

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
 Data File : BP017248.D
 Acq On : 20 Sep 2023 20:17
 Operator : MA/JU
 Sample : 04330-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYX6

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-BP092023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017248.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	35	38	43	rVB	21993	24790	0.98%	0.087%
2	3.852	127	130	134	rVB	44742	53700	2.13%	0.188%
3	3.934	139	144	151	rVB2	68282	103008	4.08%	0.360%
4	4.052	161	164	169	rVB2	17366	25086	0.99%	0.088%
5	4.275	199	202	207	rVB2	25035	33904	1.34%	0.119%
6	4.440	227	230	235	rVB5	12059	18548	0.73%	0.065%
7	4.634	260	263	269	rVB8	13623	18822	0.75%	0.066%
8	4.693	269	273	279	rBV3	16362	25954	1.03%	0.091%
9	4.987	319	323	328	rVB	38471	55258	2.19%	0.193%
10	5.111	340	344	347	rBV5	9435	11944	0.47%	0.042%
11	5.240	361	366	370	rBV8	5679	10882	0.43%	0.038%
12	5.752	448	453	459	rBV4	8088	11602	0.46%	0.041%
13	5.893	469	477	484	rVB5	17144	33170	1.31%	0.116%
14	6.781	624	628	636	rVB5	13225	18134	0.72%	0.063%
15	6.893	643	647	652	rBV7	8966	15432	0.61%	0.054%
16	7.093	675	681	695	rBV	435578	775954	30.74%	2.715%
17	7.281	705	713	721	rBV	368634	571747	22.65%	2.000%
18	7.463	733	744	753	rBV	406925	680748	26.97%	2.382%
19	7.616	766	770	778	rVB10	9637	20008	0.79%	0.070%
20	7.940	817	825	834	rBV	828375	1295596	51.33%	4.533%
21	8.652	939	946	955	rBV2	236392	405719	16.07%	1.420%
22	8.793	962	970	977	rBV	170551	275249	10.90%	0.963%
23	8.934	985	994	1001	rBV2	18094	44421	1.76%	0.155%
24	9.010	1004	1007	1011	rVB5	7938	12536	0.50%	0.044%
25	9.140	1015	1029	1039	rBV	407930	693418	27.47%	2.426%
26	9.222	1039	1043	1045	rVV3	13418	19215	0.76%	0.067%
27	9.257	1045	1049	1052	rVB5	10850	17361	0.69%	0.061%
28	9.340	1059	1063	1069	rVB8	5457	12296	0.49%	0.043%
29	9.616	1104	1110	1116	rBV8	7208	18899	0.75%	0.066%
30	9.852	1136	1150	1160	rBV	291341	518857	20.56%	1.815%
31	9.999	1168	1175	1180	rBV2	96801	189232	7.50%	0.662%
32	10.222	1206	1213	1219	rVB2	40265	72479	2.87%	0.254%
33	10.381	1232	1240	1250	rVB	400915	698144	27.66%	2.443%
34	10.587	1271	1275	1280	rVB8	8916	14534	0.58%	0.051%
35	10.657	1281	1287	1290	rVB3	17575	26373	1.04%	0.092%
36	10.763	1298	1305	1312	rBV	1061197	1738217	68.86%	6.082%

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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-BP092023.MA.M
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37	10.828	1312	1316	1324	rVB3	21465	38529	1.53%	0.135%
38	10.928	1326	1333	1342	rBV	275159	508991	20.16%	1.781%
39	11.175	1370	1375	1381	rVB5	16199	29394	1.16%	0.103%
40	11.328	1393	1401	1405	rBV10	15489	39052	1.55%	0.137%

41	11.522	1430	1434	1440	rBV7	12094	26315	1.04%	0.092%
42	11.593	1441	1446	1451	rVV2	19733	41258	1.63%	0.144%
43	11.698	1458	1464	1468	rBV2	39309	59004	2.34%	0.206%
44	11.751	1470	1473	1478	rVV6	13186	23846	0.94%	0.083%
45	11.804	1478	1482	1487	rVB8	8478	14860	0.59%	0.052%

46	11.898	1495	1498	1504	rBV6	13976	22064	0.87%	0.077%
47	11.969	1504	1510	1516	rBV4	24481	45953	1.82%	0.161%
48	12.110	1529	1534	1538	rVB	41304	56844	2.25%	0.199%
49	12.169	1539	1544	1547	rBV4	11359	20388	0.81%	0.071%
50	12.234	1552	1555	1559	rBV6	7271	12130	0.48%	0.042%

51	12.422	1583	1587	1591	rBV6	15297	26840	1.06%	0.094%
52	12.516	1599	1603	1604	rBV4	14612	18075	0.72%	0.063%
53	12.640	1622	1624	1630	rBV6	7720	12867	0.51%	0.045%
54	12.745	1637	1642	1649	rBV10	20066	43754	1.73%	0.153%
55	12.916	1669	1671	1677	rVB6	11113	14846	0.59%	0.052%

56	13.057	1690	1695	1700	rVB2	93789	133990	5.31%	0.469%
57	13.351	1736	1745	1751	rBV2	105482	222629	8.82%	0.779%
58	13.440	1756	1760	1765	rVB4	35966	71419	2.83%	0.250%
59	13.745	1808	1812	1815	rBV5	30007	51158	2.03%	0.179%
60	13.969	1847	1850	1851	rBV2	40174	43776	1.73%	0.153%

61	14.016	1852	1858	1863	rVB	915521	1397368	55.36%	4.889%
62	14.298	1900	1906	1913	rBV	1050994	1646106	65.21%	5.760%
63	14.428	1925	1928	1937	rVB	154132	217278	8.61%	0.760%
64	14.598	1952	1957	1969	rVB	1393165	2144158	84.94%	7.502%
65	14.834	1994	1997	2001	rBV3	60411	90280	3.58%	0.316%

66	14.928	2010	2013	2018	rVB	57736	61800	2.45%	0.216%
67	15.045	2029	2033	2040	rBV4	87698	153349	6.08%	0.537%
68	15.387	2088	2091	2096	rVB	236953	265817	10.53%	0.930%
69	15.598	2122	2127	2134	rVB	1288200	1834867	72.69%	6.420%
70	15.787	2154	2159	2164	rBV	320107	510123	20.21%	1.785%

71	16.269	2235	2241	2246	rBV	884094	1478528	58.57%	5.173%
72	17.057	2372	2375	2378	rBV	295810	343227	13.60%	1.201%
73	17.098	2378	2382	2387	rVB	742060	1076802	42.66%	3.768%
74	17.375	2424	2429	2441	rVB	1652324	2524211	100.00%	8.832%
75	17.475	2441	2446	2451	rBV	1220196	1781734	70.59%	6.234%

76	17.804	2499	2502	2508	rVB	320972	394005	15.61%	1.379%
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ClientSampleId :
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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-BP092023.MA.M
Title : SVOA CALIBRATION

77	19.716	2823	2827	2832	rBV	1840952	2521778	99.90%	8.823%
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Sum of corrected areas: 28580650

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Misc :
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Instrument :
BNA_P
ClientSampleId :
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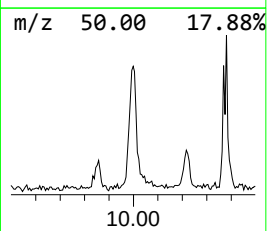
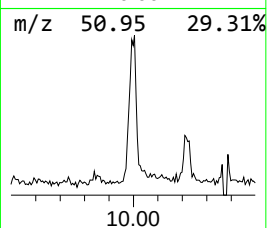
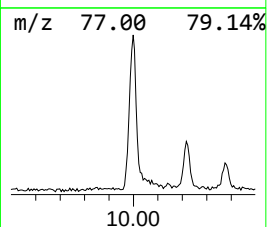
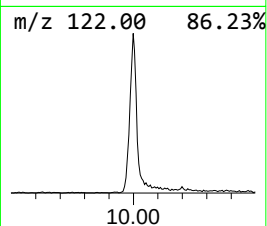
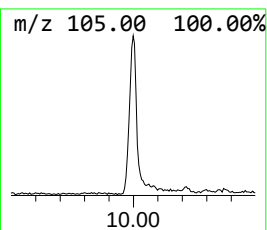
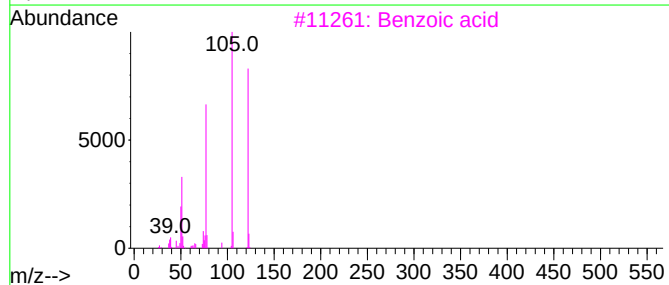
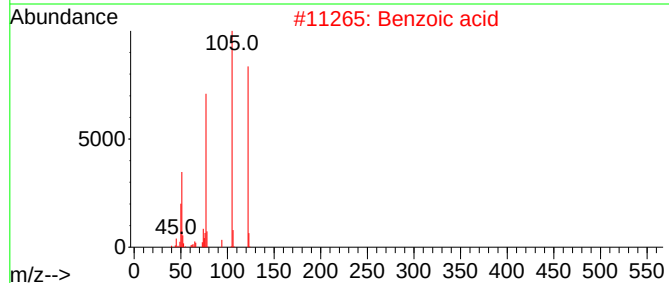
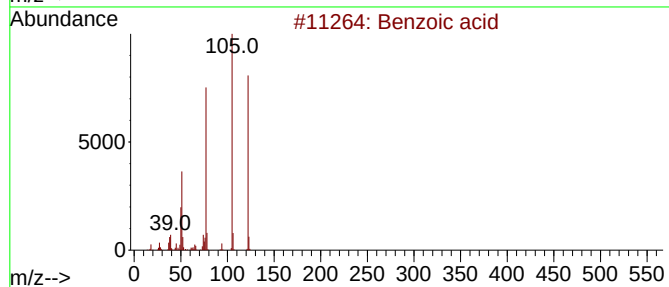
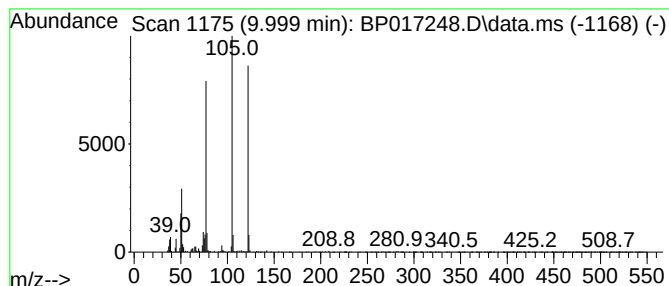
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.999	2.18 ng/ul	189232	Naphthalene-d8	10.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid	122	C7H6O2	000065-85-0	97
2			Benzoic acid	122	C7H6O2	000065-85-0	96
3			Benzoic acid	122	C7H6O2	000065-85-0	95
4			Benzoic acid	122	C7H6O2	000065-85-0	95
5			.beta.-1,5-O-Dibenzoyl-ribofuranose	358	C19H18O7	1000129-71-8	91



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ALS Vial : 17 Sample Multiplier: 1

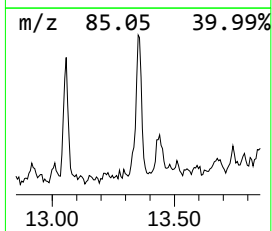
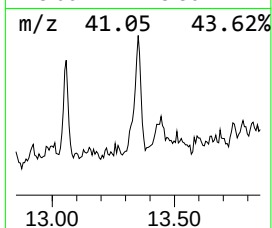
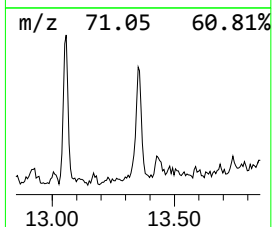
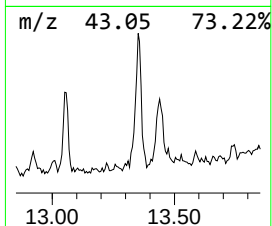
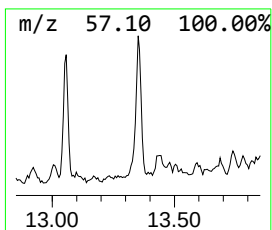
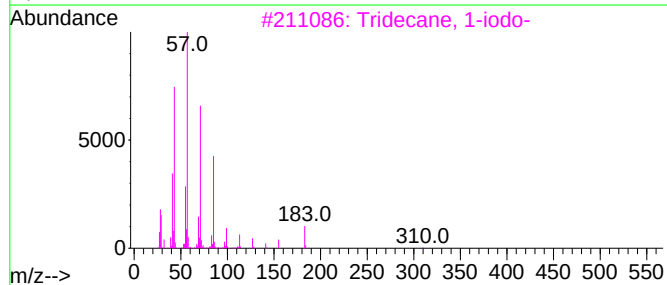
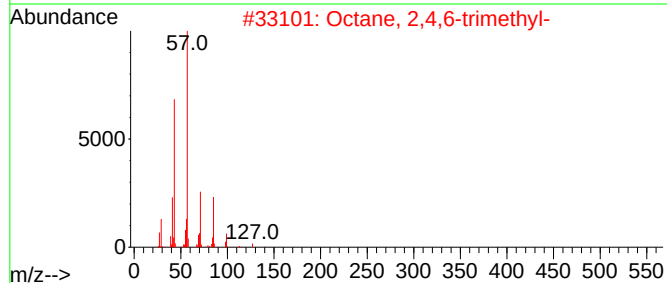
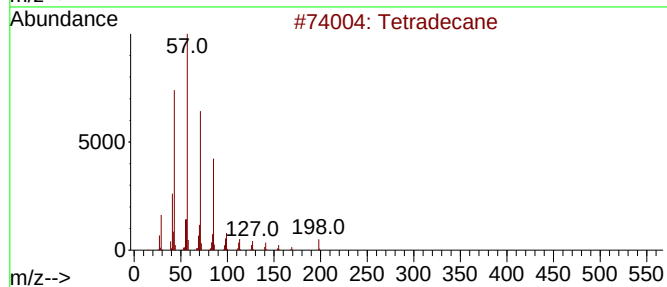
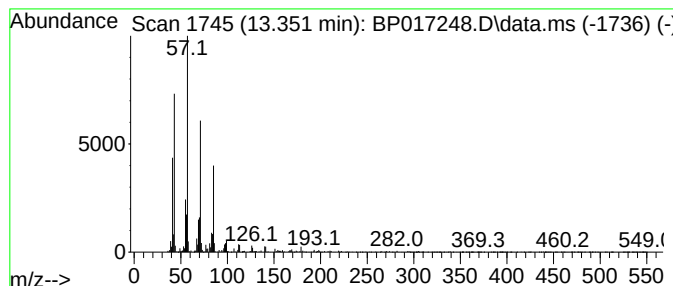
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.351	2.08 ng/ul	222629	Acenaphthene-d10	14.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	86
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
4	10-Methylnonadecane	282	C20H42	056862-62-5	72
5	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	72



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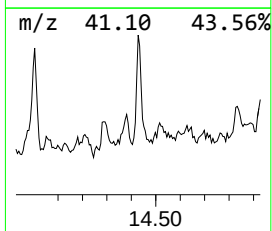
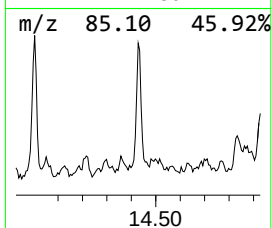
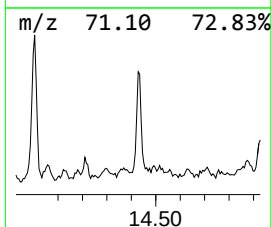
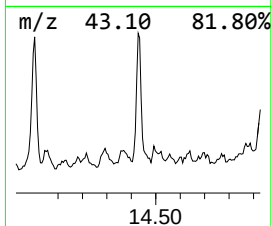
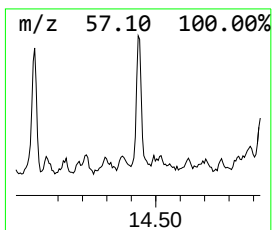
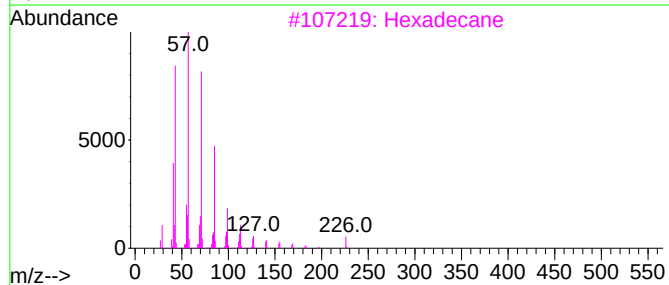
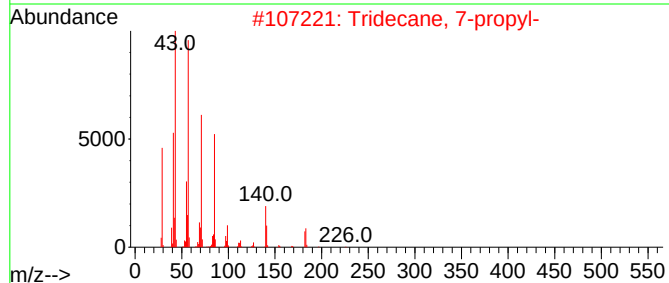
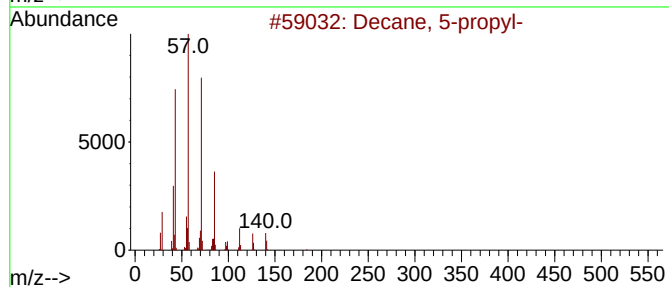
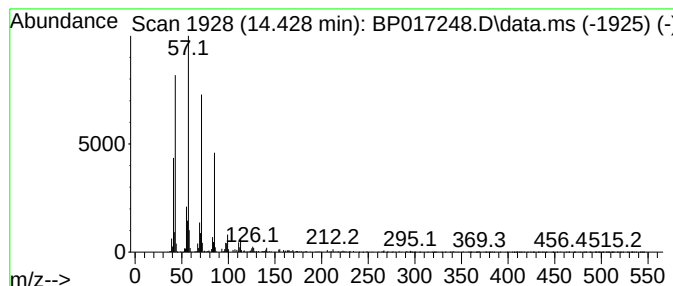
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.428	2.03 ng/ul	217278	Acenaphthene-d10	14.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5-propyl-	184	C13H28	017312-62-8	86
2	Tridecane, 7-propyl-	226	C16H34	055045-09-5	81
3	Hexadecane	226	C16H34	000544-76-3	81
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	80
5	Hexadecane	226	C16H34	000544-76-3	76



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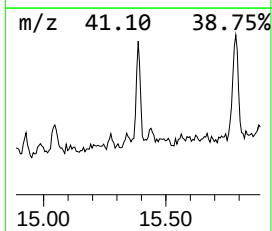
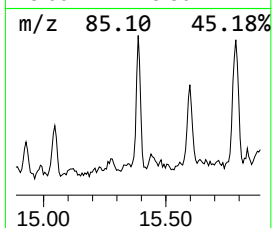
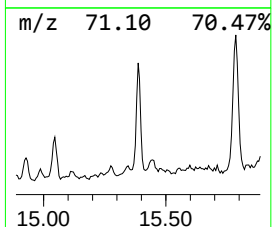
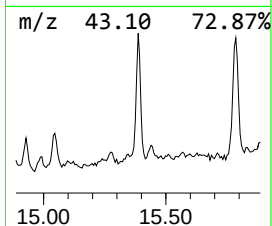
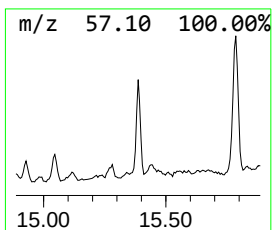
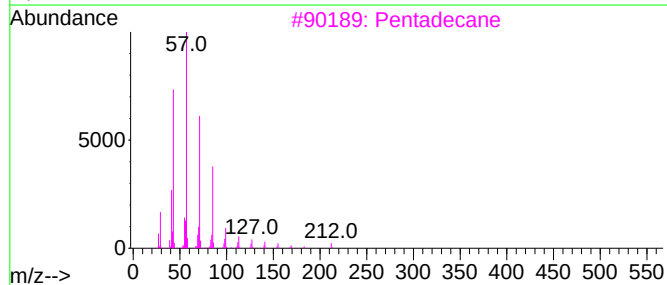
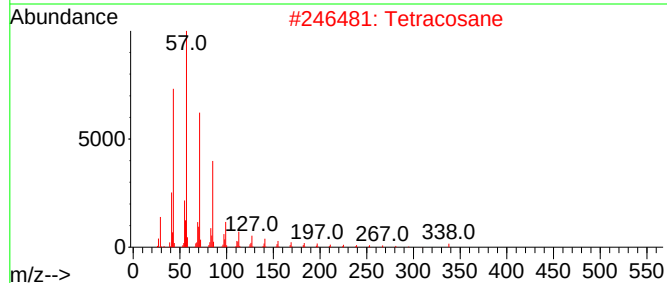
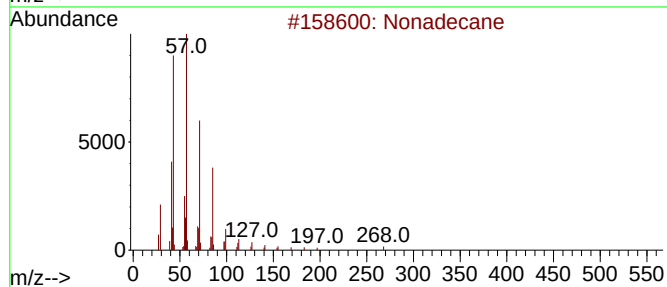
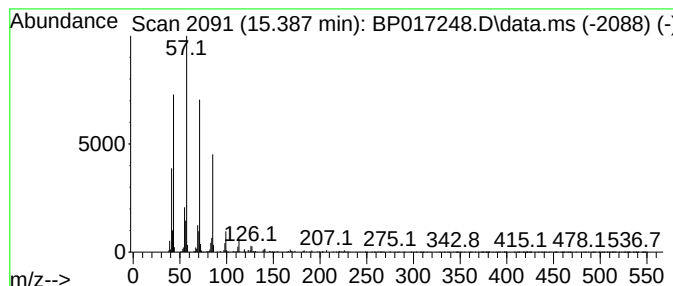
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.387	2.48 ng/ul	265817	Acenaphthene-d10	14.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	90
2			Tetracosane	338	C24H50	000646-31-1	90
3			Pentadecane	212	C15H32	000629-62-9	86
4			Hexadecane	226	C16H34	000544-76-3	80
5			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017248.D
Acq On : 20 Sep 2023 20:17
Operator : MA/JU
Sample : 04330-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZYX6

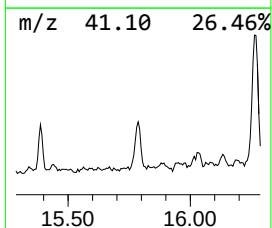
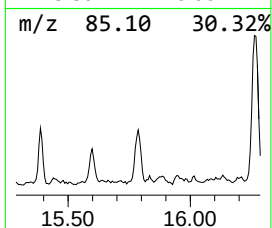
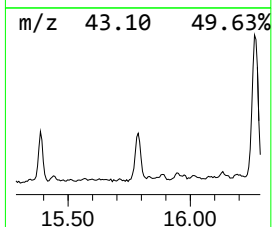
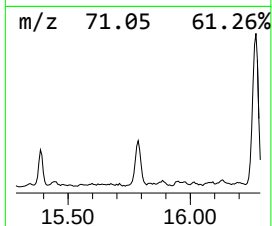
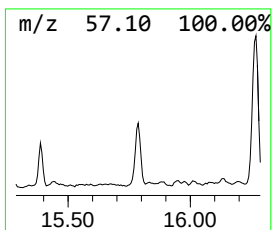
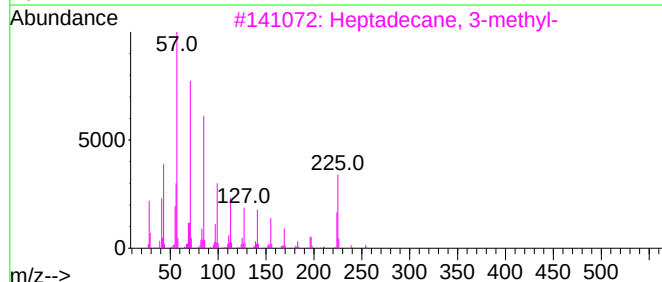
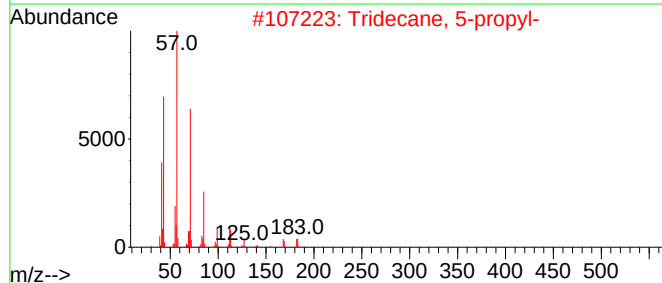
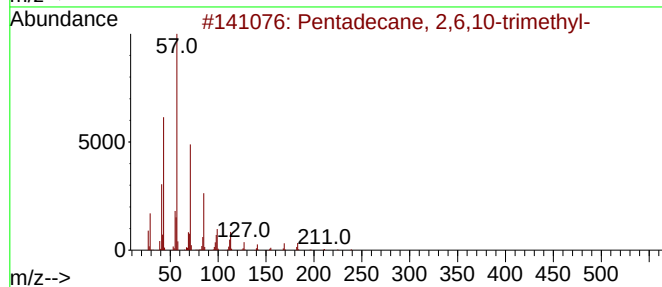
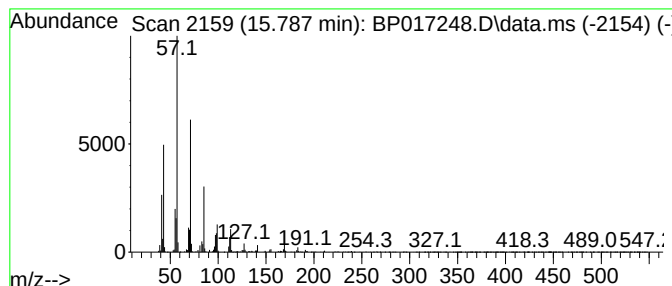
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.787	4.76 ng/ul	510123	Acenaphthene-d10	14.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	94
2			Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
3			Heptadecane, 3-methyl-	254	C18H38	006418-44-6	80
4			Heneicosane	296	C21H44	000629-94-7	72
5			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017248.D
Acq On : 20 Sep 2023 20:17
Operator : MA/JU
Sample : 04330-05
Misc :
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Instrument :
BNA_P
ClientSampleId :
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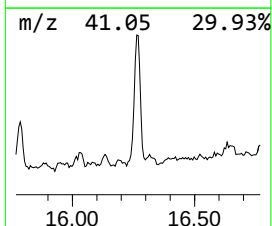
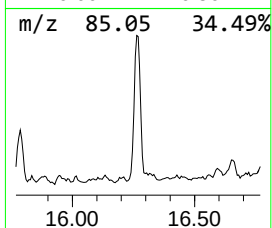
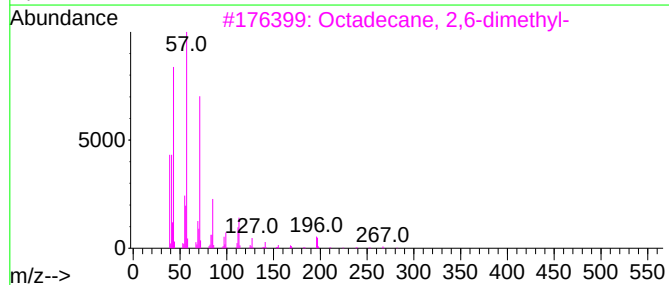
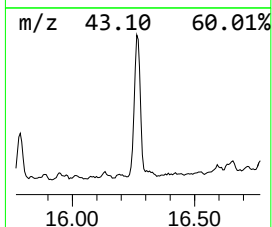
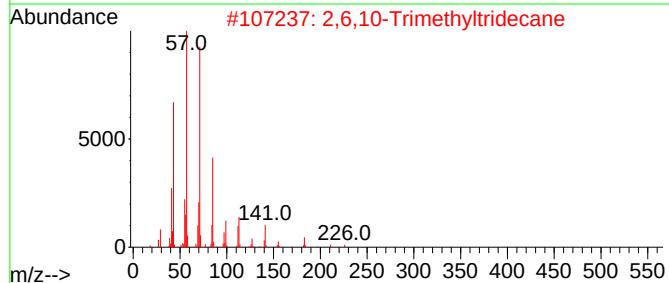
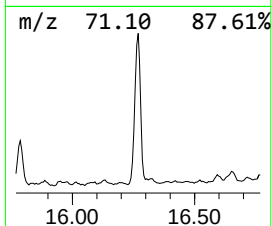
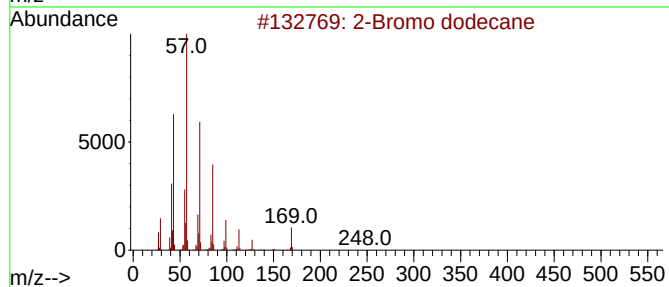
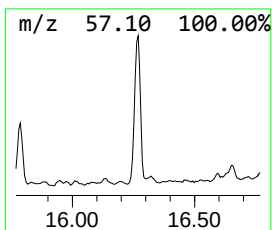
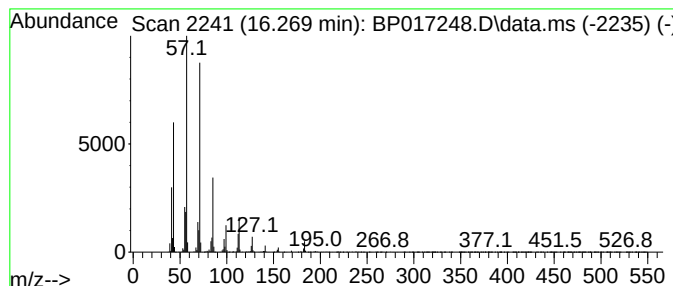
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Bromo dodecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.269	11.71 ng/ul	1478530	Phenanthrene-d10	17.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Bromo dodecane	248	C12H25Br	013187-99-0	91
2			2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	90
3			Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	90
4			Heptacosane	380	C27H56	000593-49-7	86
5			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017248.D
Acq On : 20 Sep 2023 20:17
Operator : MA/JU
Sample : 04330-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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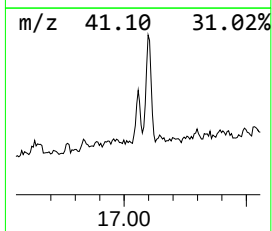
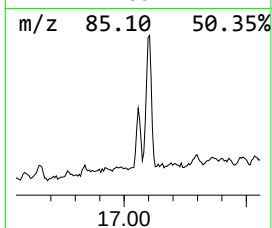
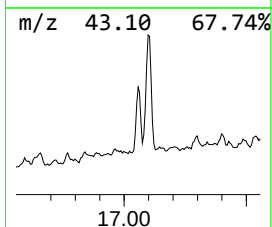
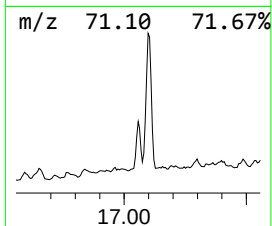
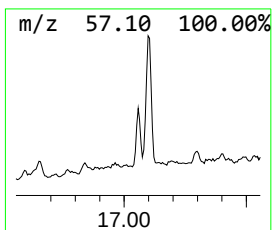
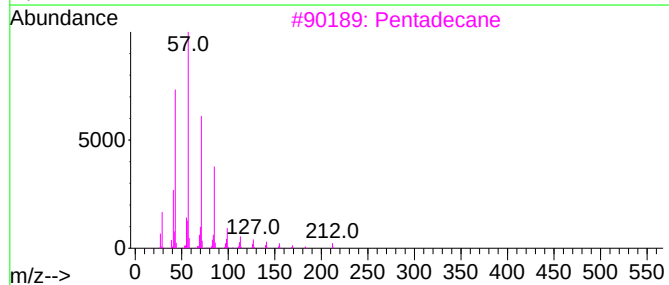
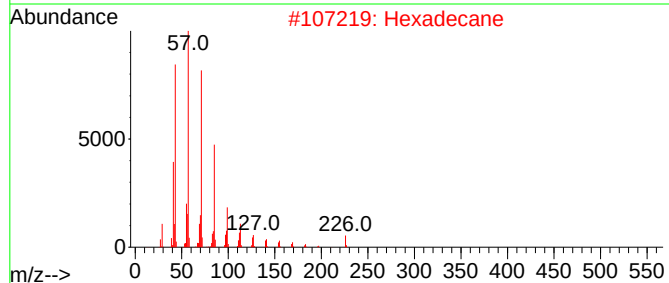
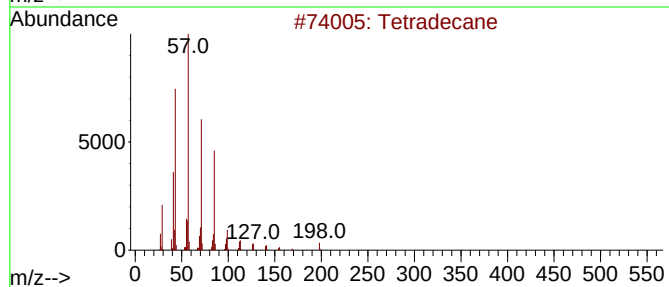
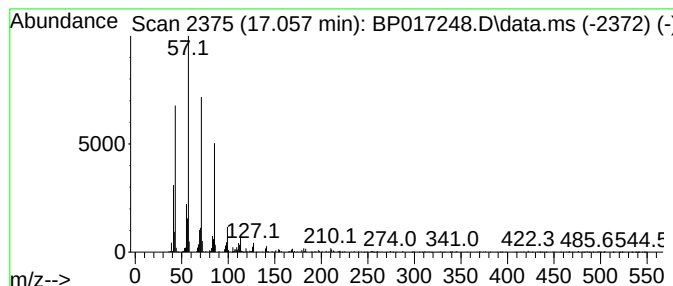
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.057	2.72 ng/ul	343227	Phenanthrene-d10	17.375

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	91
2		Hexadecane	226	C16H34	000544-76-3	90
3		Pentadecane	212	C15H32	000629-62-9	83
4		Nonadecane	268	C19H40	000629-92-5	83
5		Nonadecane	268	C19H40	000629-92-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017248.D
Acq On : 20 Sep 2023 20:17
Operator : MA/JU
Sample : 04330-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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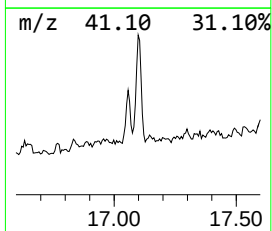
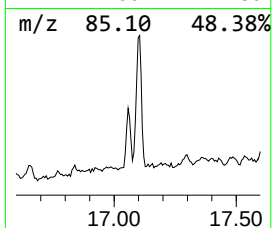
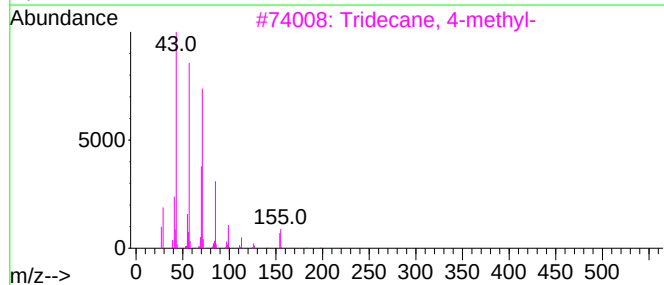
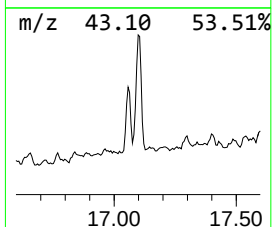
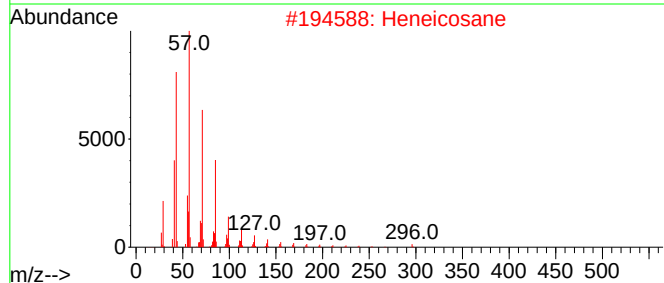
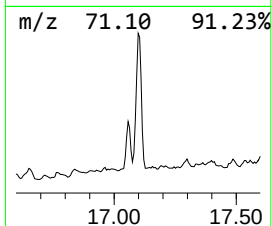
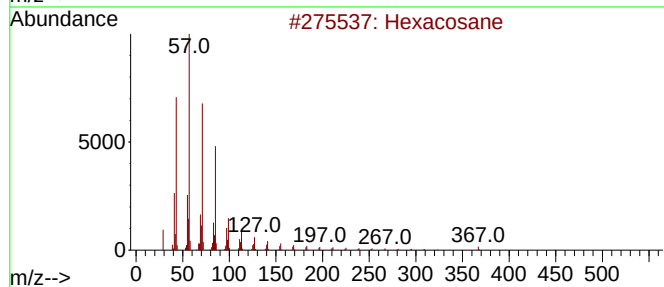
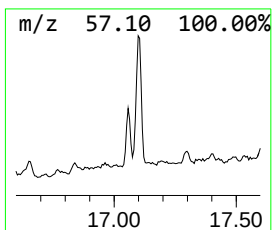
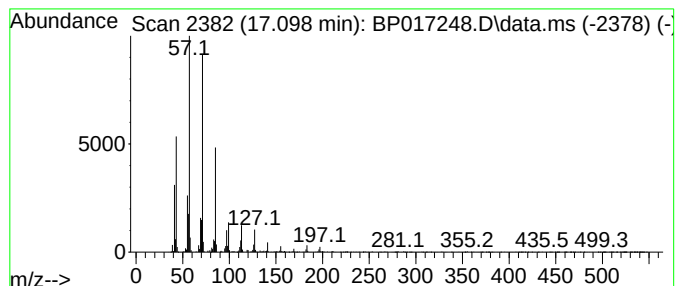
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.098	8.53 ng/ul	1076800	Phenanthrene-d10	17.375

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tridecane, 4-methyl-	198	C14H30	026730-12-1	87
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017248.D
Acq On : 20 Sep 2023 20:17
Operator : MA/JU
Sample : 04330-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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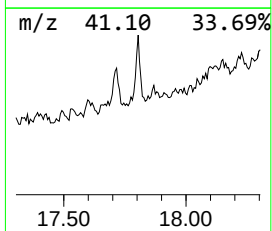
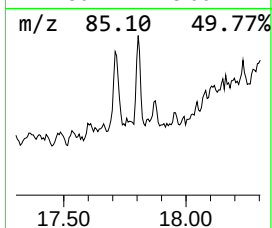
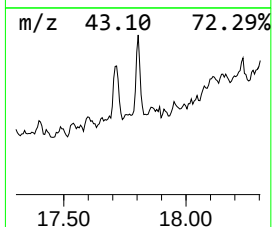
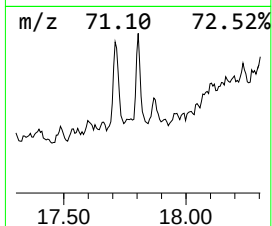
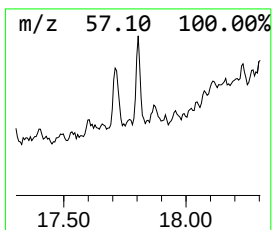
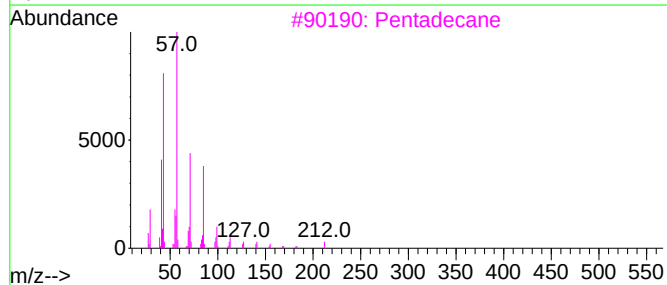
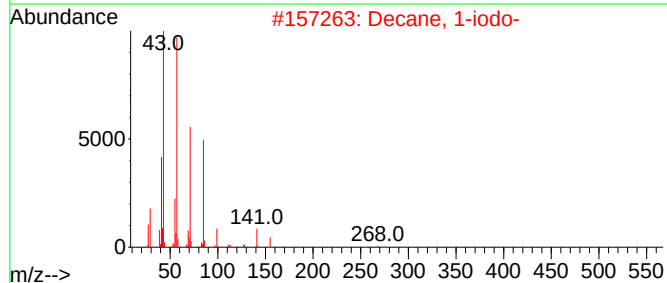
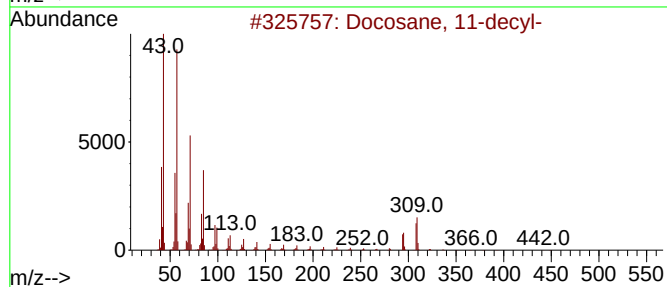
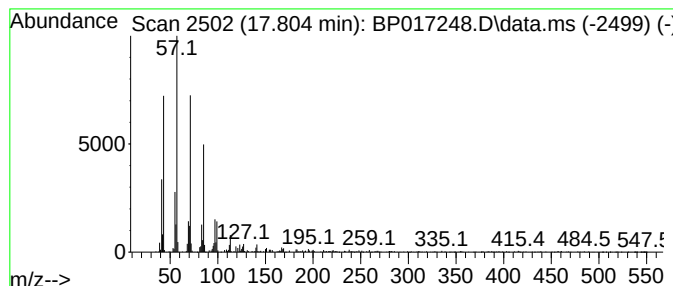
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.804	3.12 ng/ul	394005	Phenanthrene-d10	17.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 11-decyl-	451	C32H66	055401-55-3	89
2			Decane, 1-iodo-	268	C10H21I	002050-77-3	81
3			Pentadecane	212	C15H32	000629-62-9	81
4			Eicosane	282	C20H42	000112-95-8	80
5			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092023\
Data File : BP017248.D
Acq On : 20 Sep 2023 20:17
Operator : MA/JU
Sample : 04330-05
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: S...	13.351	2.1	ng/ul	222629	3	14.604	2144160	20.0
(DEL) Alkane: S...	14.428	2.0	ng/ul	217278	3	14.604	2144160	20.0
(DEL) Alkane: S...	15.387	2.5	ng/ul	265817	3	14.604	2144160	20.0
(DEL) Alkane: S...	15.787	4.8	ng/ul	510123	3	14.604	2144160	20.0
2-Bromo dodecane	16.269	11.7	ng/ul	1478530	4	17.375	2524210	20.0
(DEL) Alkane: S...	17.057	2.7	ng/ul	343227	4	17.375	2524210	20.0
(DEL) Alkane: S...	17.098	8.5	ng/ul	1076800	4	17.375	2524210	20.0
(DEL) Alkane: S...	17.804	3.1	ng/ul	394005	4	17.375	2524210	20.0