

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007507.D
 Acq On : 20 Oct 2021 00:59
 Operator : CG/JU
 Sample : M4207-09
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B22

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-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007507.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.652	58	62	68	rVB4	28399	46243	1.41%	0.124%
2	3.805	86	88	99	rBV	64977	98071	2.99%	0.263%
3	4.040	123	128	137	rVV	58688	89941	2.74%	0.242%
4	4.134	139	144	151	rVB	46065	71498	2.18%	0.192%
5	5.593	387	392	399	rVV	58356	97736	2.98%	0.262%
6	5.969	450	456	464	rVV3	90300	186815	5.69%	0.502%
7	6.334	514	518	527	rVB	23001	41757	1.27%	0.112%
8	6.452	532	538	545	rVV5	34998	81739	2.49%	0.219%
9	6.599	556	563	566	rVV5	28497	59060	1.80%	0.159%
10	6.934	614	620	629	rBV3	28580	75010	2.28%	0.201%
11	7.111	643	650	659	rVB	212236	357670	10.89%	0.960%
12	7.281	669	679	685	rBV	197961	330644	10.07%	0.888%
13	7.346	685	690	694	rVB2	27520	41657	1.27%	0.112%
14	7.481	700	713	720	rBV	234537	380766	11.59%	1.022%
15	7.599	727	733	740	rVB2	90046	169775	5.17%	0.456%
16	7.952	786	793	799	rVV	706306	1106486	33.68%	2.971%
17	8.075	809	814	822	rVV2	28713	51429	1.57%	0.138%
18	8.658	907	913	923	rVV	193909	321176	9.78%	0.862%
19	8.769	923	932	941	rVV2	31732	79373	2.42%	0.213%
20	9.105	974	989	996	rBV	217785	429285	13.07%	1.153%
21	9.205	1001	1006	1015	rVB	76755	125141	3.81%	0.336%
22	9.834	1105	1113	1121	rBV	156659	273203	8.32%	0.734%
23	10.375	1197	1205	1213	rVV	280934	489734	14.91%	1.315%
24	10.540	1227	1233	1241	rBV	145911	239934	7.30%	0.644%
25	10.757	1262	1270	1275	rVV	968047	1652868	50.32%	4.438%
26	10.810	1275	1279	1285	rVV	147860	243043	7.40%	0.653%
27	10.887	1285	1292	1298	rVV	233983	396745	12.08%	1.065%
28	10.946	1298	1302	1308	rVB	31992	47434	1.44%	0.127%
29	11.810	1442	1449	1455	rBV	59530	101724	3.10%	0.273%
30	12.199	1509	1515	1523	rBV2	94929	148622	4.52%	0.399%
31	12.416	1544	1552	1561	rVB	238116	387325	11.79%	1.040%
32	12.634	1583	1589	1595	rVV	189925	298898	9.10%	0.803%
33	13.151	1672	1677	1681	rVB	46347	64485	1.96%	0.173%
34	13.434	1716	1725	1730	rBV2	137318	222144	6.76%	0.596%
35	13.498	1730	1736	1742	rVB	36412	57316	1.74%	0.154%
36	13.622	1750	1757	1761	rBV3	26689	52858	1.61%	0.142%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
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37	13.757	1774	1780	1781	rBV	51778	80501	2.45%	0.216%
38	13.910	1802	1806	1811	rVV2	125412	210856	6.42%	0.566%
39	13.963	1811	1815	1817	rVV	113302	163032	4.96%	0.438%
40	13.998	1817	1821	1830	rVV	530563	760239	23.14%	2.041%
41	14.093	1832	1837	1841	rVV3	81202	106885	3.25%	0.287%
42	14.145	1841	1846	1850	rVV2	82819	140931	4.29%	0.378%
43	14.281	1863	1869	1879	rVB	629612	986679	30.04%	2.649%
44	14.504	1902	1907	1912	rBV	124987	163443	4.98%	0.439%
45	14.587	1912	1921	1928	rBV	1416500	2189879	66.66%	5.880%
46	14.769	1946	1952	1963	rBV2	132359	252888	7.70%	0.679%
47	14.987	1984	1989	1997	rVB2	117846	216953	6.60%	0.583%
48	15.063	1997	2002	2007	rVB2	41550	68361	2.08%	0.184%
49	15.128	2009	2013	2019	rBV4	36499	57385	1.75%	0.154%
50	15.210	2021	2027	2032	rBV3	40280	75165	2.29%	0.202%
51	15.263	2032	2036	2040	rVB2	42620	56062	1.71%	0.151%
52	15.334	2040	2048	2051	rBV4	27037	46624	1.42%	0.125%
53	15.381	2051	2056	2060	rBV3	53214	99591	3.03%	0.267%
54	15.457	2065	2069	2074	rBV	134272	155695	4.74%	0.418%
55	15.581	2084	2090	2095	rVV	826509	1174802	35.76%	3.154%
56	15.628	2095	2098	2103	rVB	118498	160419	4.88%	0.431%
57	15.692	2103	2109	2114	rBV2	113927	174917	5.32%	0.470%
58	15.863	2133	2138	2143	rVV3	73085	145082	4.42%	0.390%
59	15.934	2146	2150	2159	rVV	84244	153515	4.67%	0.412%
60	16.016	2159	2164	2168	rVV5	30192	51154	1.56%	0.137%
61	16.081	2168	2175	2182	rVV2	109843	237647	7.23%	0.638%
62	16.163	2183	2189	2199	rVB3	43123	92709	2.82%	0.249%
63	16.322	2211	2216	2218	rVV	181822	245761	7.48%	0.660%
64	16.351	2218	2221	2226	rVB	241288	328726	10.01%	0.883%
65	16.881	2306	2311	2317	rBV3	69181	153144	4.66%	0.411%
66	16.992	2326	2330	2338	rVB4	43751	78838	2.40%	0.212%
67	17.075	2338	2344	2348	rVV2	51227	102969	3.13%	0.276%
68	17.122	2348	2352	2356	rVV	184422	248996	7.58%	0.669%
69	17.175	2357	2361	2368	rVB3	56286	112661	3.43%	0.303%
70	17.339	2383	2389	2394	rBV	1741262	2624166	79.88%	7.046%
71	17.381	2394	2396	2401	rVV	280641	354384	10.79%	0.952%
72	17.439	2401	2406	2417	rVB	944352	1389364	42.29%	3.731%
73	17.534	2418	2422	2427	rVB7	39371	60843	1.85%	0.163%
74	17.657	2439	2443	2448	rVB4	53452	66944	2.04%	0.180%
75	17.857	2469	2477	2484	rVV	207315	281063	8.56%	0.755%
76	18.022	2501	2505	2510	rBV2	62022	77894	2.37%	0.209%

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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

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 Peak separation: 5

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77	18.210	2533	2537	2538	rBV	88677	102853	3.13%	0.276%
78	18.251	2538	2544	2549	rVB2	201035	428720	13.05%	1.151%
79	18.386	2562	2567	2571	rBV2	125447	198928	6.06%	0.534%
80	18.428	2571	2574	2580	rVB	111625	153712	4.68%	0.413%
81	18.528	2587	2591	2600	rBV	215872	264080	8.04%	0.709%
82	18.692	2615	2619	2622	rBV	67753	91037	2.77%	0.244%
83	19.092	2683	2687	2691	rBV	92615	143086	4.36%	0.384%
84	19.139	2691	2695	2700	rVV2	275964	431614	13.14%	1.159%
85	19.181	2700	2702	2706	rVB	74141	84243	2.56%	0.226%
86	19.228	2707	2710	2712	rBV2	46498	58062	1.77%	0.156%
87	19.345	2726	2730	2738	rVB	270869	356798	10.86%	0.958%
88	19.469	2748	2751	2753	rVB2	62562	62630	1.91%	0.168%
89	19.669	2780	2785	2789	rBV	1292370	1879128	57.20%	5.046%
90	19.698	2789	2790	2799	rVB	418180	359020	10.93%	0.964%
91	19.775	2800	2803	2808	rVB2	57582	88132	2.68%	0.237%
92	20.216	2875	2878	2881	rBV	216622	214413	6.53%	0.576%
93	20.698	2957	2960	2965	rBV	204303	207008	6.30%	0.556%
94	21.151	3033	3037	3041	rBV2	295463	402532	12.25%	1.081%
95	21.427	3078	3084	3088	rBV	2338936	3284993	100.00%	8.821%
96	21.598	3110	3113	3119	rVB	188802	218012	6.64%	0.585%
97	22.069	3190	3193	3200	rVB	233114	321917	9.80%	0.864%
98	22.586	3278	3281	3287	rVB2	161790	225131	6.85%	0.605%
99	23.716	3468	3473	3479	rVB2	806169	1422705	43.31%	3.820%
100	23.874	3494	3500	3509	rVB2	1520274	3178934	96.77%	8.536%

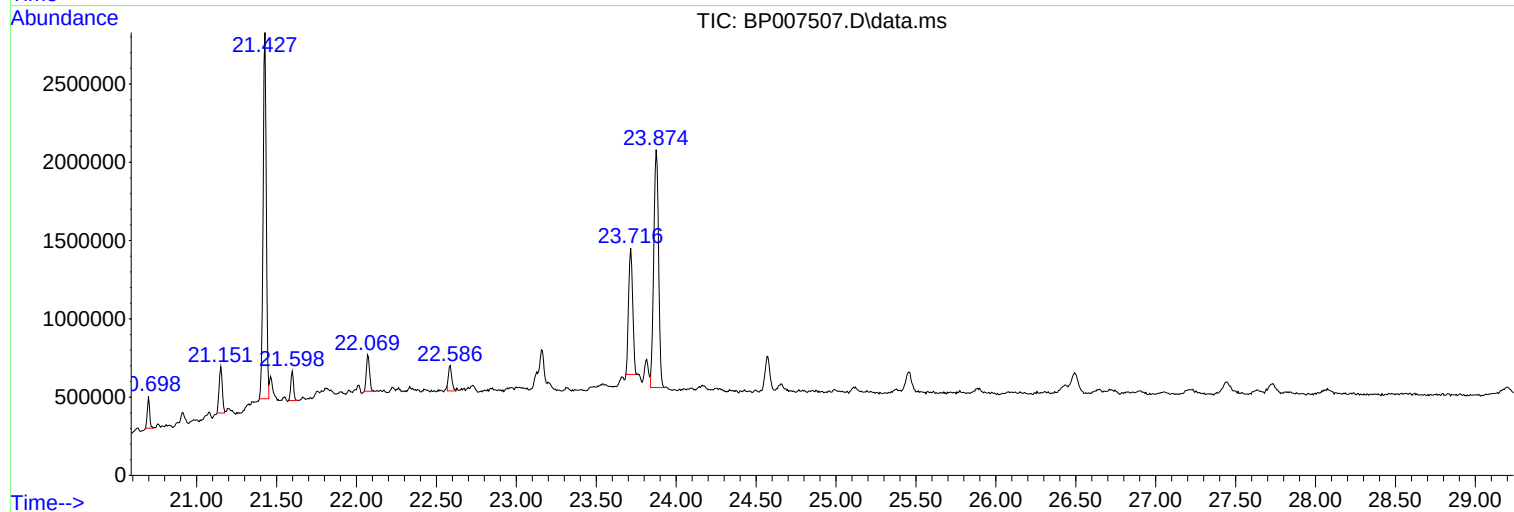
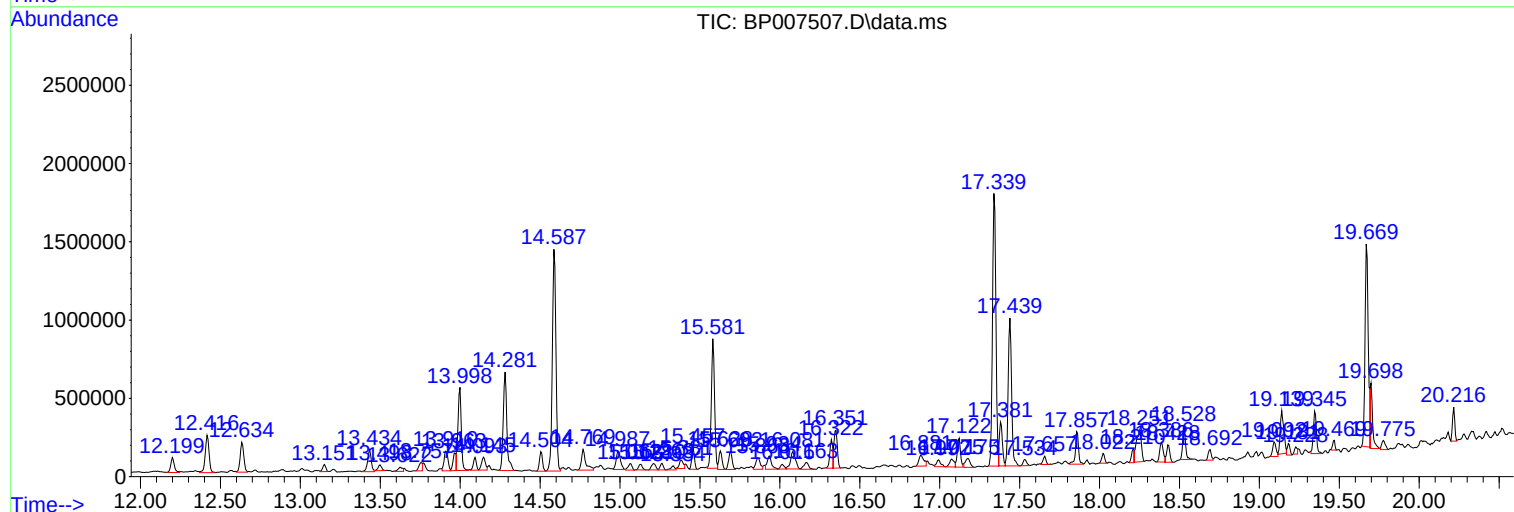
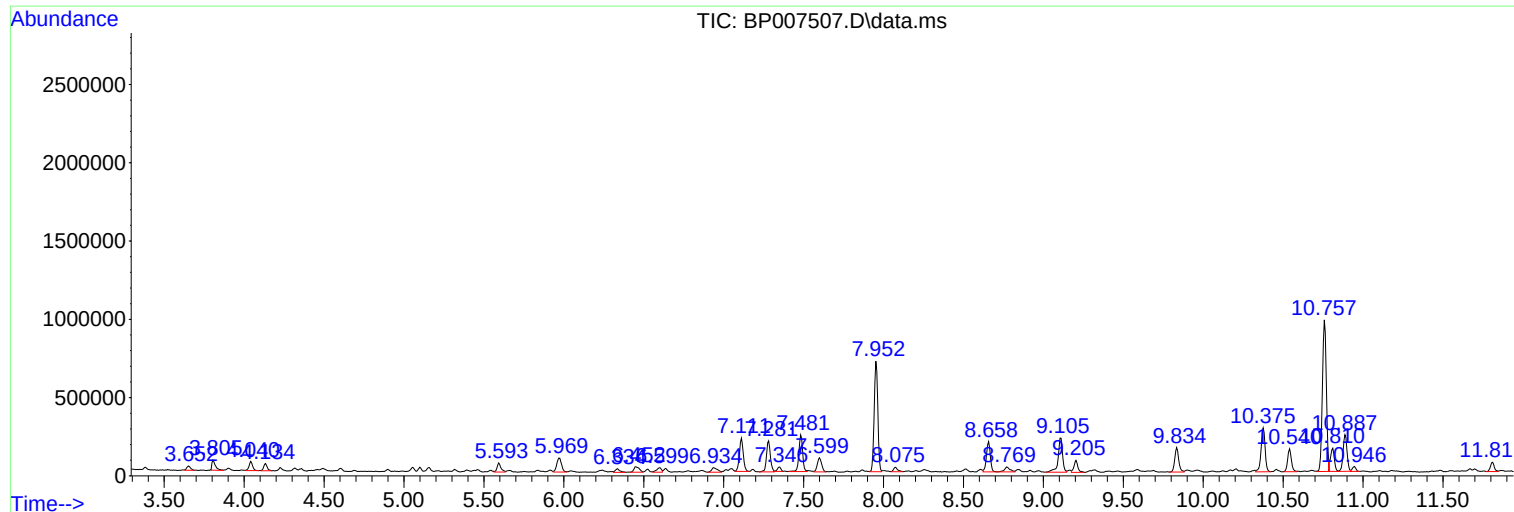
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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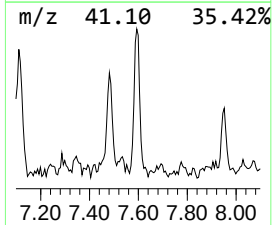
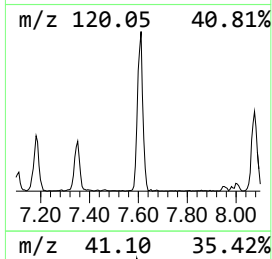
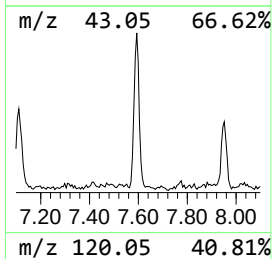
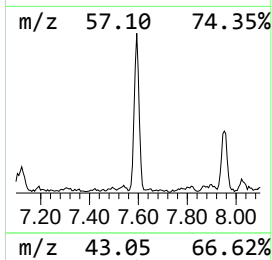
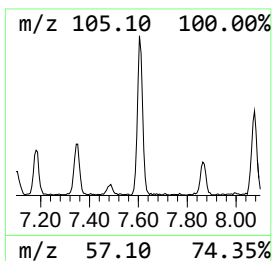
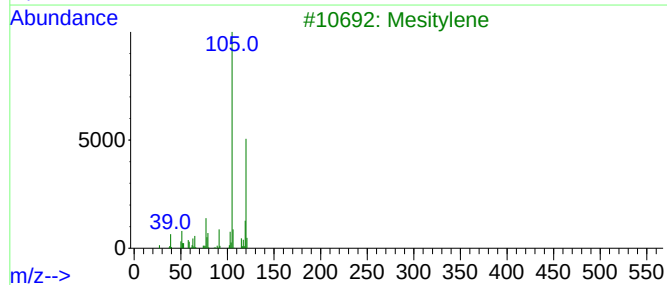
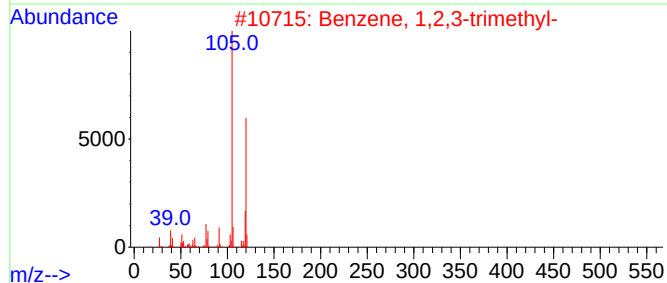
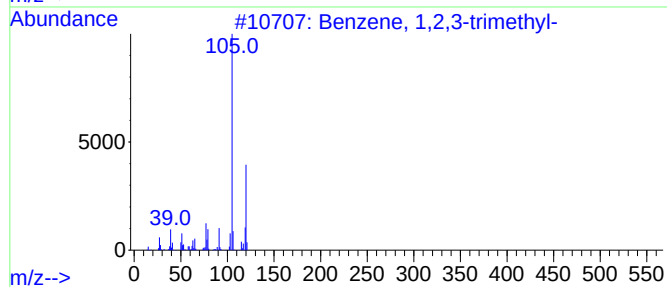
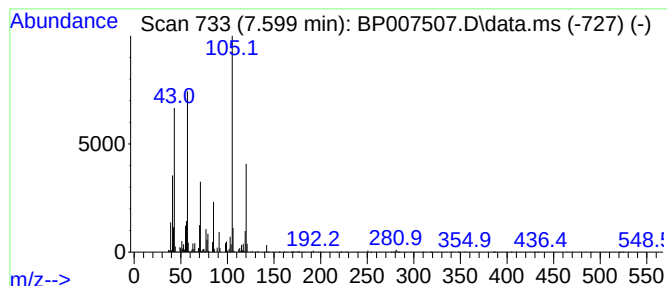
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.599	3.07 ng/ul	169775	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	90
4			Mesitylene	120	C9H12	000108-67-8	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	60



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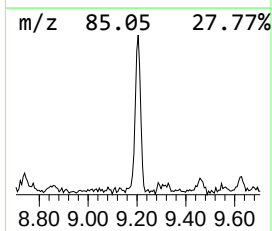
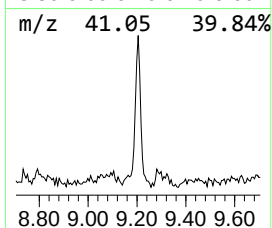
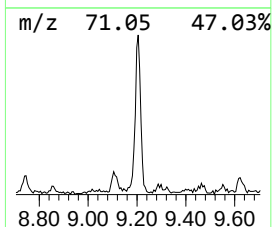
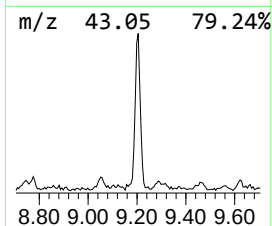
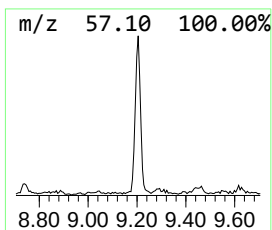
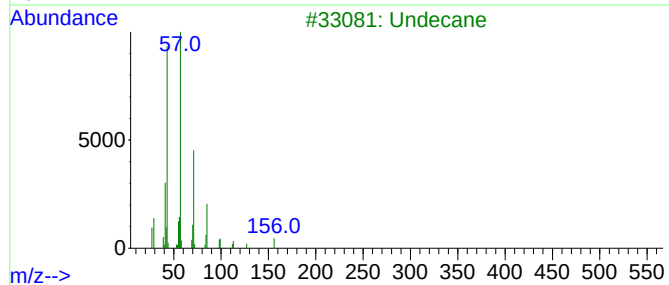
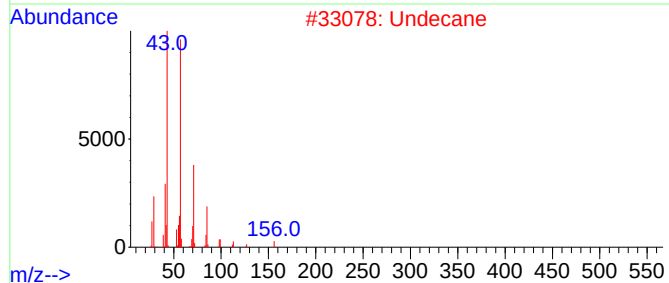
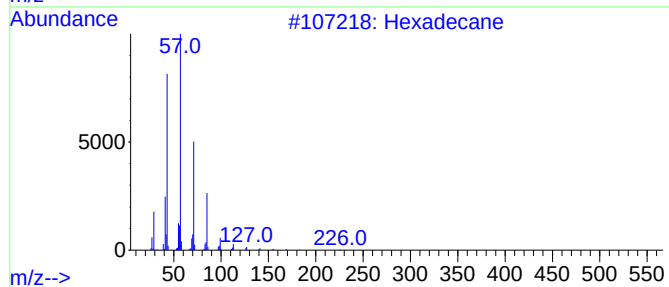
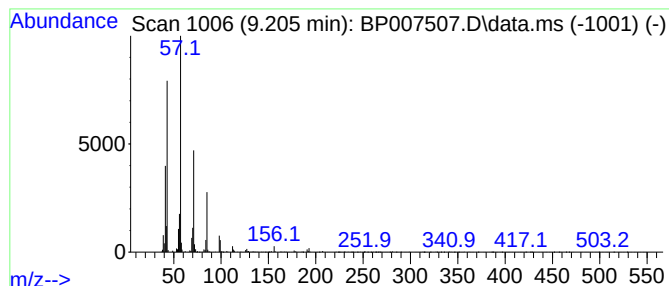
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.205	2.26 ng/ul	125141	1,4-Dichlorobenzene-d4	7.952

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	94
2		Undecane	156	C11H24	001120-21-4	90
3		Undecane	156	C11H24	001120-21-4	87
4		Tridecane	184	C13H28	000629-50-5	86
5		Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	80



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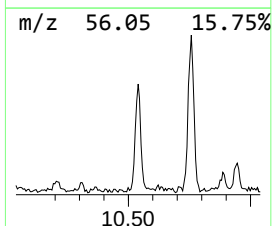
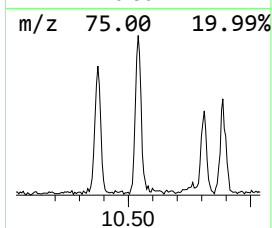
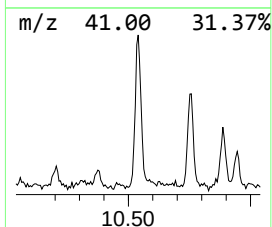
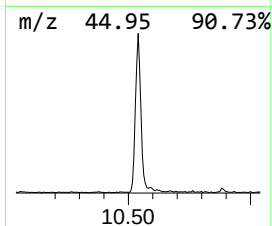
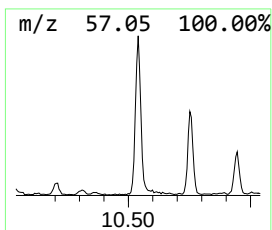
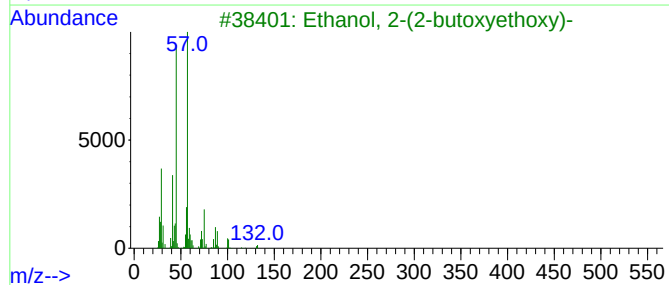
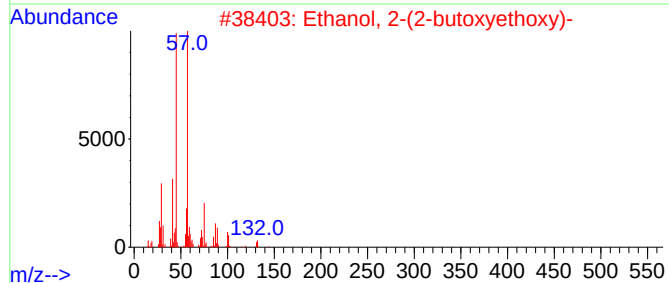
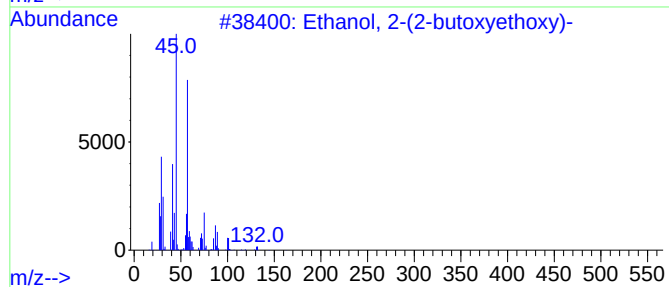
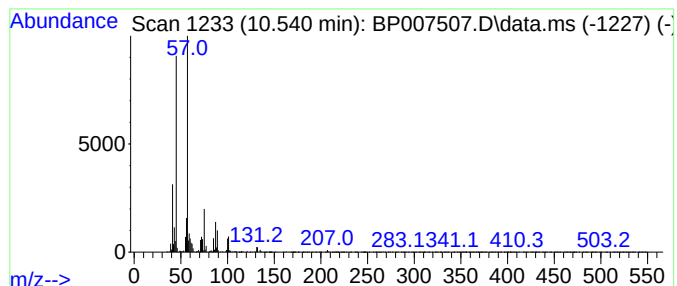
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.540	2.90 ng/ul	239934	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	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			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007507.D
Acq On : 20 Oct 2021 00:59
Operator : CG/JU
Sample : M4207-09
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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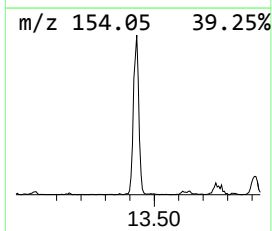
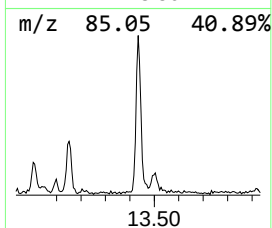
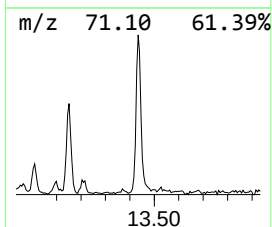
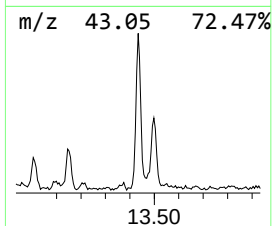
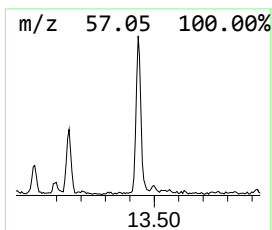
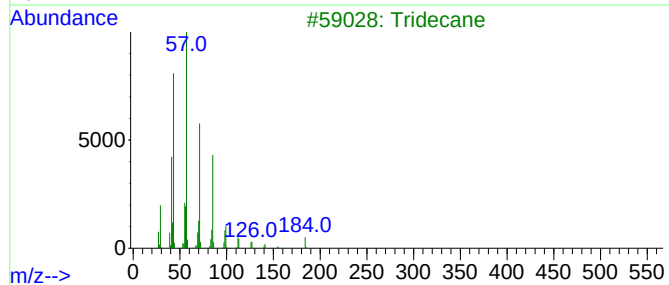
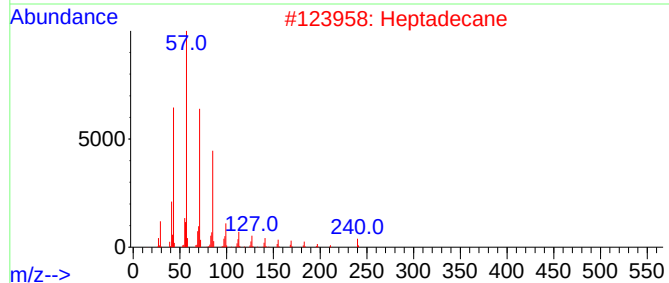
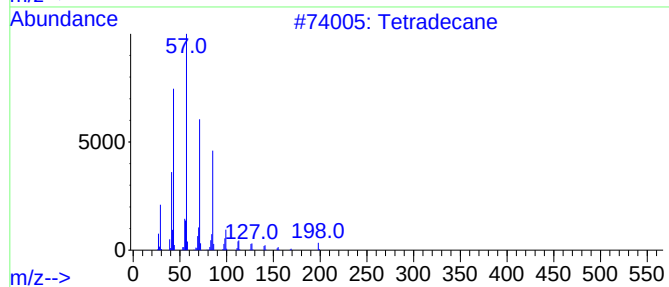
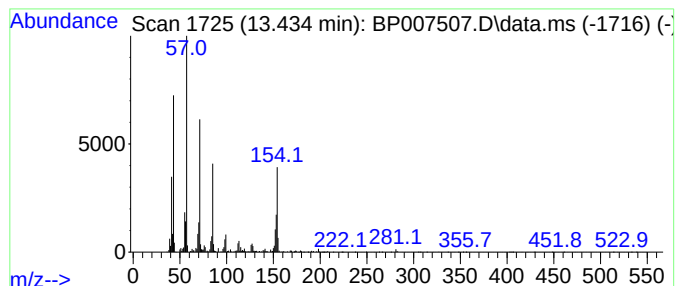
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.434	2.03 ng/ul	222144	Acenaphthene-d10	14.587

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	90
2			Heptadecane	240	C17H36	000629-78-7	60
3			Tridecane	184	C13H28	000629-50-5	53
4			Tridecane, 6-propyl-	226	C16H34	055045-10-8	53
5			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007507.D
Acq On : 20 Oct 2021 00:59
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Sample : M4207-09
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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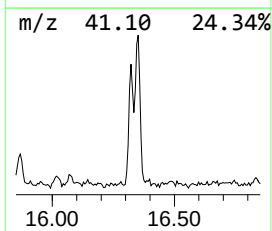
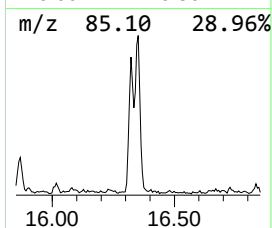
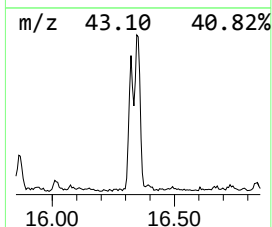
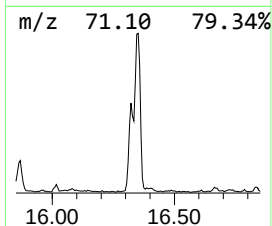
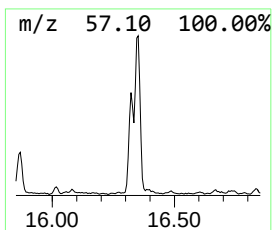
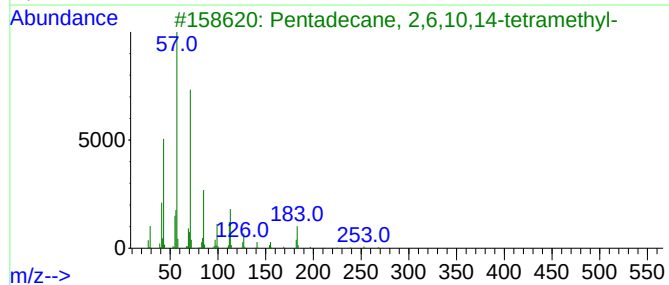
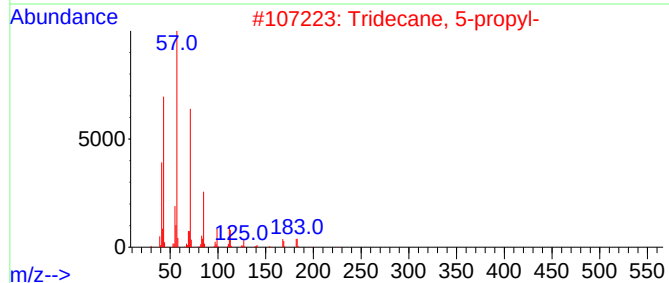
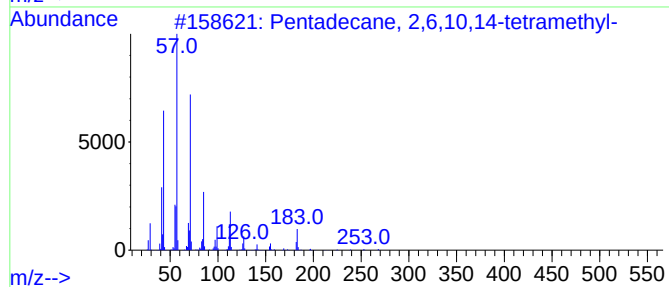
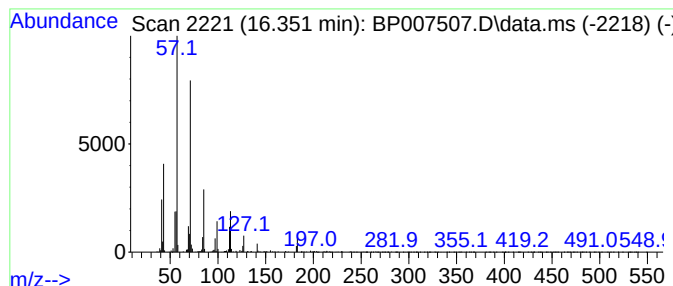
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.351	2.51 ng/ul	328726	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
2			Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
3			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
4			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	83
5			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007507.D
Acq On : 20 Oct 2021 00:59
Operator : CG/JU
Sample : M4207-09
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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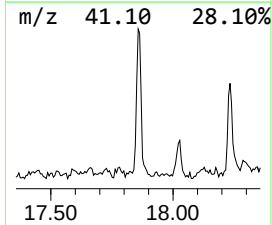
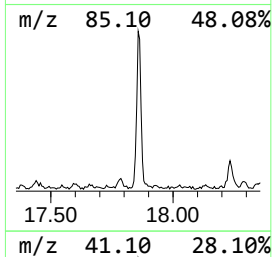
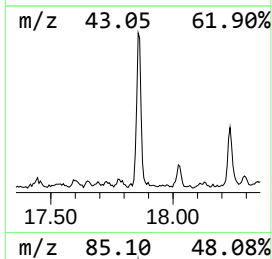
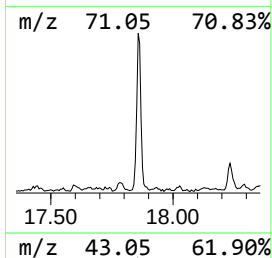
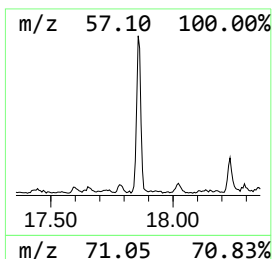
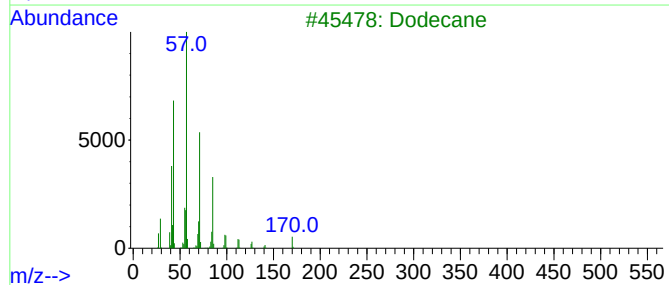
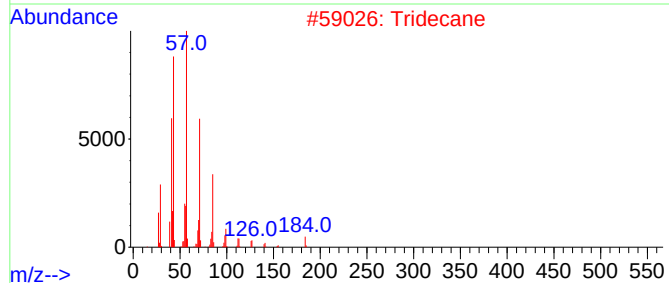
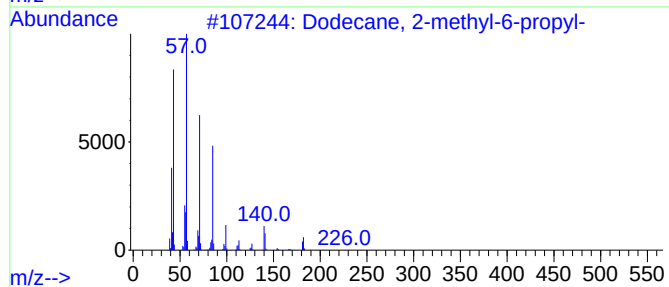
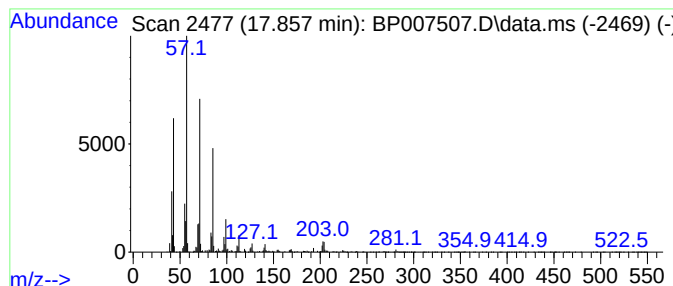
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.857	2.14 ng/ul	281063	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	91
2			Tridecane	184	C13H28	000629-50-5	90
3			Dodecane	170	C12H26	000112-40-3	87
4			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
5			Heptadecane	240	C17H36	000629-78-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007507.D
Acq On : 20 Oct 2021 00:59
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Sample : M4207-09
Misc :
ALS Vial : 60 Sample Multiplier: 1

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ClientSampleId :
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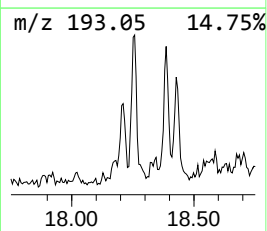
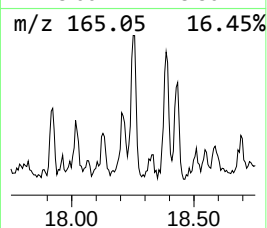
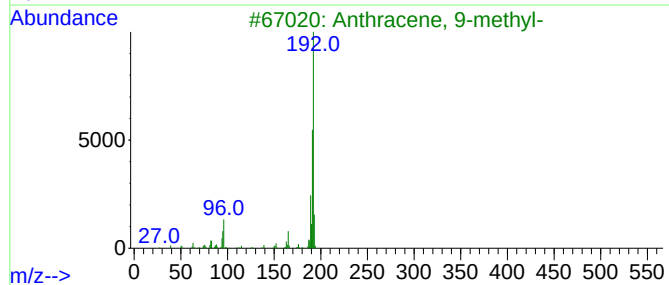
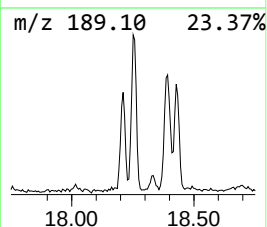
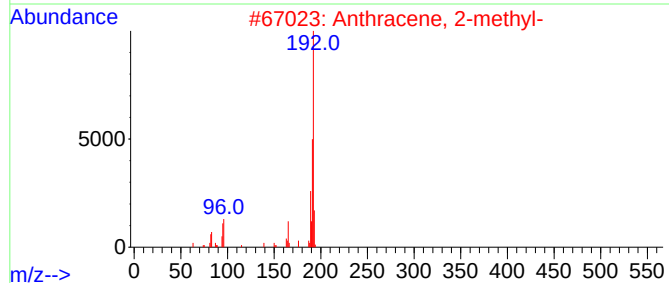
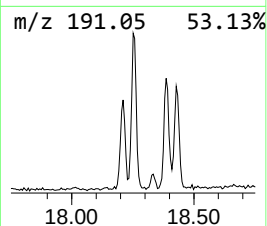
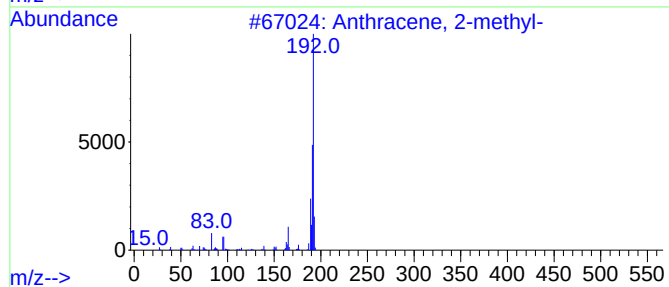
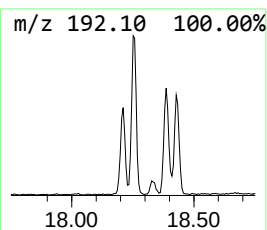
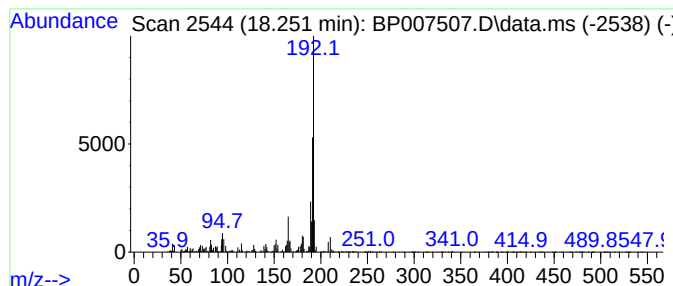
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Anthracene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	3.27 ng/ul	428720	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	81
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	81
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007507.D
Acq On : 20 Oct 2021 00:59
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Sample : M4207-09
Misc :
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Instrument :
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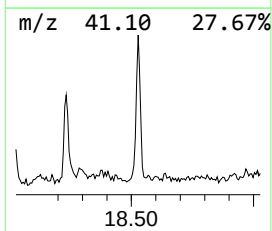
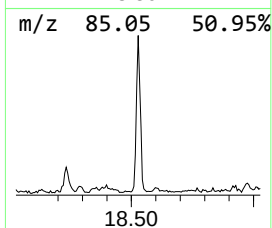
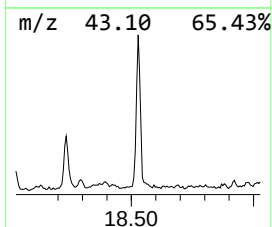
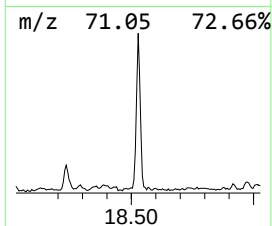
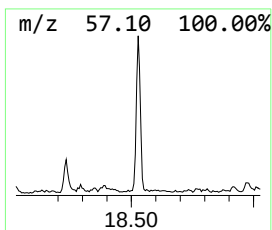
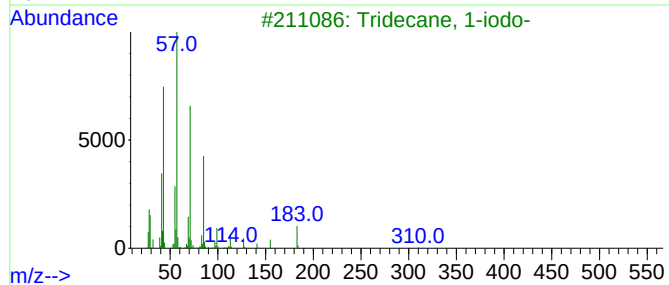
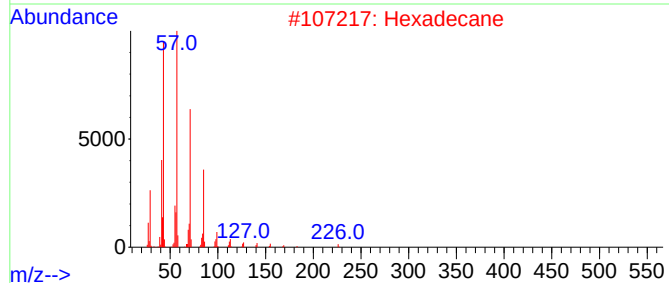
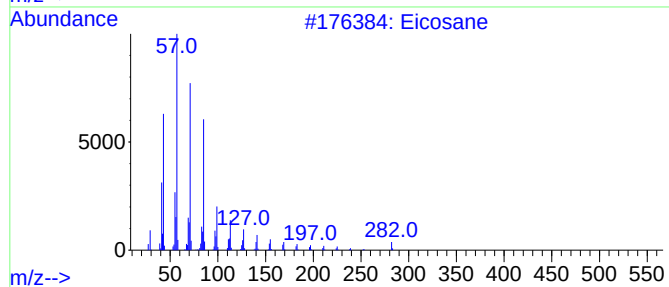
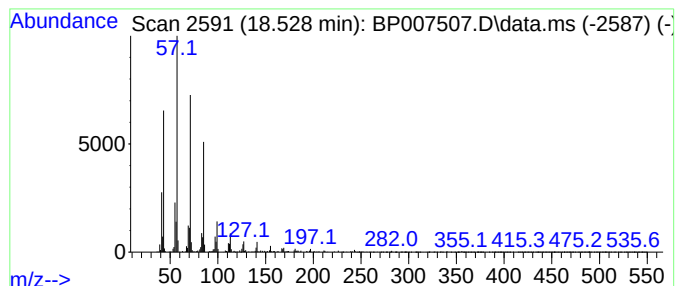
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.528	2.01 ng/ul	264080	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Hexadecane	226	C16H34	000544-76-3	96
3			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
4			Heptadecane	240	C17H36	000629-78-7	91
5			Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
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Misc :
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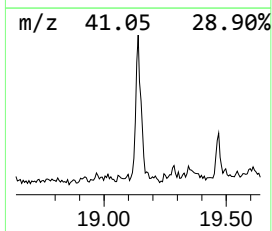
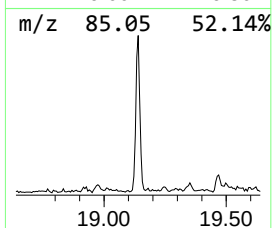
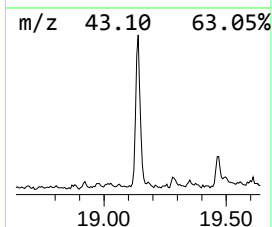
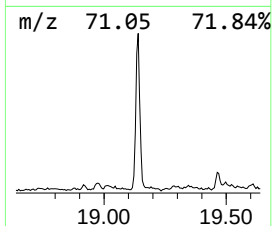
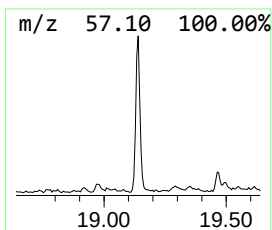
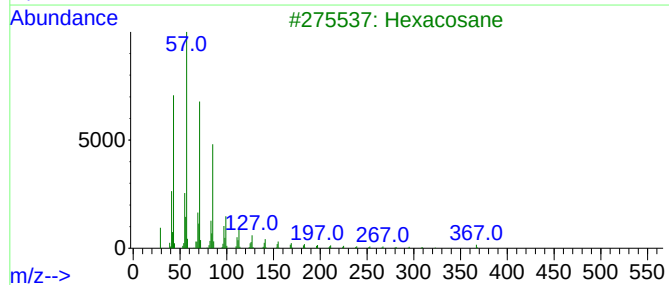
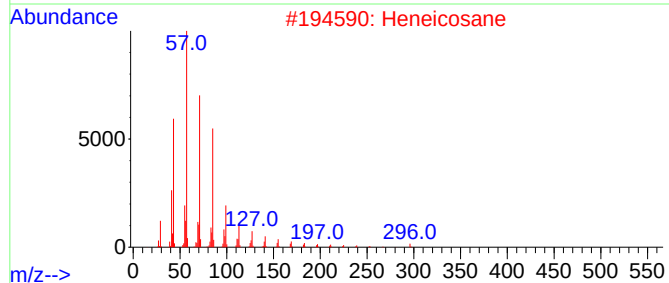
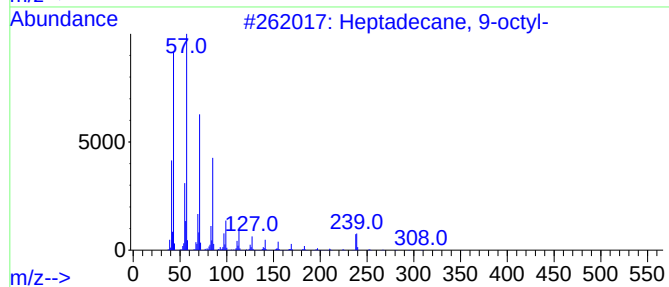
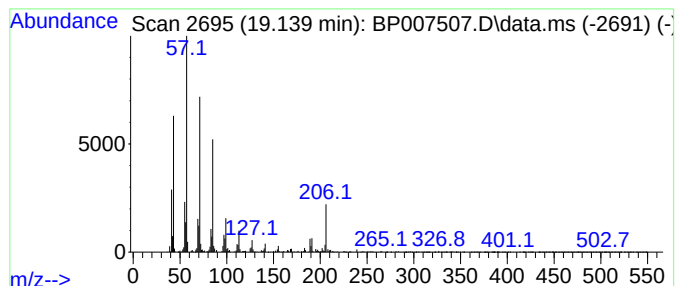
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.139	3.29 ng/ul	431614	Phenanthrene-d10	17.339	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	89
2	Heneicosane	296	C21H44	000629-94-7	81
3	Hexacosane	366	C26H54	000630-01-3	81
4	Octadecane	254	C18H38	000593-45-3	74
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007507.D
Acq On : 20 Oct 2021 00:59
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Sample : M4207-09
Misc :
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Instrument :
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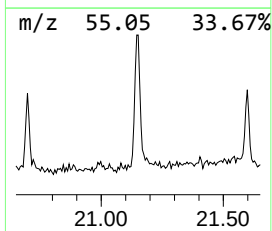
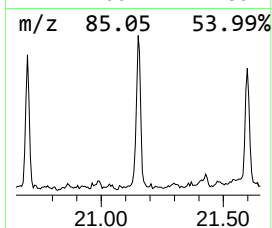
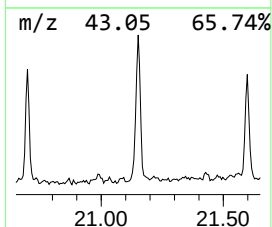
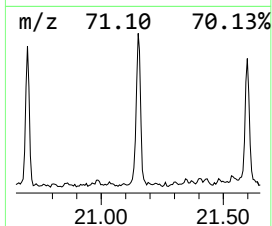
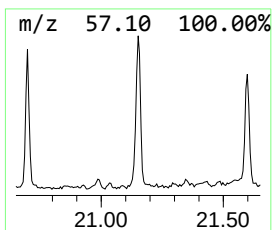
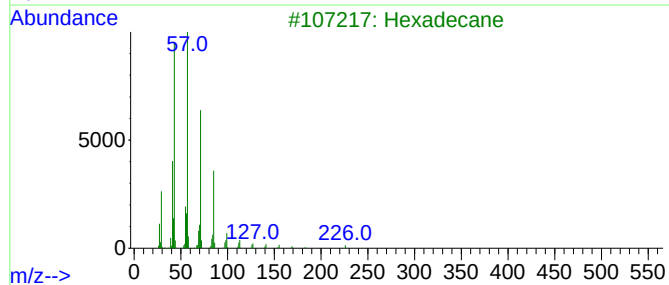
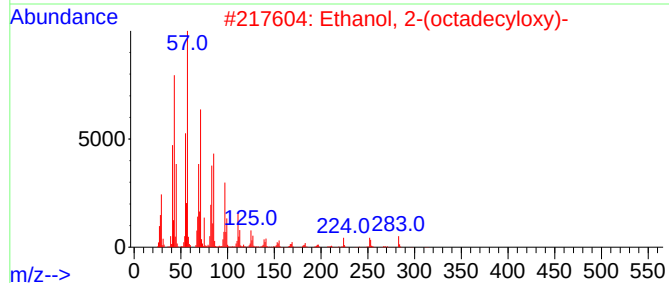
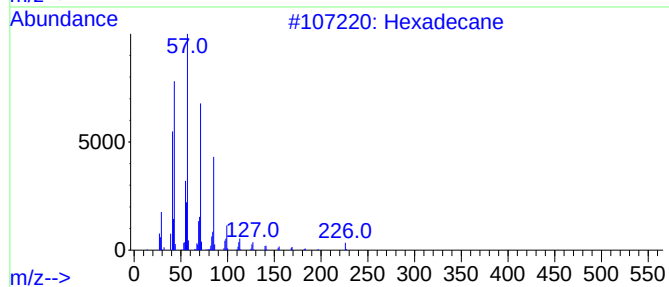
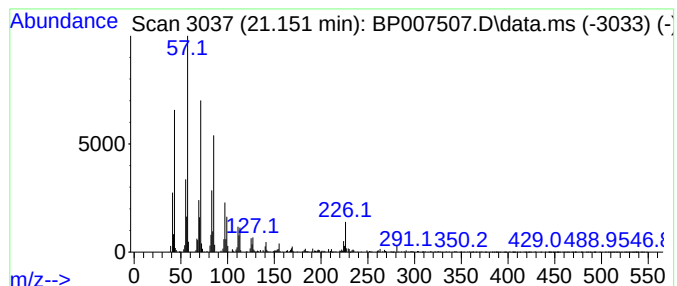
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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.
21.151	2.45 ng/ul	402532	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	95
2			Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	93
3			Hexadecane	226	C16H34	000544-76-3	92
4			Hexadecane	226	C16H34	000544-76-3	90
5			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007507.D
 Acq On : 20 Oct 2021 00:59
 Operator : CG/JU
 Sample : M4207-09
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.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
Benzene, 1,2,3-...	7.599	3.1	ng/u1	169775	1	7.952	1106490	20.0
(DEL) Alkane: S...	9.205	2.3	ng/u1	125141	1	7.952	1106490	20.0
Ethanol, 2-(2-b...	10.540	2.9	ng/u1	239934	2	10.757	1652870	20.0
(DEL) Alkane: S...	13.434	2.0	ng/u1	222144	3	14.587	2189880	20.0
(DEL) Alkane: S...	16.351	2.5	ng/u1	328726	4	17.339	2624170	20.0
(DEL) Alkane: S...	17.857	2.1	ng/u1	281063	4	17.339	2624170	20.0
Anthracene, 2-m...	18.251	3.3	ng/u1	428720	4	17.339	2624170	20.0
(DEL) Alkane: S...	18.528	2.0	ng/u1	264080	4	17.339	2624170	20.0
(DEL) Alkane: S...	19.139	3.3	ng/u1	431614	4	17.339	2624170	20.0
(DEL) Alkane: S...	21.151	2.5	ng/u1	402532	5	21.427	3284990	20.0