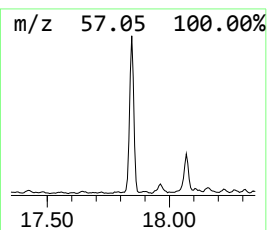
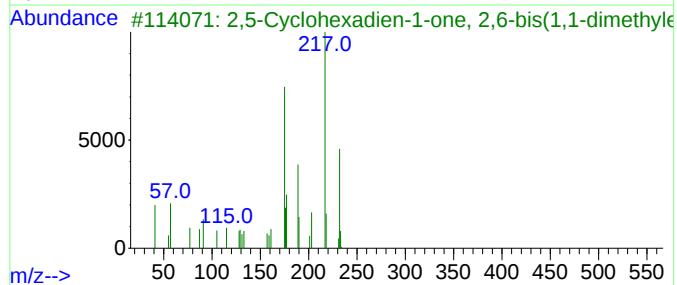
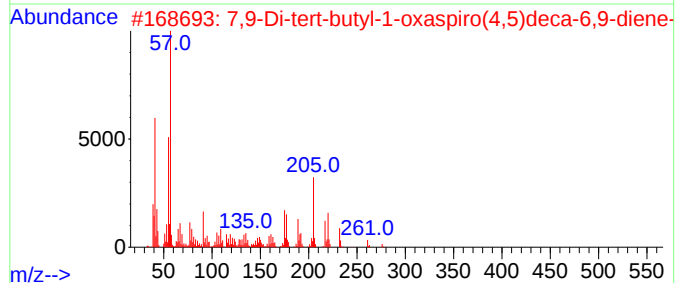
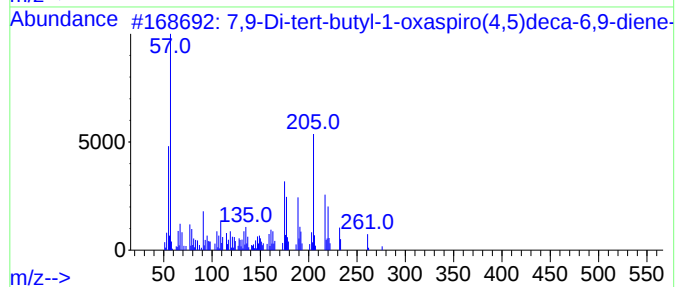
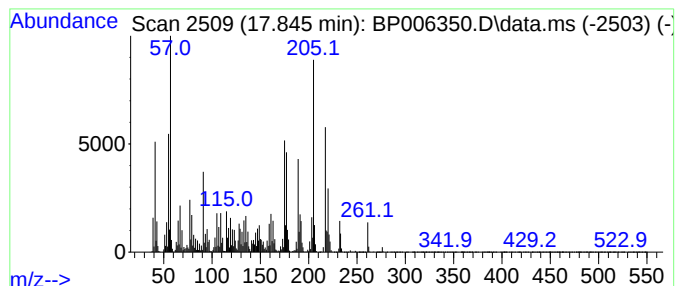
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	13.48 ng/ul	2067920	Phenanthrene-d10	17.163

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	83		
4	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	68		
5	1,3-Pentadiene, 1,1-diphenyl-, (Z)-	220	C17H16	015295-31-5	51		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
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Acq On : 26 Jul 2021 22:00
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ClientSampleId :
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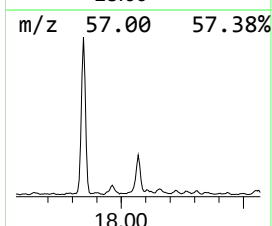
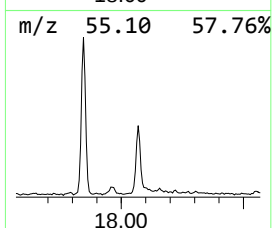
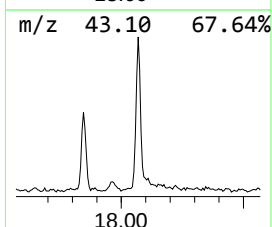
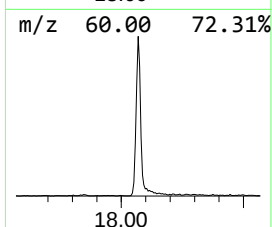
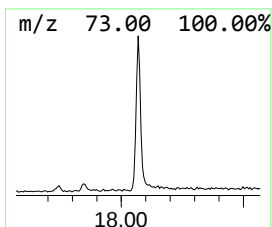
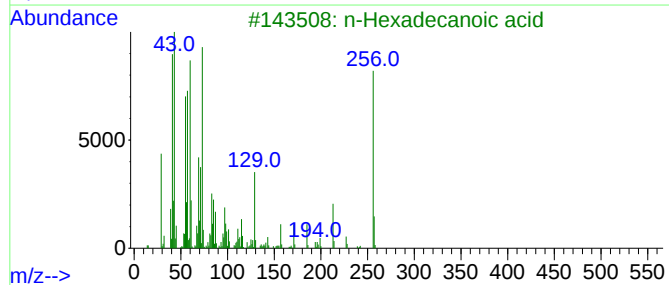
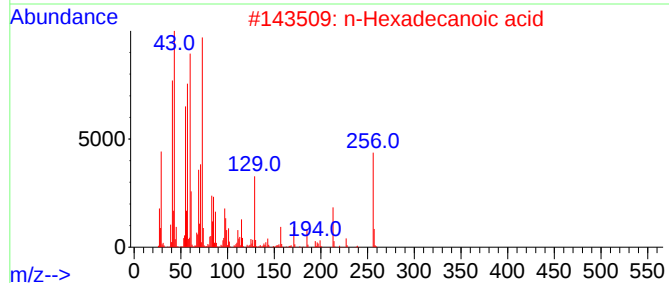
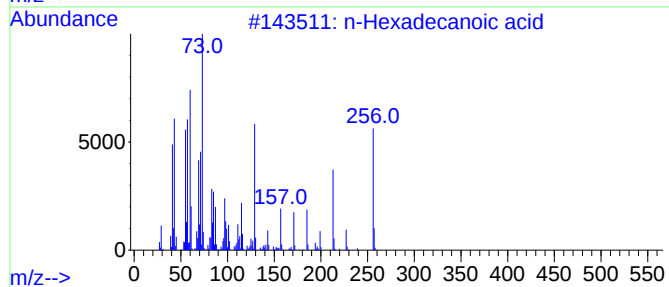
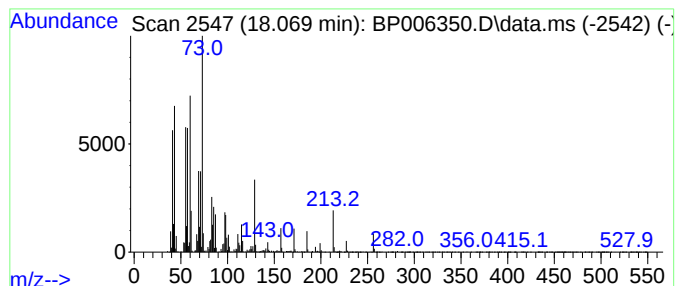
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	3.79 ng/ul	581650	Phenanthrene-d10	17.163

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
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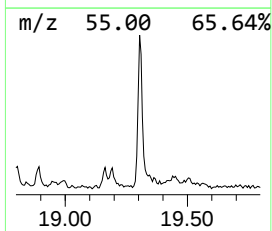
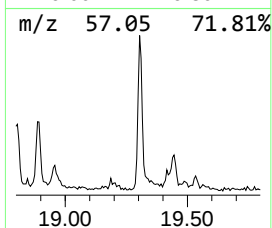
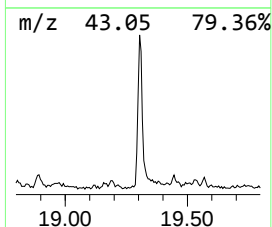
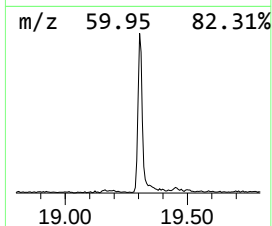
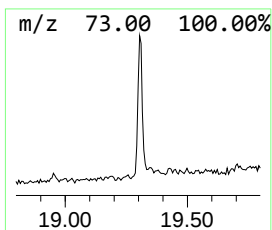
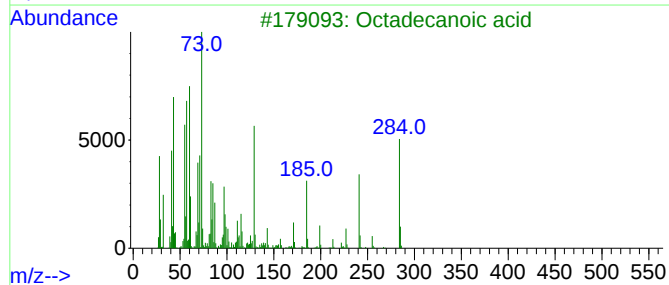
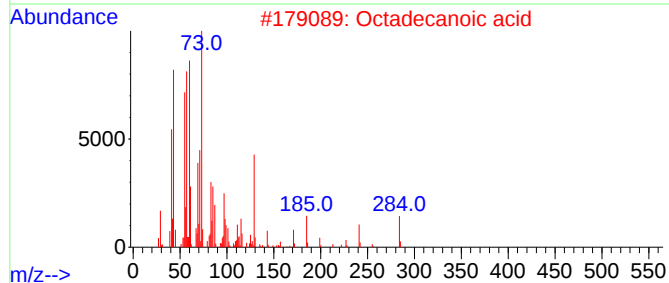
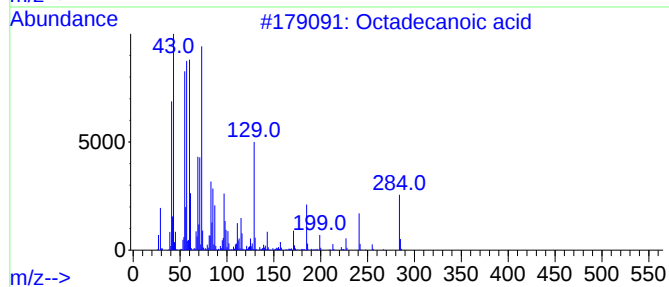
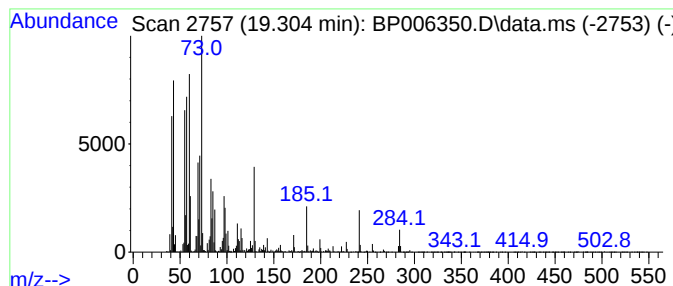
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.300	2.34 ng/ul	393035	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
Data File : BP006350.D
Acq On : 26 Jul 2021 22:00
Operator : CG/JU
Sample : M3091-20
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0058

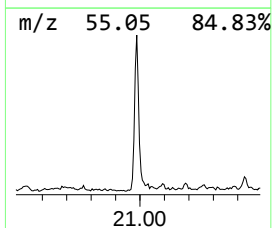
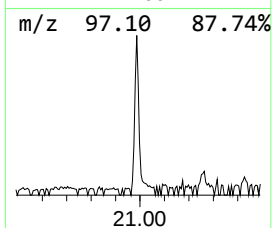
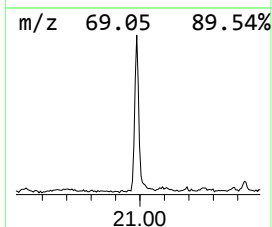
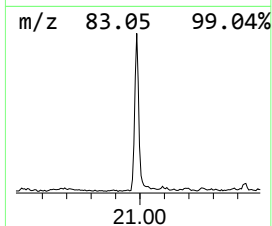
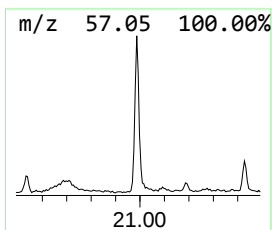
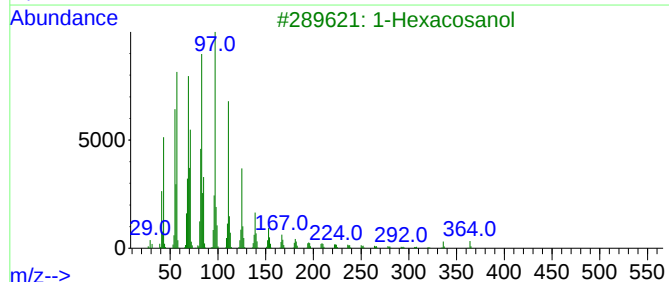
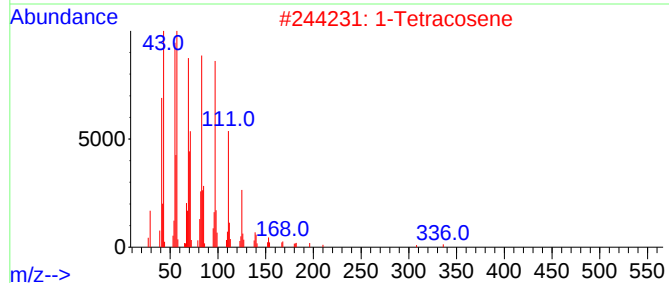
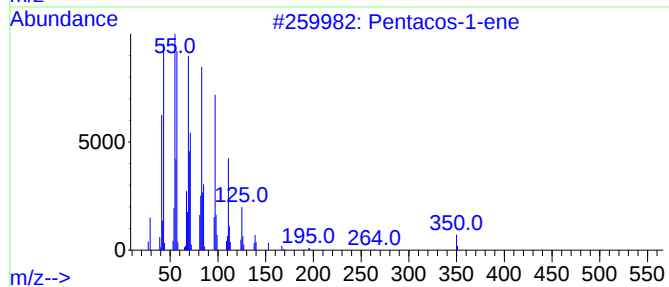
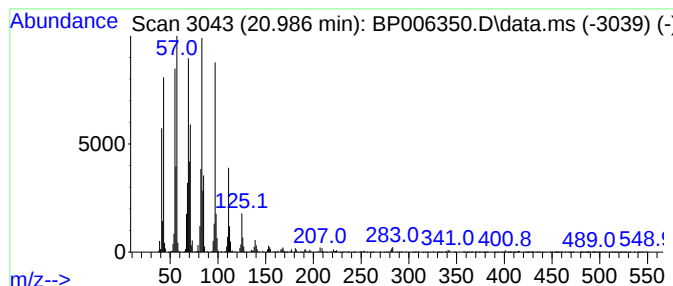
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	3.35 ng/ul	562198	Chrysene-d12	21.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacos-1-ene	350	C25H50	016980-85-1	94
2			1-Tetracosene	336	C24H48	010192-32-2	91
3			1-Hexacosanol	382	C26H54O	000506-52-5	91
4			Behenic alcohol	326	C22H46O	000661-19-8	91
5			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
 Data File : BP006350.D
 Acq On : 26 Jul 2021 22:00
 Operator : CG/JU
 Sample : M3091-20
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0058

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.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
Phenol, 2-methoxy-	8.880	6.0	ng/ul	458214	1	7.787	1525500	20.0
1,6-Octadien-3-...	9.040	2.3	ng/ul	174275	1	7.787	1525500	20.0
Dipropylene gly...	9.840	10.1	ng/ul	1085900	2	10.569	2151310	20.0
Pentane, 2-meth...	10.080	9.3	ng/ul	1001840	2	10.569	2151310	20.0
Ethanol, 1-(2-b...	10.360	6.3	ng/ul	676908	2	10.569	2151310	20.0
2-Propanol, 1-(...	11.120	2.3	ng/ul	246175	2	10.569	2151310	20.0
Methane, diethoxy-	11.180	2.5	ng/ul	272155	2	10.569	2151310	20.0
7,9-Di-tert-but...	17.850	13.5	ng/ul	2067920	4	17.163	3068980	20.0
n-Hexadecanoic ...	18.070	3.8	ng/ul	581650	4	17.163	3068980	20.0
Octadecanoic acid	19.300	2.3	ng/ul	393035	5	21.257	3355020	20.0
(DEL) Alkane: S...	20.990	3.4	ng/ul	562198	5	21.257	3355020	20.0