

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
 Data File : BP006326.D
 Acq On : 25 Jul 2021 03:03
 Operator : CG/JU
 Sample : M3063-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AW9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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: BP006326.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	21	24	28	rVB	12715	17777	0.22%	0.021%
2	3.305	33	37	42	rBV	62686	85620	1.06%	0.102%
3	3.711	102	106	123	rBV	256701	455209	5.64%	0.542%
4	3.928	139	143	151	rVB2	25758	45633	0.57%	0.054%
5	4.028	155	160	169	rVB	20166	33011	0.41%	0.039%
6	4.387	216	221	227	rVV	88030	121383	1.50%	0.144%
7	4.852	296	300	304	rBV4	10690	18182	0.23%	0.022%
8	5.664	434	438	443	rVB3	14177	21257	0.26%	0.025%
9	5.869	468	473	474	rBV5	15370	22418	0.28%	0.027%
10	5.893	474	477	485	rVB	86756	127712	1.58%	0.152%
11	6.787	619	629	634	rBV6	19584	48872	0.61%	0.058%
12	6.905	644	649	651	rBV	46873	67753	0.84%	0.081%
13	6.958	651	658	670	rVB2	390923	847249	10.50%	1.008%
14	7.134	682	688	698	rVB	1178454	1829119	22.67%	2.176%
15	7.322	713	720	731	rVB	1227123	1881993	23.33%	2.239%
16	7.428	734	738	744	rVV7	17132	31505	0.39%	0.037%
17	7.487	744	748	754	rVB8	13754	19801	0.25%	0.024%
18	7.787	791	799	806	rBV	1061041	1684686	20.88%	2.004%
19	7.858	806	811	818	rVB	58651	90892	1.13%	0.108%
20	8.028	834	840	845	rBV4	17723	43197	0.54%	0.051%
21	8.293	878	885	889	rBV2	20547	36314	0.45%	0.043%
22	8.405	899	904	911	rVB3	19187	36530	0.45%	0.043%
23	8.493	911	919	931	rBV	1019625	1695909	21.02%	2.017%
24	8.610	931	939	945	rVB	59449	98466	1.22%	0.117%
25	8.881	979	985	990	rBV	234148	381156	4.72%	0.453%
26	8.946	990	996	1002	rVB	1362446	2193041	27.18%	2.609%
27	9.116	1020	1025	1032	rVB3	52318	81217	1.01%	0.097%
28	9.505	1084	1091	1097	rVB10	15348	35321	0.44%	0.042%
29	9.657	1109	1117	1133	rBV	1061720	1788626	22.17%	2.128%
30	9.793	1136	1140	1144	rVV6	7373	15016	0.19%	0.018%
31	9.846	1144	1149	1152	rVV5	18840	32959	0.41%	0.039%
32	9.893	1152	1157	1167	rVB3	64466	117582	1.46%	0.140%
33	9.999	1170	1175	1182	rBV10	9348	18710	0.23%	0.022%
34	10.110	1186	1194	1201	rVV9	9470	28093	0.35%	0.033%
35	10.193	1201	1208	1217	rVV	1640257	2593657	32.15%	3.085%
36	10.363	1230	1237	1246	rBV	522841	801072	9.93%	0.953%

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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

37	10.487	1252	1258	1265	rVV4	82446	150112	1.86%	0.179%
38	10.575	1265	1273	1279	rVV	1468114	2418352	29.98%	2.877%
39	10.640	1279	1284	1289	rVV2	67579	117012	1.45%	0.139%
40	10.716	1289	1297	1307	rVB	1498933	2586277	32.06%	3.077%
41	10.887	1318	1326	1332	rBV	238577	377901	4.68%	0.450%
42	10.969	1336	1340	1349	rVB3	27698	51676	0.64%	0.061%
43	11.122	1361	1366	1372	rBV2	15294	23240	0.29%	0.028%
44	11.222	1379	1383	1389	rBV2	21016	38408	0.48%	0.046%
45	11.369	1403	1408	1417	rVB2	57203	100443	1.25%	0.119%
46	11.475	1422	1426	1431	rVV3	9037	14946	0.19%	0.018%
47	11.557	1431	1440	1447	rVV9	14568	43425	0.54%	0.052%
48	11.751	1469	1473	1478	rBV4	17057	28780	0.36%	0.034%
49	11.857	1487	1491	1496	rVB2	31580	44955	0.56%	0.053%
50	12.163	1536	1543	1549	rVB	34699	60857	0.75%	0.072%
51	12.334	1562	1572	1576	rBV8	13899	38415	0.48%	0.046%
52	12.616	1615	1620	1624	rBV5	13475	20756	0.26%	0.025%
53	12.698	1626	1634	1635	rBV8	13778	23157	0.29%	0.028%
54	12.781	1643	1648	1654	rVB	54514	83080	1.03%	0.099%
55	12.881	1660	1665	1669	rBV	22705	34573	0.43%	0.041%
56	13.034	1685	1691	1696	rBV	116516	161129	2.00%	0.192%
57	13.263	1726	1730	1735	rVB	11090	15175	0.19%	0.018%
58	13.340	1735	1743	1749	rBV	400371	595192	7.38%	0.708%
59	13.651	1790	1796	1802	rBV2	26888	50951	0.63%	0.061%
60	13.787	1815	1819	1821	rBV2	19730	25072	0.31%	0.030%
61	13.834	1821	1827	1835	rBV	3127060	4323166	53.59%	5.143%
62	14.104	1863	1873	1880	rBV	3546057	5034642	62.41%	5.989%
63	14.275	1898	1902	1907	rBV	13386	19772	0.25%	0.024%
64	14.416	1917	1926	1933	rBV	2116934	3117130	38.64%	3.708%
65	14.504	1936	1941	1945	rVB2	32059	43637	0.54%	0.052%
66	14.610	1949	1959	1969	rBV	369792	565735	7.01%	0.673%
67	14.840	1994	1998	2002	rVB5	26556	31179	0.39%	0.037%
68	15.016	2021	2028	2034	rBV2	36900	56834	0.70%	0.068%
69	15.098	2034	2042	2047	rVV	61465	102257	1.27%	0.122%
70	15.145	2047	2050	2052	rVV2	14487	20581	0.26%	0.024%
71	15.169	2052	2054	2059	rVB	15665	18643	0.23%	0.022%
72	15.245	2060	2067	2072	rBV2	105227	173926	2.16%	0.207%
73	15.410	2087	2095	2106	rBV	4447194	6326232	78.42%	7.525%
74	15.522	2108	2114	2128	rVV	1408515	2000021	24.79%	2.379%
75	15.651	2131	2136	2142	rVB2	55717	95698	1.19%	0.114%
76	15.781	2153	2158	2166	rVB2	48956	74407	0.92%	0.089%

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 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	15.863	2167	2172	2176	rVB5	16609	23575	0.29%	0.028%
78	16.434	2265	2269	2273	rBV	11165	16743	0.21%	0.020%
79	16.739	2314	2321	2323	rBV4	11153	22719	0.28%	0.027%
80	16.769	2323	2326	2330	rVB3	13993	18791	0.23%	0.022%
81	16.945	2351	2356	2364	rBV2	17619	29657	0.37%	0.035%
82	17.163	2385	2393	2403	rVB	2442057	3517979	43.61%	4.185%
83	17.263	2403	2410	2423	rBV2	5078095	7098529	87.99%	8.444%
84	17.351	2423	2425	2432	rVB7	8102	15466	0.19%	0.018%
85	17.469	2442	2445	2450	rVB	17799	22438	0.28%	0.027%
86	17.539	2453	2457	2466	rBV5	10900	25733	0.32%	0.031%
87	17.845	2503	2509	2518	rBV2	1115624	1470689	18.23%	1.749%
88	18.069	2542	2547	2554	rBV	336942	472872	5.86%	0.563%
89	18.133	2554	2558	2571	rVB3	92039	211968	2.63%	0.252%
90	18.310	2585	2588	2592	rVB4	31520	38641	0.48%	0.046%
91	18.628	2639	2642	2646	rVB2	22496	23452	0.29%	0.028%
92	19.080	2715	2719	2724	rBV2	67152	93699	1.16%	0.111%
93	19.151	2727	2731	2735	rBV	61903	95159	1.18%	0.113%
94	19.310	2754	2758	2765	rBV	186941	250159	3.10%	0.298%
95	19.504	2785	2791	2801	rBV	5973708	7969995	98.79%	9.481%
96	19.733	2828	2830	2835	rBV2	17330	23337	0.29%	0.028%
97	20.986	3040	3043	3050	rBV2	326459	411433	5.10%	0.489%
98	21.263	3085	3090	3098	rBV	2967714	3752579	46.52%	4.464%
99	23.457	3456	3463	3471	rVB2	4304872	8067457	100.00%	9.597%
100	23.604	3482	3488	3498	rVB2	1961938	3836470	47.55%	4.564%

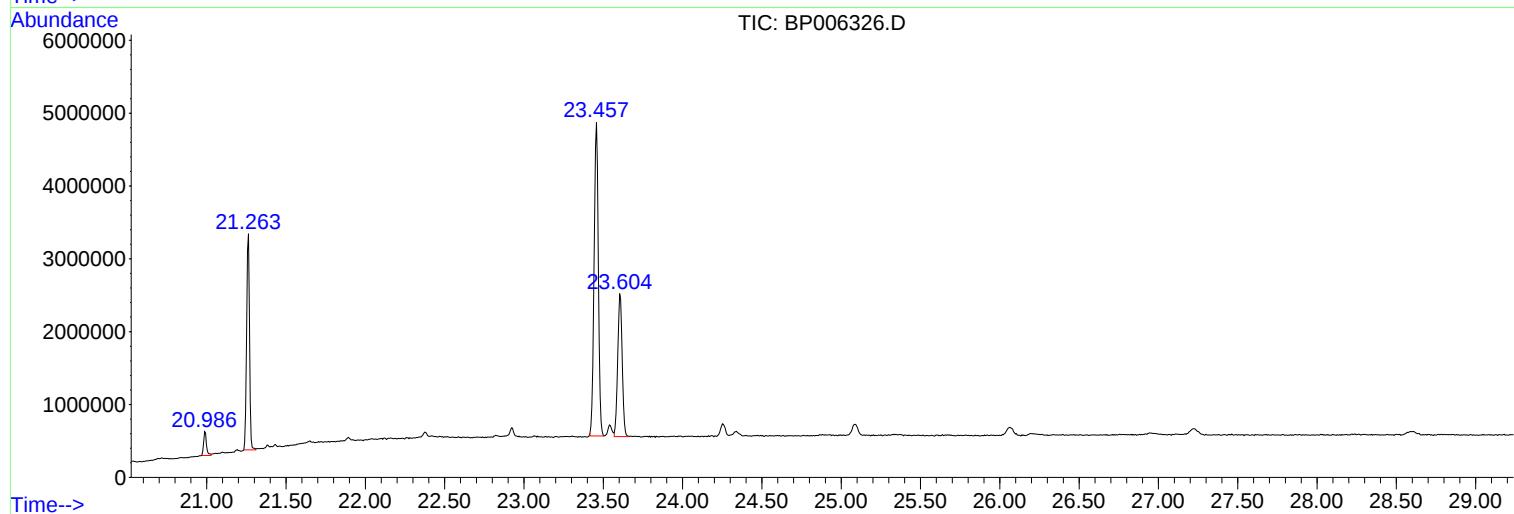
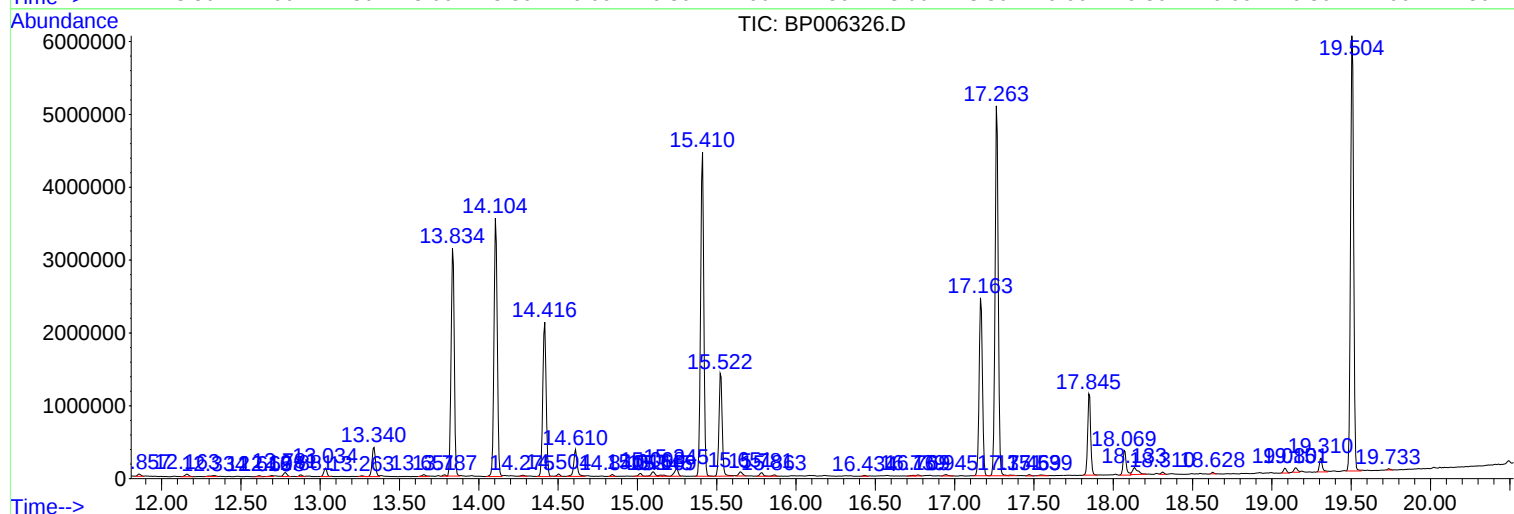
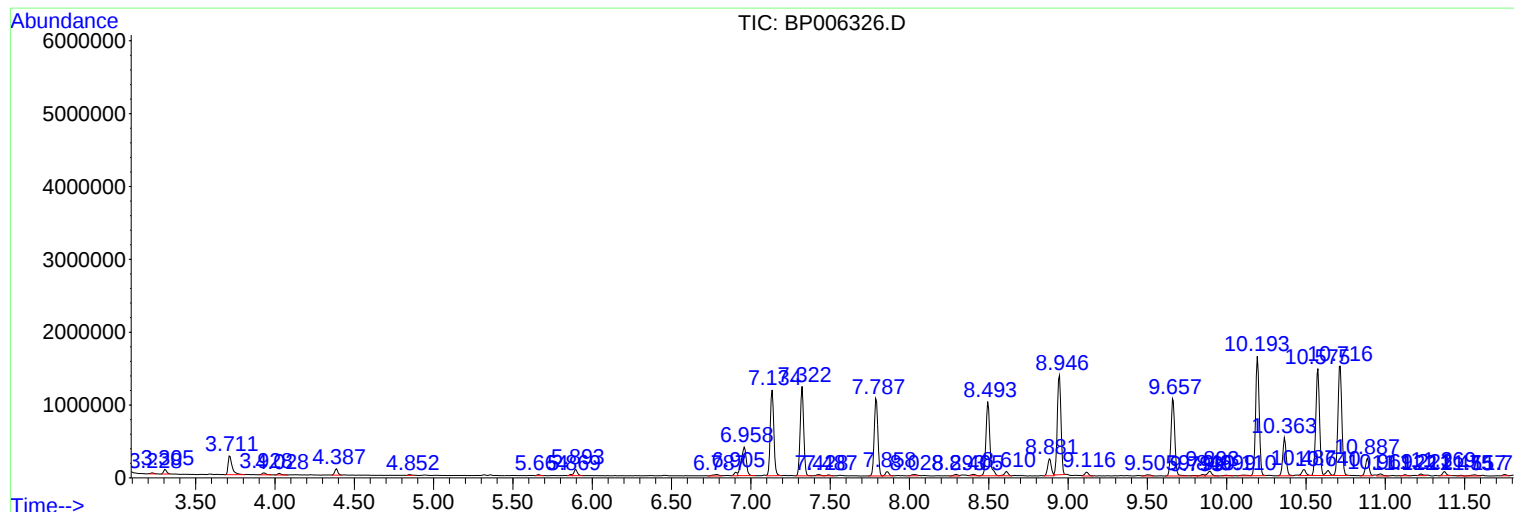
Sum of corrected areas: 84064222

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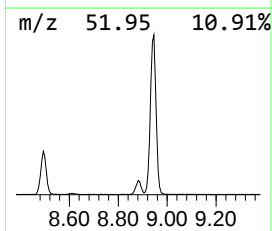
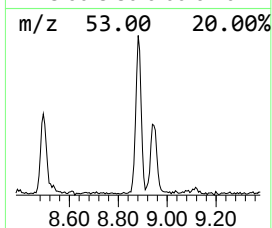
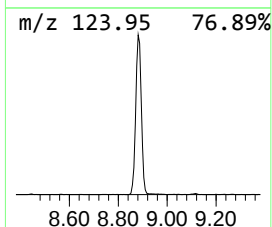
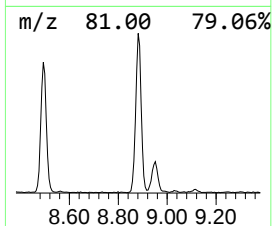
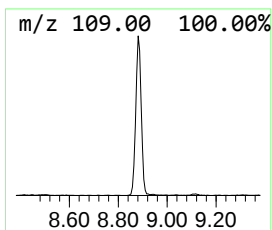
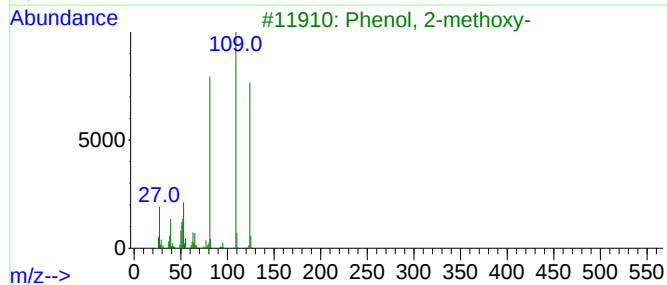
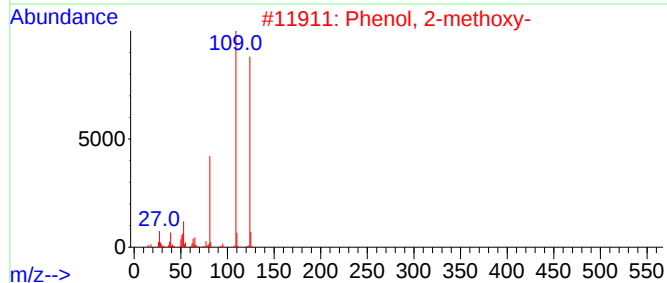
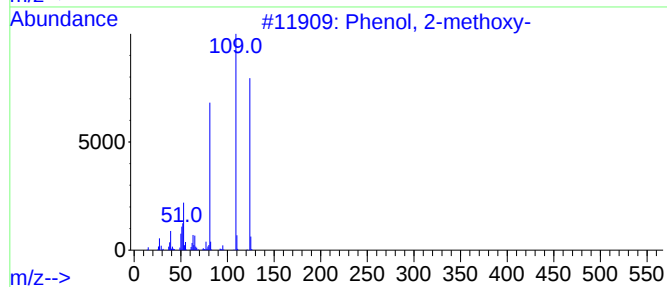
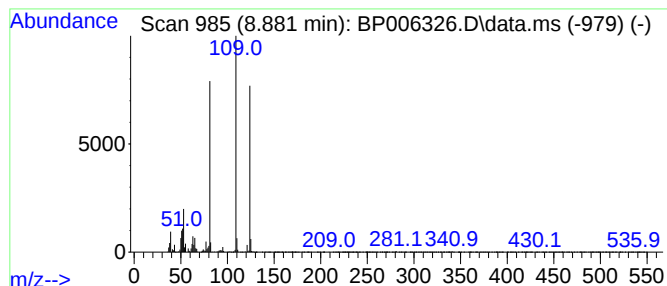
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 2-methoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.880	4.52 ng/ul	381156	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methoxy-		124	C7H8O2		000090-05-1	96
2	Phenol, 2-methoxy-		124	C7H8O2		000090-05-1	95
3	Phenol, 2-methoxy-		124	C7H8O2		000090-05-1	91
4	Mequinol		124	C7H8O2		000150-76-5	91
5	Phenol, 2-methoxy-		124	C7H8O2		000090-05-1	90



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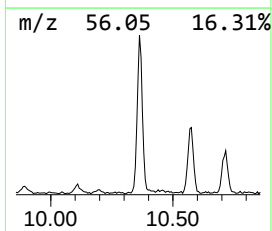
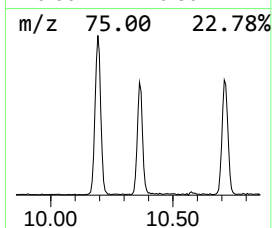
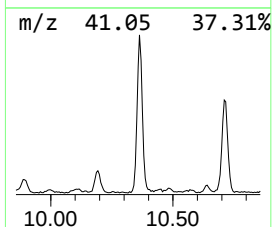
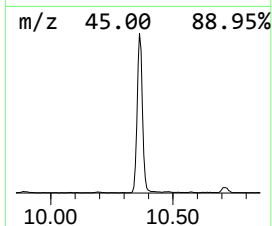
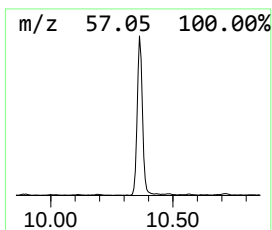
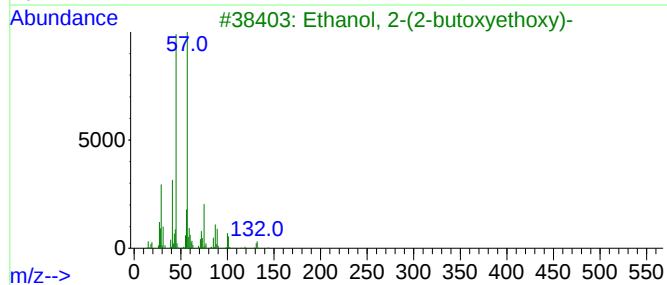
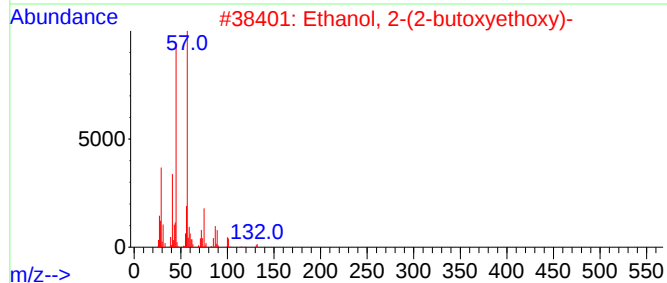
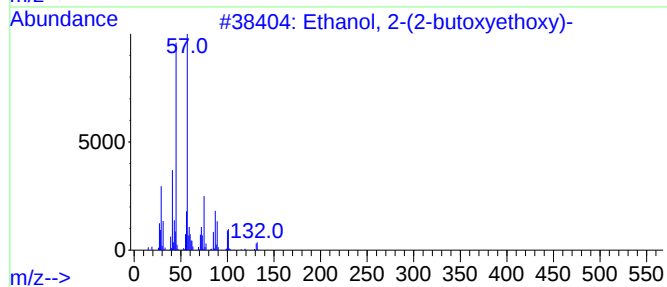
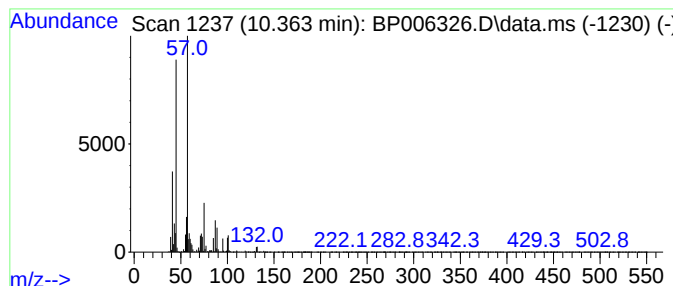
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	6.62 ng/ul	801072	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	64



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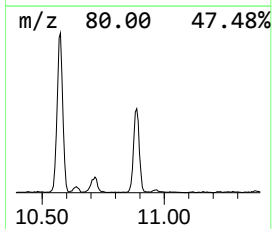
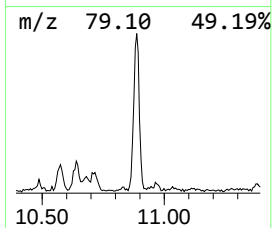
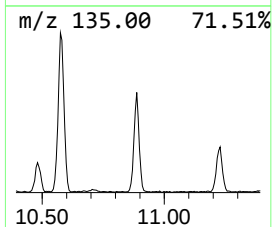
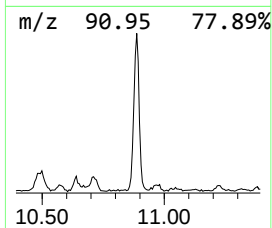
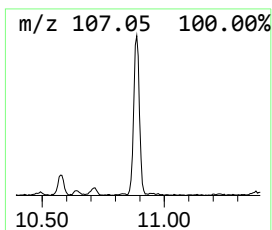
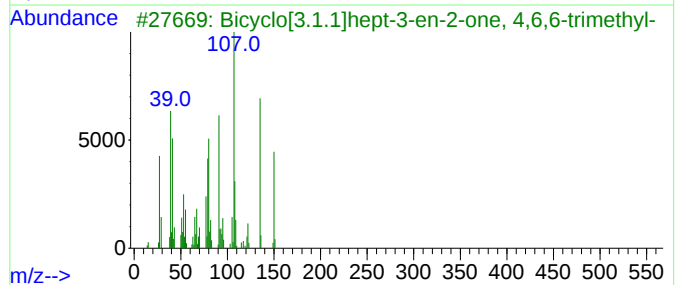
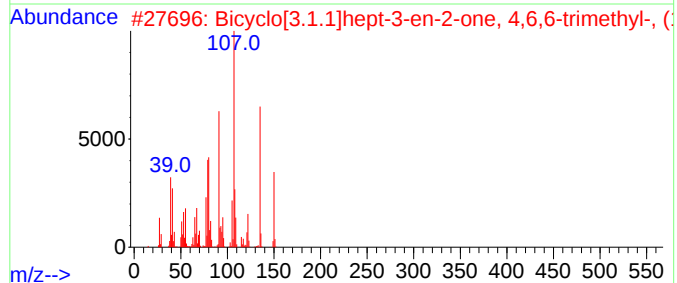
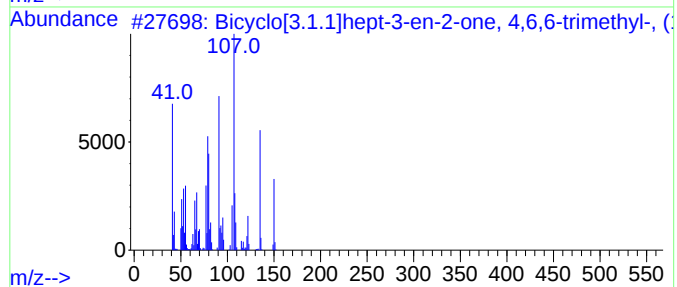
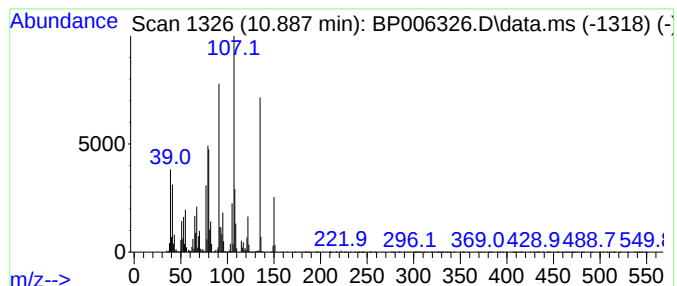
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Bicyclo[3.1.1]hept-3-en-2-o... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.890	3.13 ng/ul	377901	Naphthalene-d8	10.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	001196-01-6	96
2			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	001196-01-6	96
3			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	000080-57-9	94
4			Bicyclo[3.1.1]hept-3-en-2-one, 4...	150	C10H14O	001196-01-6	90
5			Bicyclo[3.1.1]hept-2-en-6-one, 2...	150	C10H14O	000473-06-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006326.D
Acq On : 25 Jul 2021 03:03
Operator : CG/JU
Sample : M3063-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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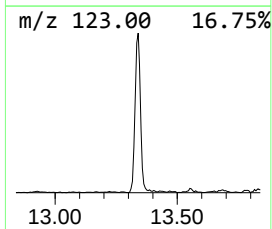
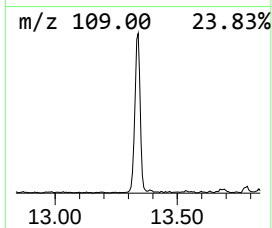
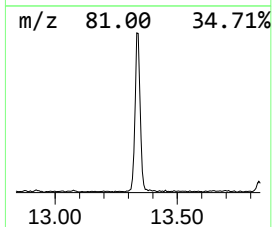
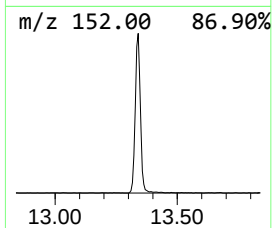
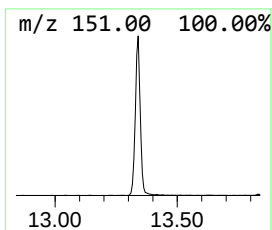
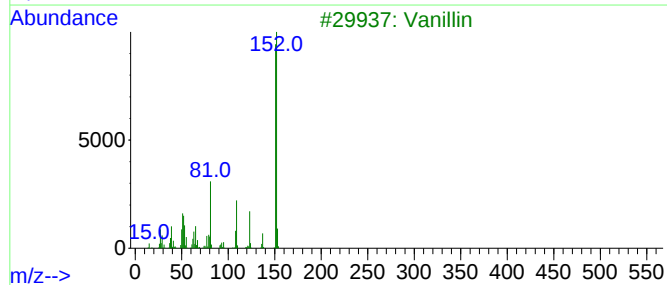
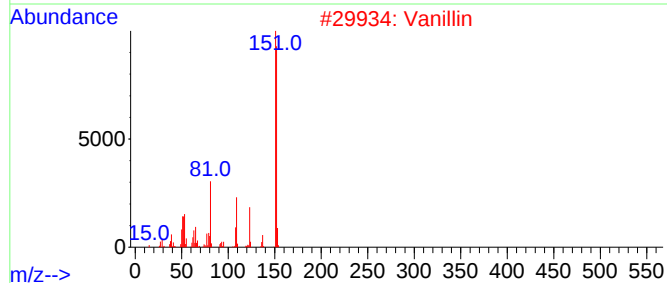
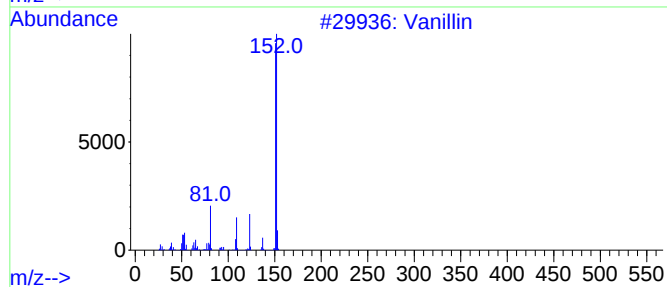
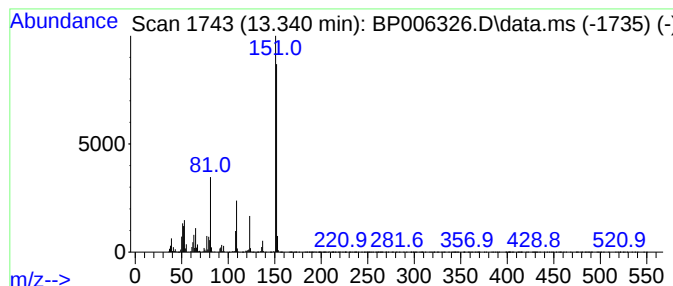
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Vanillin Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.340	3.82 ng/ul	595192	Acenaphthene-d10	14.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin			152	C8H8O3	000121-33-5	97
2	Vanillin			152	C8H8O3	000121-33-5	97
3	Vanillin			152	C8H8O3	000121-33-5	97
4	Benzaldehyde, 3-hydroxy-4-methoxy-			152	C8H8O3	000621-59-0	97
5	Benzaldehyde, 3-hydroxy-4-methoxy-			152	C8H8O3	000621-59-0	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006326.D
Acq On : 25 Jul 2021 03:03
Operator : CG/JU
Sample : M3063-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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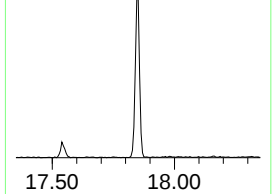
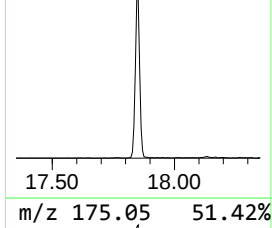
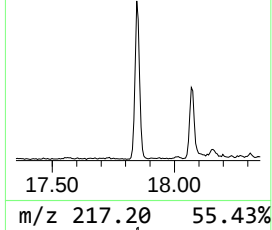
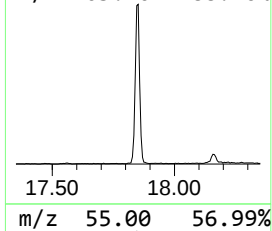
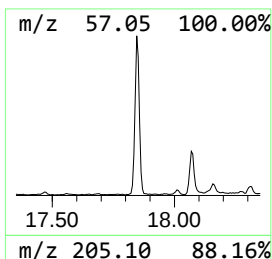
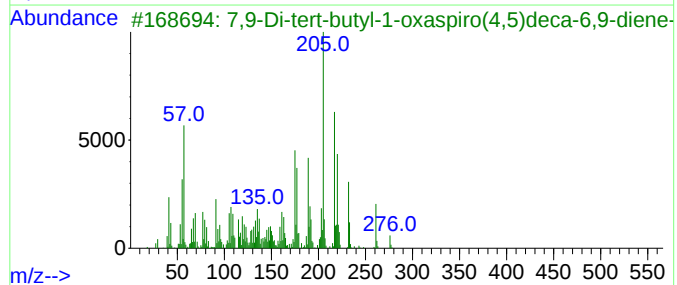
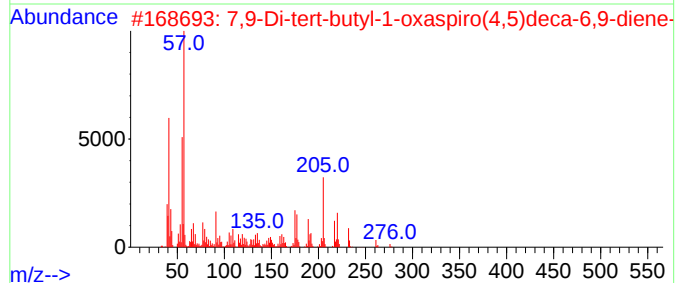
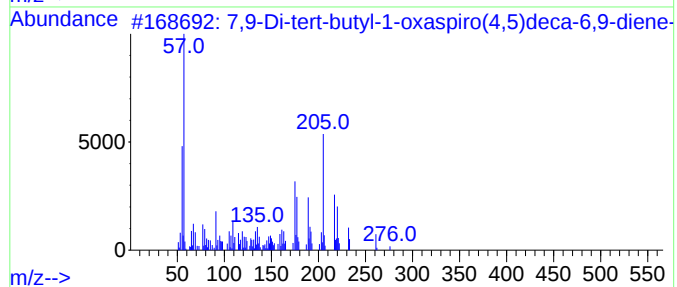
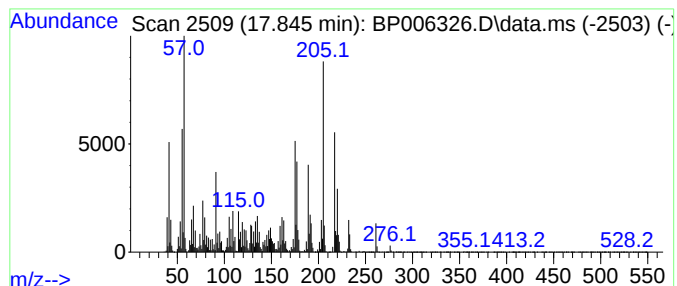
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	8.36 ng/ul	1470690	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	97		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95		
4	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	83		
5	1,3-Pentadiene, 1,1-diphenyl-, (Z)-	220	C17H16	015295-31-5	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006326.D
Acq On : 25 Jul 2021 03:03
Operator : CG/JU
Sample : M3063-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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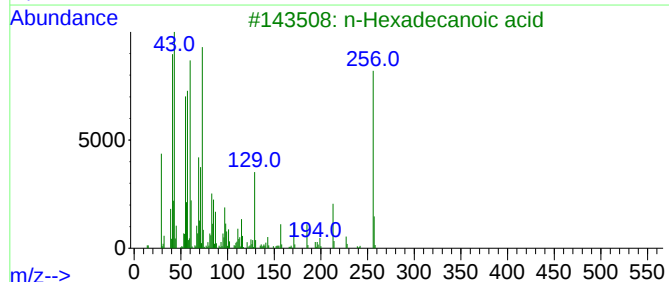
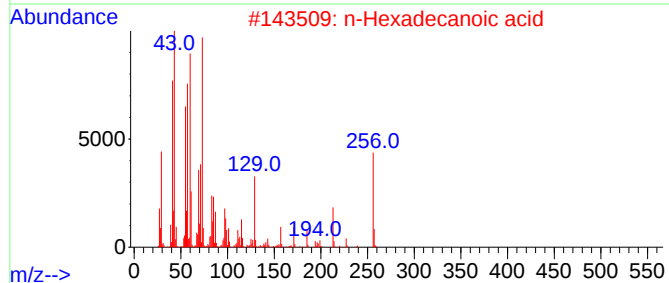
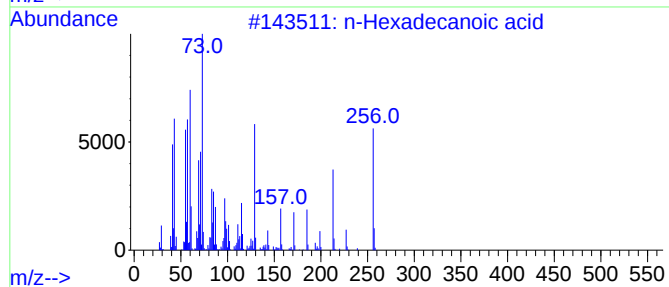
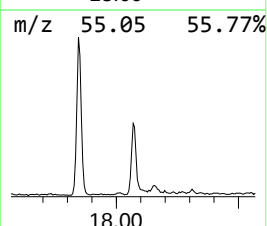
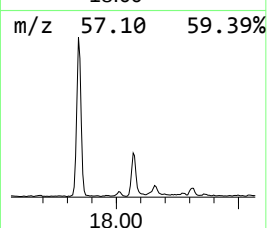
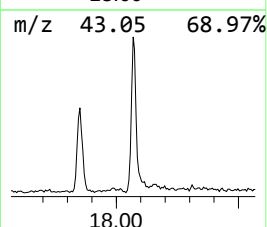
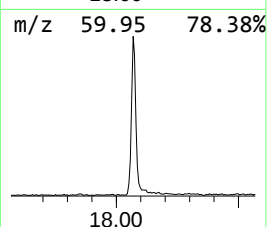
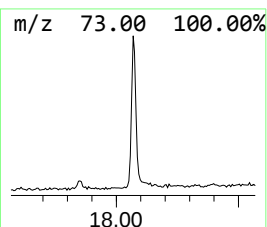
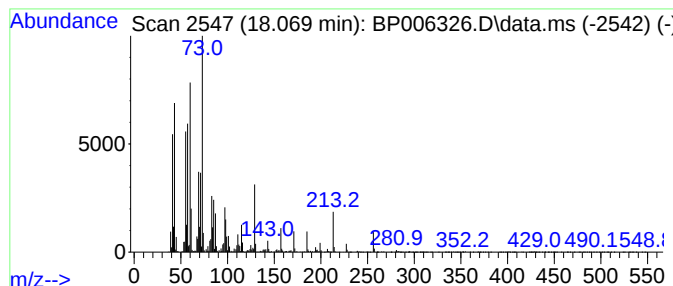
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	2.69 ng/ul	472872	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006326.D
Acq On : 25 Jul 2021 03:03
Operator : CG/JU
Sample : M3063-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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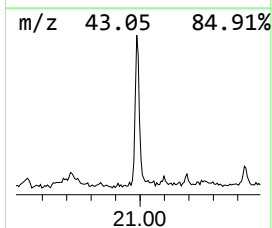
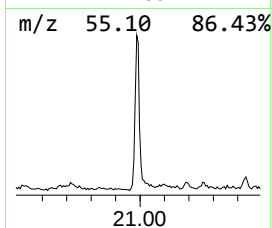
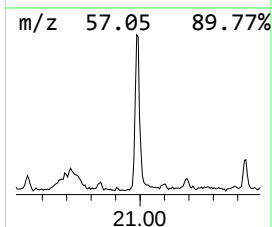
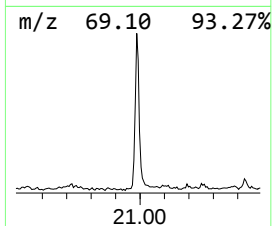
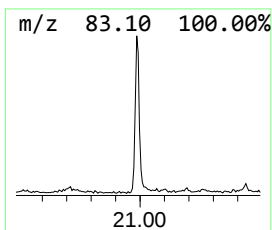
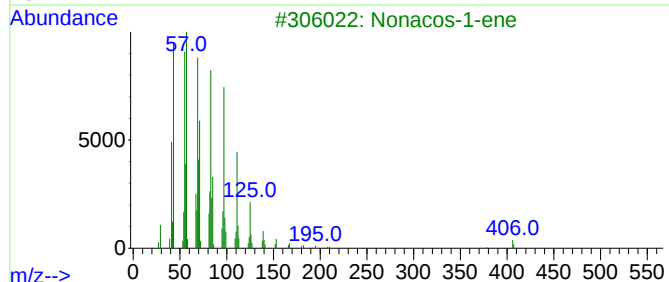
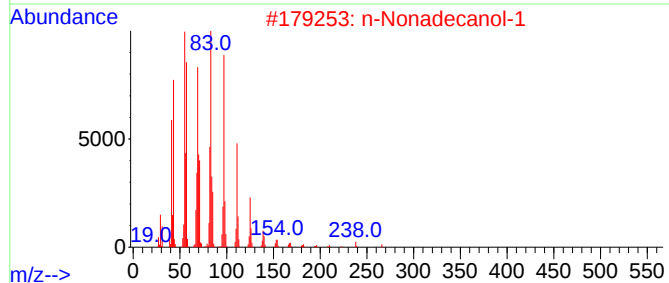
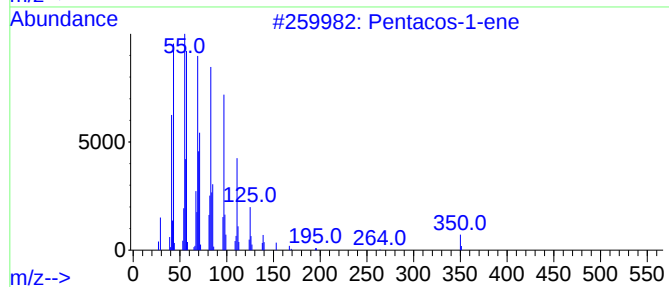
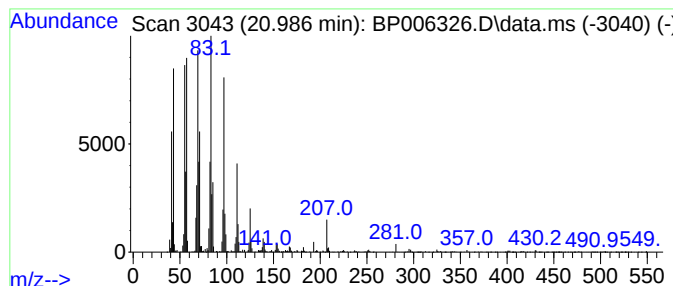
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	2.19 ng/ul	411433	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacos-1-ene	350	C25H50	016980-85-1	95
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
3			Nonacos-1-ene	406	C29H58	018835-35-3	94
4			1-Heneicosanol	312	C21H44O	015594-90-8	93
5			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072421\
Data File : BP006326.D
Acq On : 25 Jul 2021 03:03
Operator : CG/JU
Sample : M3063-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AW9

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	4.5	ng/ul	381156	1	7.787	1684690	20.0
Ethanol, 2-(2-b...	10.360	6.6	ng/ul	801072	2	10.575	2418350	20.0
Bicyclo[3.1.1]h...	10.890	3.1	ng/ul	377901	2	10.575	2418350	20.0
Vanillin	13.340	3.8	ng/ul	595192	3	14.416	3117130	20.0
7,9-Di-tert-but...	17.850	8.4	ng/ul	1470690	4	17.163	3517980	20.0
n-Hexadecanoic ...	18.070	2.7	ng/ul	472872	4	17.163	3517980	20.0
(DEL) Alkane: S...	20.990	2.2	ng/ul	411433	5	21.263	3752580	20.0