

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
 Data File : BP023455.D
 Acq On : 17 Dec 2024 05:57
 Operator : RC/JU
 Sample : P5234-15
 Misc : GCMS CONFIRMATION
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AR9

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-BP112624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023455.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	14	rVB	3046800	2719544	99.81%	10.590%
2	3.346	57	61	65	rVB2	30644	36662	1.35%	0.143%
3	3.811	136	140	146	rVB2	16307	23429	0.86%	0.091%
4	4.334	224	229	233	rVB8	4093	7649	0.28%	0.030%
5	4.623	274	278	283	rVB3	7136	9968	0.37%	0.039%
6	4.781	302	305	311	rBV6	4984	7760	0.28%	0.030%
7	5.093	354	358	365	rBV	15575	25511	0.94%	0.099%
8	5.222	375	380	388	rBV	40899	74478	2.73%	0.290%
9	5.375	403	406	410	rVB6	4597	7583	0.28%	0.030%
10	5.446	413	418	422	rBV8	4139	8538	0.31%	0.033%
11	5.587	435	442	451	rBV	101569	164433	6.03%	0.640%
12	5.828	478	483	492	rVB	4174	9210	0.34%	0.036%
13	6.046	511	520	526	rBV	37432	64693	2.37%	0.252%
14	6.528	598	602	607	rVB4	6329	9492	0.35%	0.037%
15	6.634	616	620	625	rBV2	15768	25871	0.95%	0.101%
16	6.693	625	630	638	rVB4	8689	17302	0.64%	0.067%
17	6.769	638	643	649	rBV	22946	34748	1.28%	0.135%
18	6.922	663	669	671	rBV5	7259	13259	0.49%	0.052%
19	7.011	678	684	690	rVB7	7849	19133	0.70%	0.075%
20	7.081	690	696	703	rVB10	5594	10060	0.37%	0.039%
21	7.146	703	707	708	rBV3	5739	6845	0.25%	0.027%
22	7.175	708	712	720	rVB	41772	65327	2.40%	0.254%
23	7.522	759	771	785	rBV	761345	1297929	47.64%	5.054%
24	7.652	785	793	799	rVB4	17268	45089	1.65%	0.176%
25	7.887	827	833	840	rVB	33614	58351	2.14%	0.227%
26	8.064	856	863	867	rBV10	3562	7569	0.28%	0.029%
27	8.152	871	878	888	rVB6	7521	16962	0.62%	0.066%
28	8.599	951	954	964	rVV7	7280	16105	0.59%	0.063%
29	8.716	964	974	983	rVB9	7503	23999	0.88%	0.093%
30	8.840	990	995	1002	rVB9	4593	9156	0.34%	0.036%
31	9.128	1039	1044	1048	rBV5	5100	7644	0.28%	0.030%
32	9.187	1048	1054	1059	rVB9	5412	11741	0.43%	0.046%
33	9.352	1076	1082	1088	rBV9	3252	6537	0.24%	0.025%
34	9.681	1132	1138	1144	rBV5	9667	20163	0.74%	0.079%
35	10.058	1196	1202	1207	rBV10	2851	6084	0.22%	0.024%
36	10.140	1210	1216	1220	rBV9	3213	6335	0.23%	0.025%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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37	10.275	1227	1239	1246	rBV	883253	1665800	61.14%	6.487%
38	10.428	1262	1265	1270	rVB4	7380	8243	0.30%	0.032%
39	11.281	1404	1410	1416	rBV5	9227	17301	0.63%	0.067%
40	11.540	1447	1454	1457	rBV9	4579	8225	0.30%	0.032%
41	11.693	1473	1480	1482	rBV8	5729	10672	0.39%	0.042%
42	11.952	1519	1524	1534	rBV4	7449	20345	0.75%	0.079%
43	12.169	1558	1561	1567	rVB4	10497	18351	0.67%	0.071%
44	12.475	1608	1613	1616	rBV7	4998	6525	0.24%	0.025%
45	12.546	1622	1625	1631	rVB8	7205	12162	0.45%	0.047%
46	12.646	1638	1642	1648	rVB3	12916	20554	0.75%	0.080%
47	12.710	1650	1653	1659	rVB8	4823	7458	0.27%	0.029%
48	12.893	1678	1684	1685	rBV5	5120	7365	0.27%	0.029%
49	12.946	1690	1693	1701	rVB5	9512	19620	0.72%	0.076%
50	13.193	1729	1735	1740	rBV10	8428	21587	0.79%	0.084%
51	13.322	1749	1757	1763	rBV9	10635	30089	1.10%	0.117%
52	13.463	1776	1781	1786	rBV3	12983	24879	0.91%	0.097%
53	13.528	1788	1792	1793	rBV3	13692	14367	0.53%	0.056%
54	13.616	1803	1807	1812	rBV2	21264	23547	0.86%	0.092%
55	13.657	1812	1814	1816	rBV3	6516	6265	0.23%	0.024%
56	13.875	1847	1851	1853	rBV5	9383	14092	0.52%	0.055%
57	13.963	1862	1866	1874	rVB7	20059	31454	1.15%	0.122%
58	14.040	1874	1879	1883	rBV4	14691	23188	0.85%	0.090%
59	14.140	1889	1896	1905	rVV	1234518	2194348	80.54%	8.545%
60	14.369	1930	1935	1943	rVB9	12203	31365	1.15%	0.122%
61	14.569	1961	1969	1976	rBV6	24367	64783	2.38%	0.252%
62	14.675	1984	1987	1992	rVB7	9320	13289	0.49%	0.052%
63	14.769	2000	2003	2011	rBV7	14318	28674	1.05%	0.112%
64	14.940	2029	2032	2034	rBV4	12438	14601	0.54%	0.057%
65	15.010	2042	2044	2049	rVB3	16896	21835	0.80%	0.085%
66	15.228	2077	2081	2085	rVB3	21206	31014	1.14%	0.121%
67	15.422	2110	2114	2124	rVB4	39112	81016	2.97%	0.315%
68	15.681	2155	2158	2160	rBV4	10168	14114	0.52%	0.055%
69	15.728	2164	2166	2168	rBV2	10470	11429	0.42%	0.045%
70	15.922	2189	2199	2205	rBV4	48335	141670	5.20%	0.552%
71	16.222	2247	2250	2256	rVB2	33890	51256	1.88%	0.200%
72	16.710	2330	2333	2337	rBV4	27841	32585	1.20%	0.127%
73	16.763	2339	2342	2349	rVB4	35385	56556	2.08%	0.220%
74	16.828	2349	2353	2356	rBV2	31239	44562	1.64%	0.174%
75	16.951	2368	2374	2379	rBV	1455869	2245958	82.43%	8.746%
76	16.998	2379	2382	2387	rVB	171616	261884	9.61%	1.020%

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77	17.134	2402	2405	2409	rBV6	33375	53126	1.95%	0.207%
78	17.575	2474	2480	2485	rVB4	73871	153655	5.64%	0.598%
79	17.757	2507	2511	2513	rBV5	25923	38764	1.42%	0.151%
80	17.834	2522	2524	2527	rBV4	22740	17520	0.64%	0.068%
81	17.869	2527	2530	2531	rBV	29293	31937	1.17%	0.124%
82	18.063	2558	2563	2565	rBV3	170555	249640	9.16%	0.972%
83	18.098	2565	2569	2577	rVV2	514454	868157	31.86%	3.381%
84	18.169	2579	2581	2588	rVB7	50102	79359	2.91%	0.309%
85	18.275	2593	2599	2606	rVB	1677080	2417150	88.71%	9.412%
86	18.351	2609	2612	2616	rBV2	46208	66194	2.43%	0.258%
87	18.475	2627	2633	2641	rBV	1628808	2304405	84.58%	8.973%
88	18.616	2654	2657	2664	rVB2	113497	144405	5.30%	0.562%
89	18.775	2680	2684	2690	rVB	138876	180491	6.62%	0.703%
90	18.939	2706	2712	2722	rVB4	112927	239961	8.81%	0.934%
91	19.098	2735	2739	2743	rBV	264161	352635	12.94%	1.373%
92	19.157	2745	2749	2755	rVV	158734	237019	8.70%	0.923%
93	19.239	2759	2763	2770	rVV2	160563	248328	9.11%	0.967%
94	19.710	2840	2843	2844	rBV2	71502	78617	2.89%	0.306%
95	19.986	2886	2890	2895	rBV5	81712	148978	5.47%	0.580%
96	21.386	3123	3128	3133	rBV	1261240	1865091	68.45%	7.263%
97	21.592	3161	3163	3168	rVB2	160382	207883	7.63%	0.809%
98	22.204	3264	3267	3272	rVB	237117	331167	12.15%	1.290%
99	22.886	3379	3383	3388	rVB	261863	393650	14.45%	1.533%
100	24.569	3661	3669	3686	rVB	833497	2724670	100.00%	10.610%

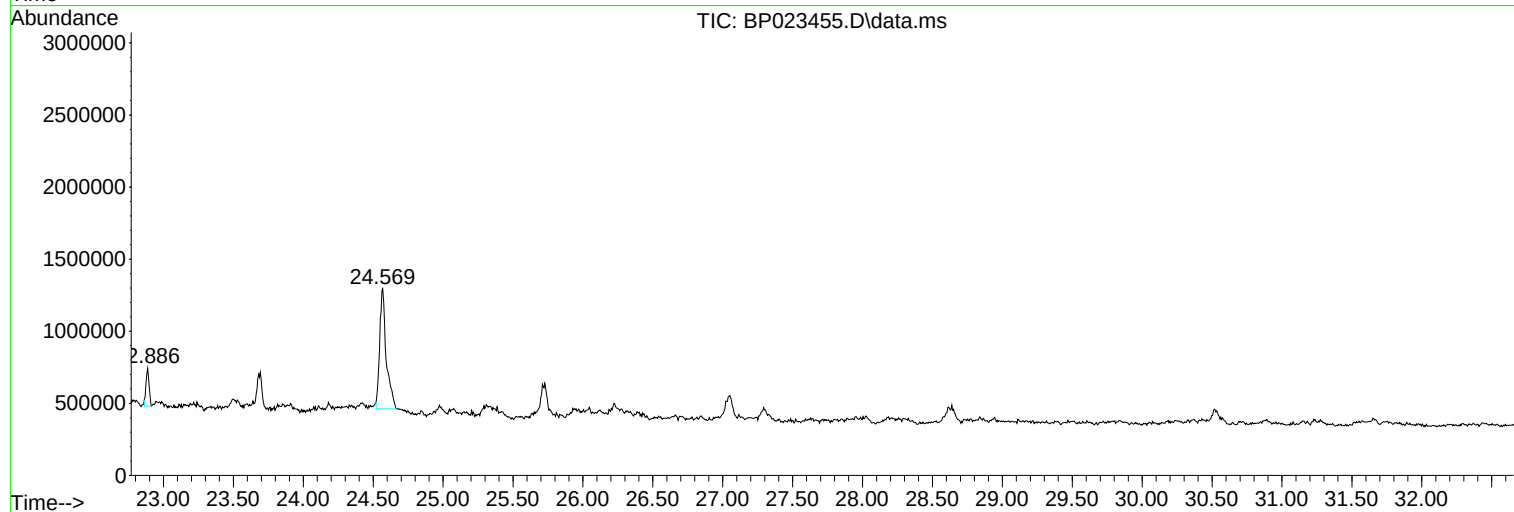
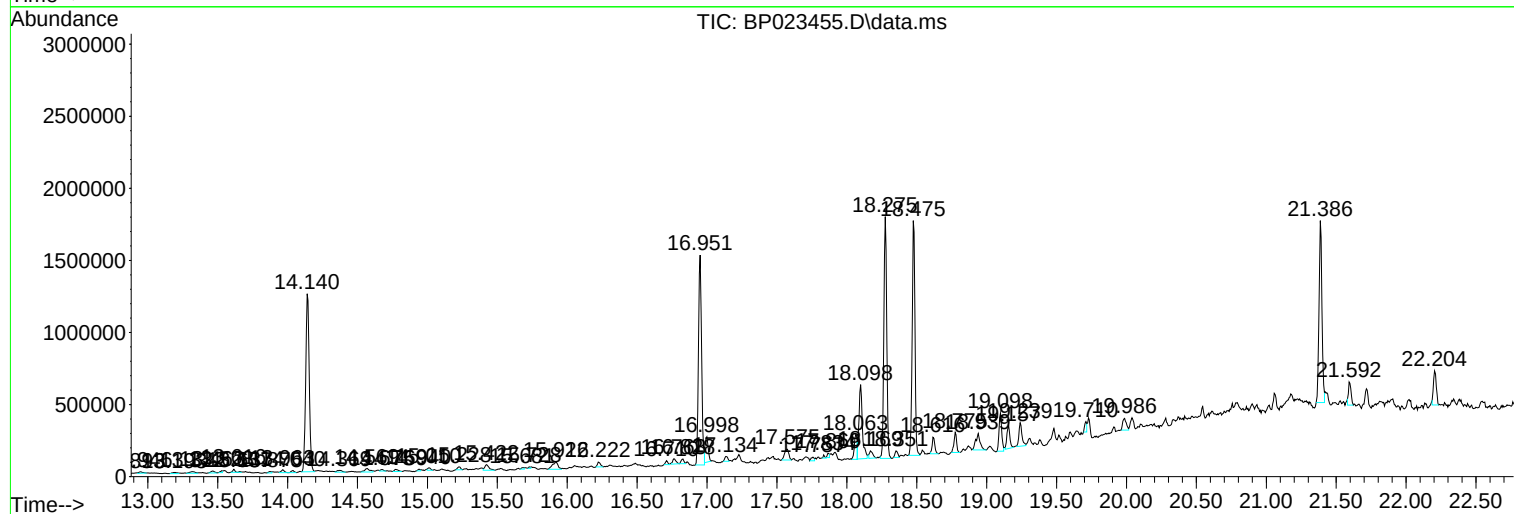
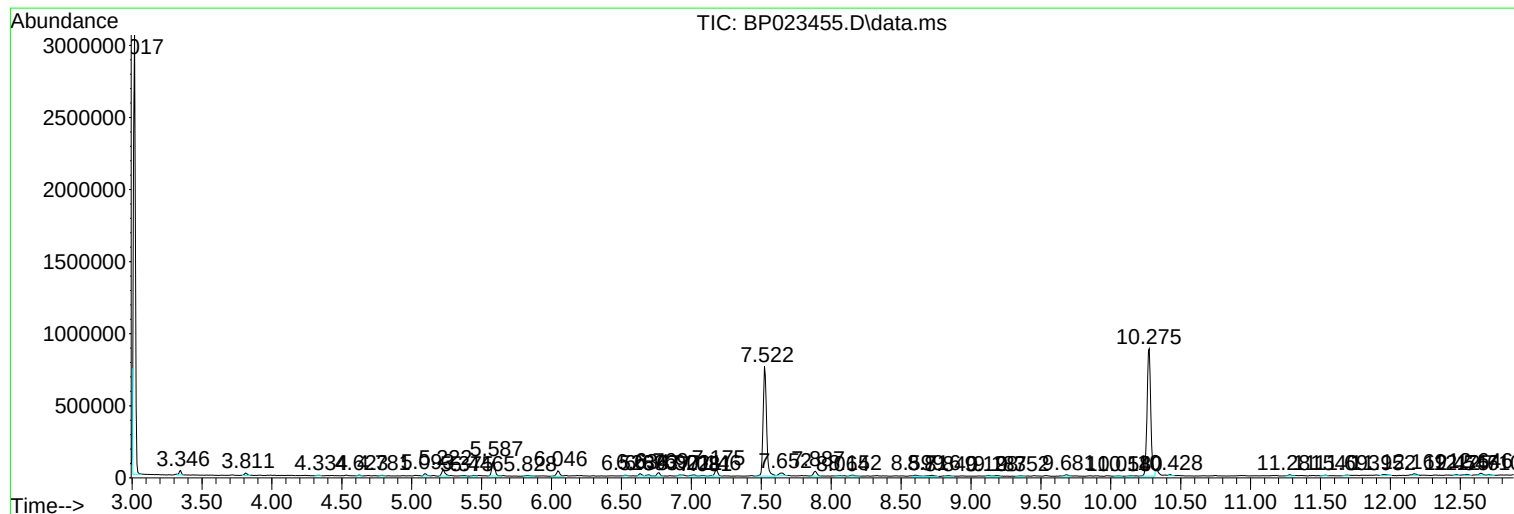
Sum of corrected areas: 25680989

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



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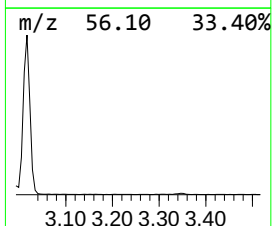
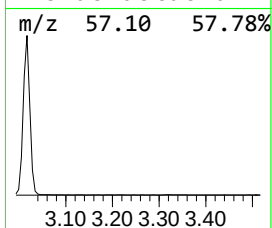
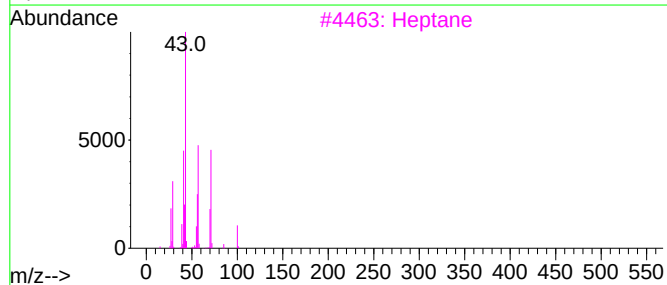
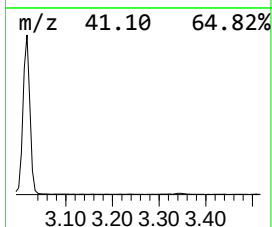
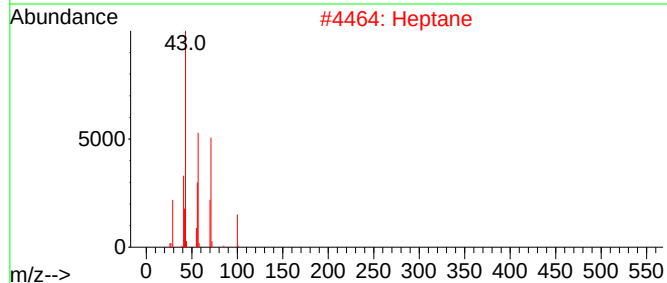
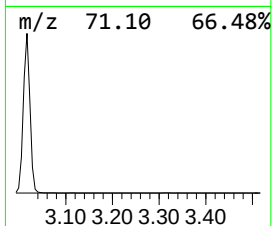
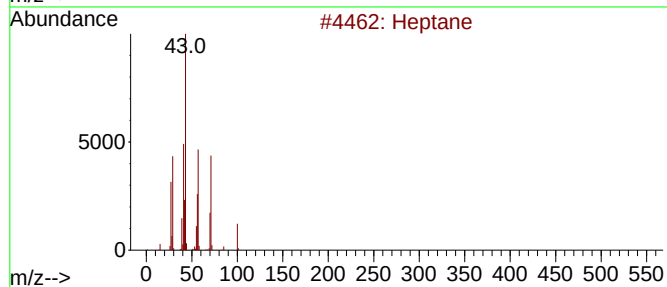
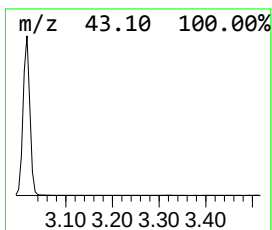
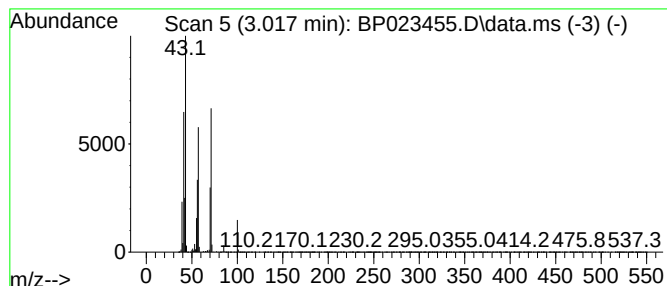
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	41.91 ng/ul	2719540	1,4-Dichlorobenzene-d4	7.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	91
2	Heptane			100	C7H16	000142-82-5	91
3	Heptane			100	C7H16	000142-82-5	87
4	Heptane			100	C7H16	000142-82-5	86
5	Hexane, 3-methyl-			100	C7H16	000589-34-4	50



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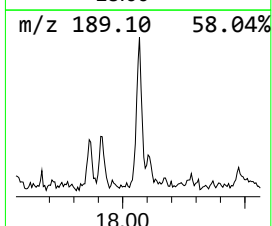
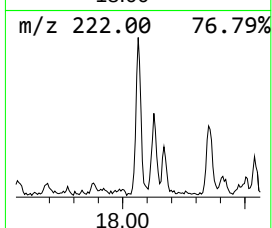
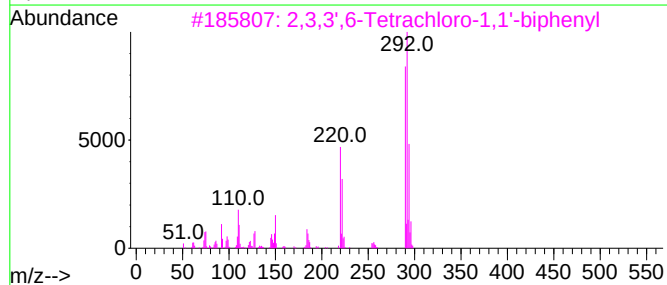
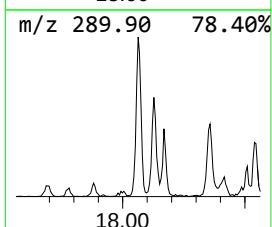
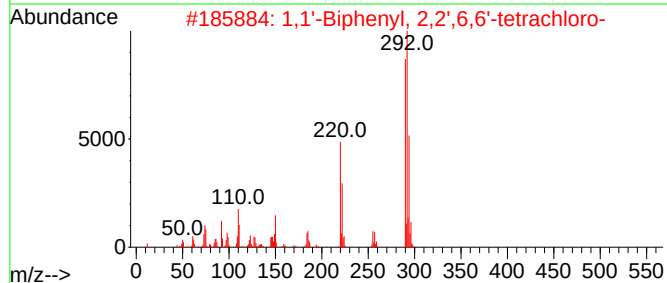
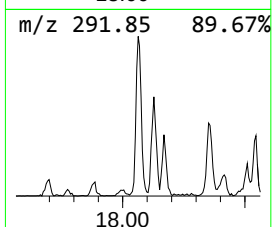
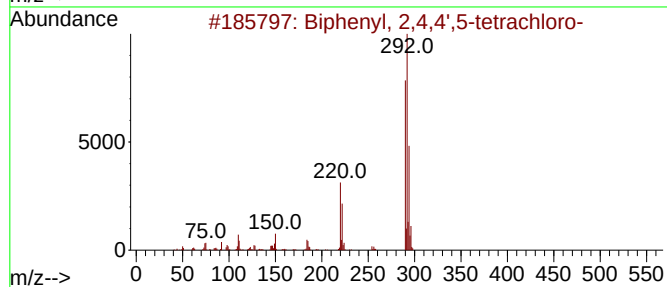
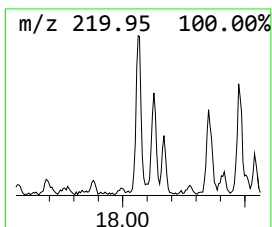
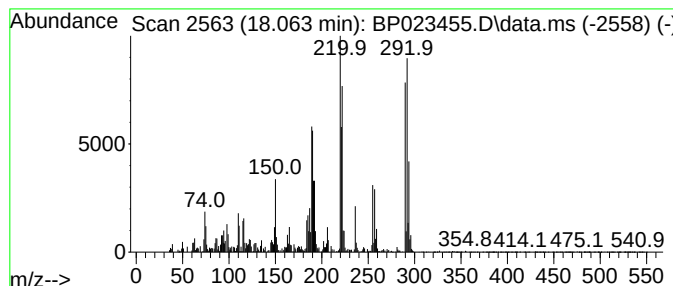
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Biphenyl, 2,4,4',5-tetrachl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.063	2.22 ng/ul	249640	Phenanthrene-d10	16.951	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
3	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	99



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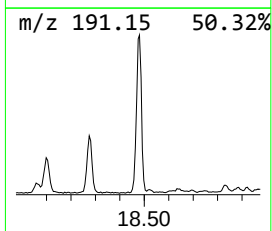
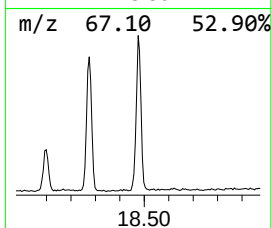
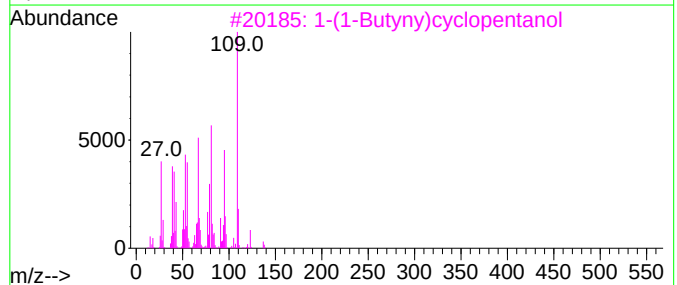
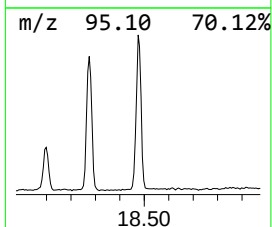
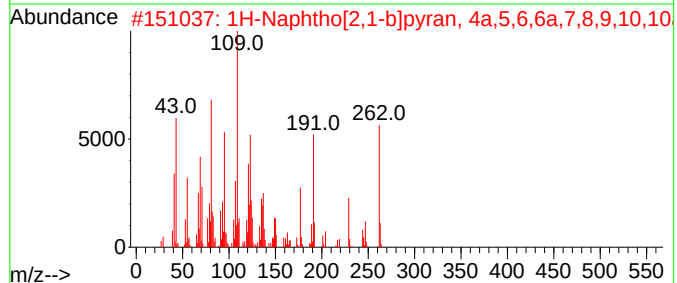
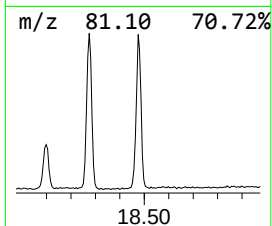
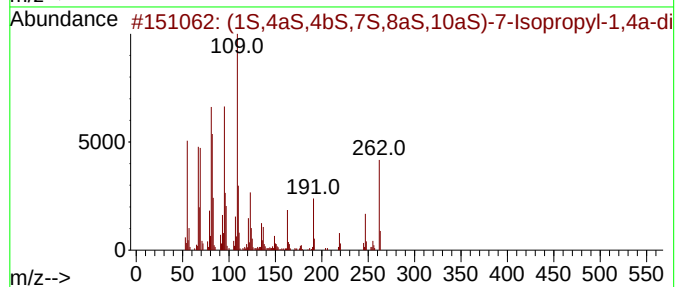
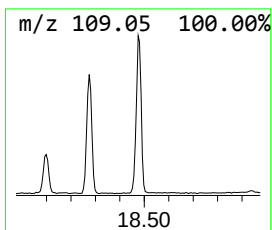
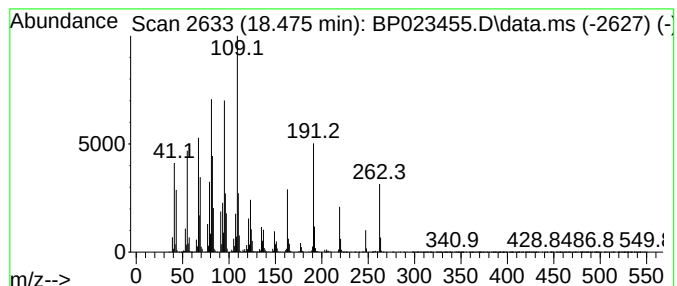
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (1S,4aS,4bS,7S,8aS,10aS)-7-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.475	20.52 ng/ul	2304410	Phenanthrene-d10	16.951

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S,4aS,4bS,7S,8aS,10aS)-7-Isopr...	262	C19H34	002221-95-6	86		
2	1H-Naphtho[2,1-b]pyran, 4a,5,6,6...	262	C18H30O	005153-92-4	41		
3	1-(1-Butyny)cyclopentanol	138	C9H14O	1000342-86-5	38		
4	(1S,4aS,4bS,7S,8aS,10aS)-7-Isopr...	262	C19H34	002221-95-6	38		
5	3-Tetradecyne	194	C14H26	060212-32-0	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
Acq On : 17 Dec 2024 05:57
Operator : RC/JU
Sample : P5234-15
Misc : GCMS CONFIRMATION
ALS Vial : 29 Sample Multiplier: 1

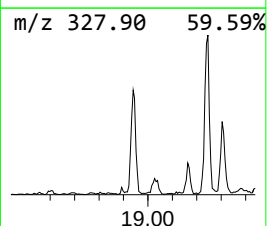
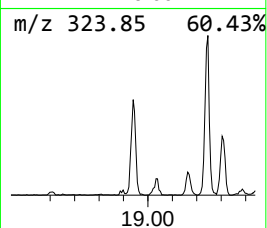
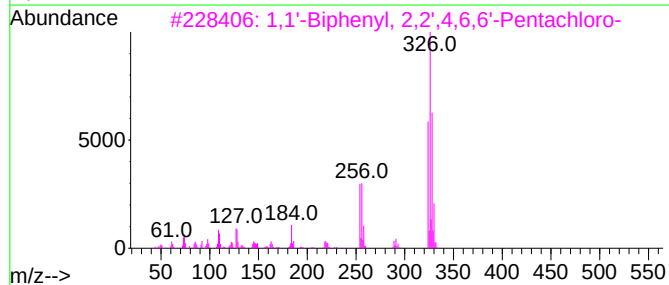
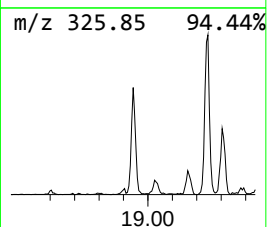
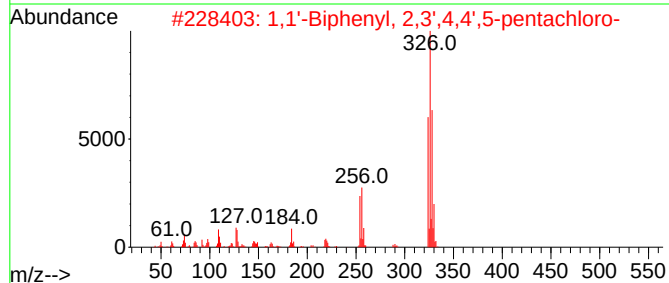
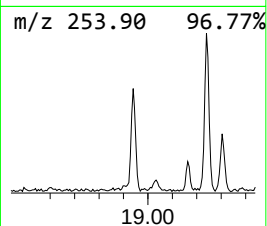
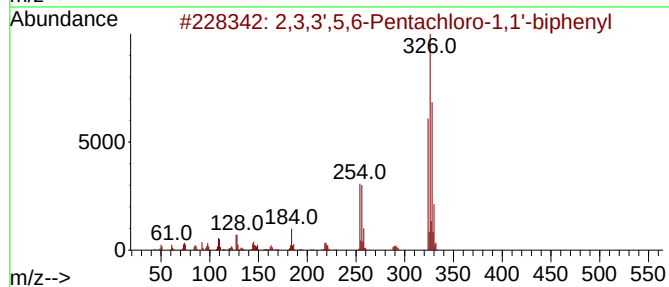
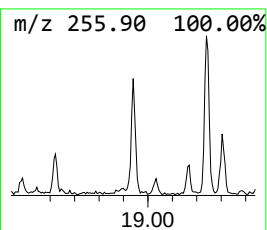
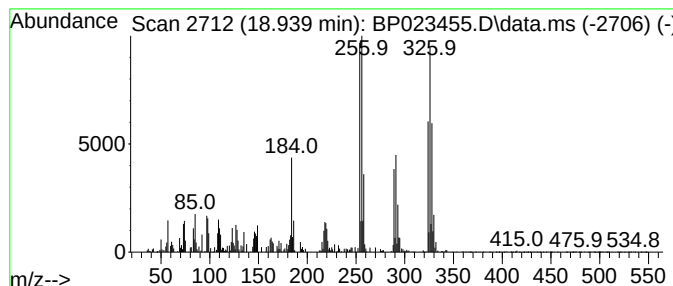
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,3,3',5,6-Pentachloro-1,1'... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.939	2.14 ng/ul	239961	Phenanthrene-d10	16.951	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99
2	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99
3	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99
4	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99
5	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
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Operator : RC/JU
Sample : P5234-15
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ALS Vial : 29 Sample Multiplier: 1

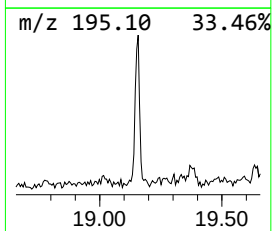
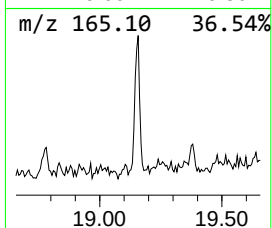
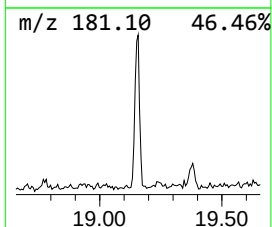
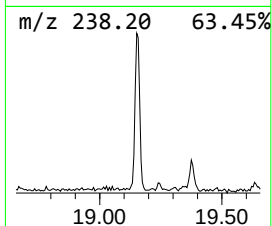
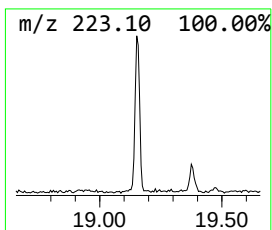
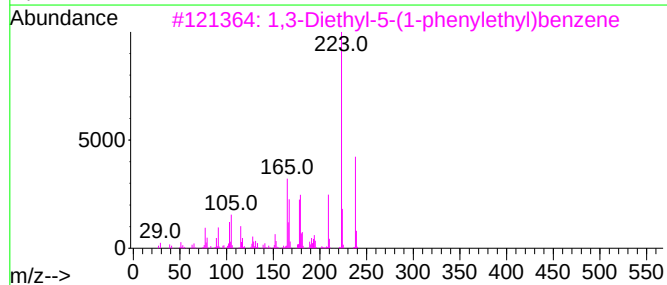
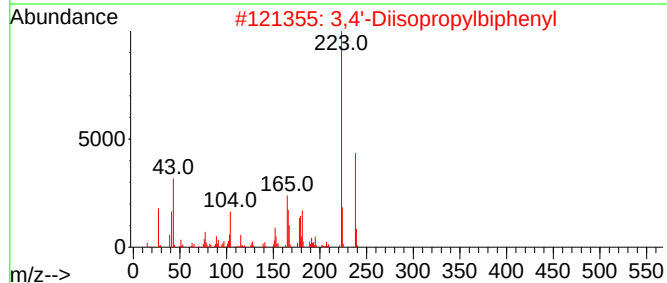
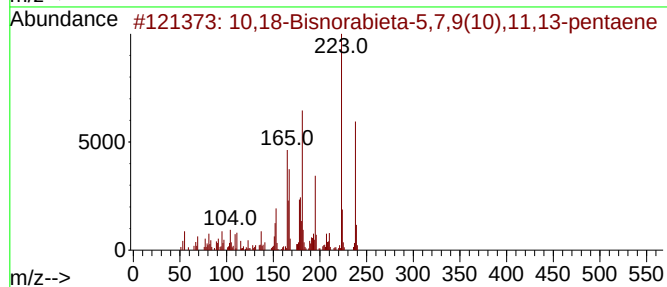
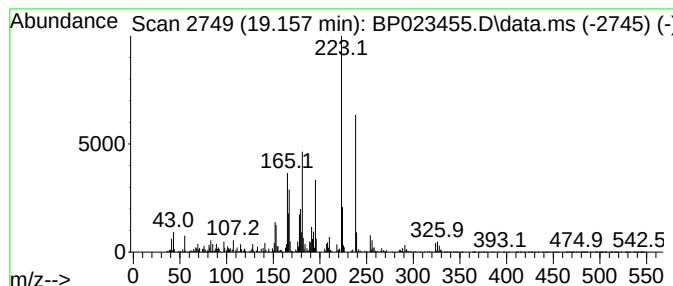
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.157	2.11 ng/ul	237019	Phenanthrene-d10	16.951	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	96
2	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	76
3	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	62
4	cyclopropanone, 2-(4-methoxyphen...	238	C16H14O2	1000401-16-2	53
5	Anthracene, 9,10-dimethoxy-	238	C16H14O2	002395-97-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
Acq On : 17 Dec 2024 05:57
Operator : RC/JU
Sample : P5234-15
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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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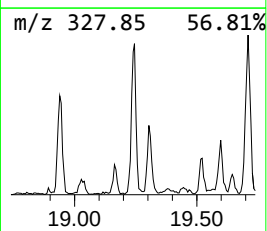
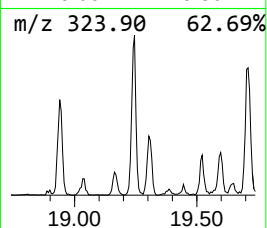
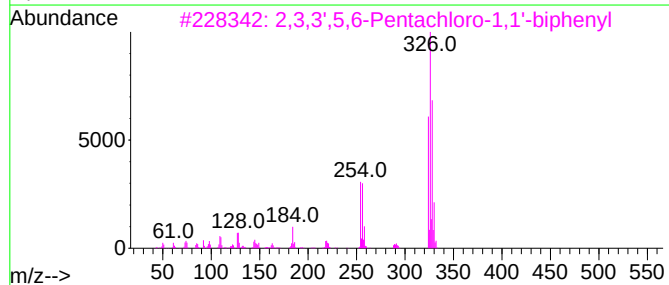
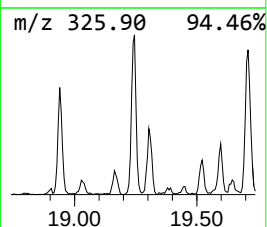
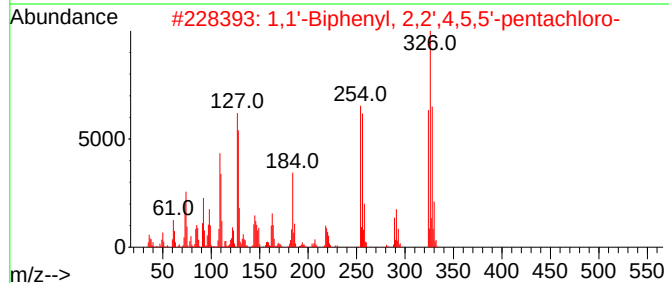
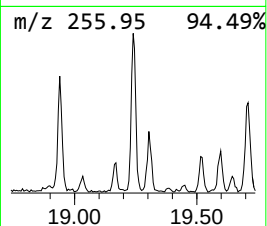
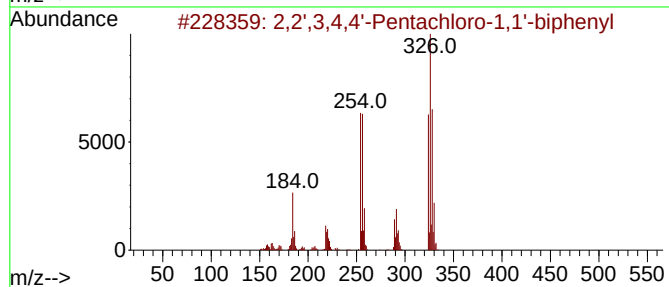
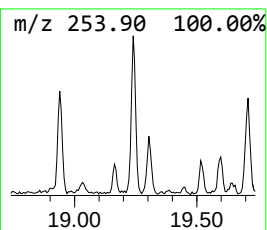
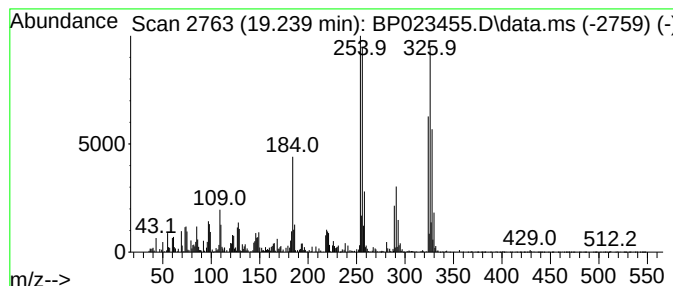
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,2',3,4,4'-Pentachloro-1,1... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	2.66 ng/ul	248328	Chrysene-d12	21.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	96		
2	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	95		
3	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	95		
4	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	95		
5	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
Acq On : 17 Dec 2024 05:57
Operator : RC/JU
Sample : P5234-15
Misc : GCMS CONFIRMATION
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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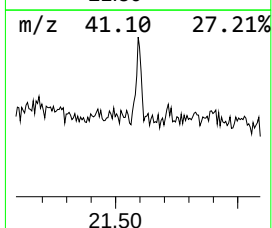
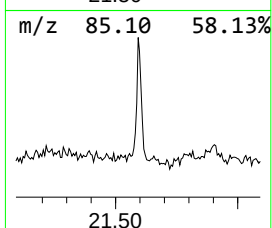
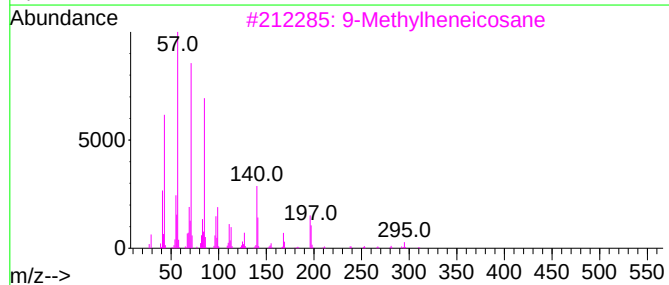
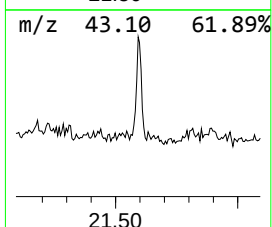
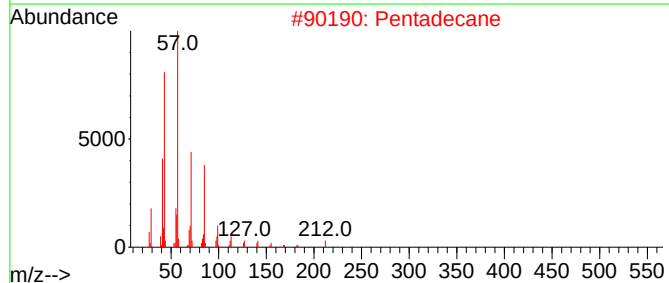
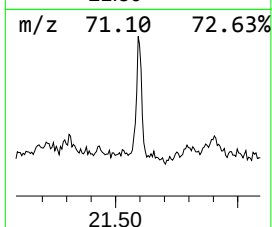
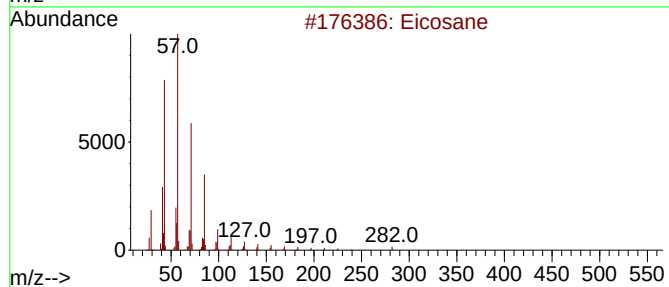
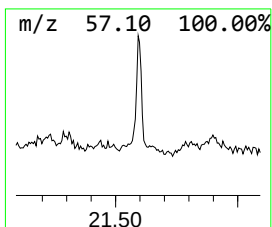
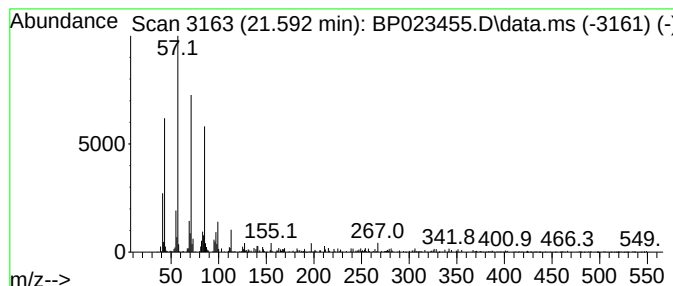
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.592	2.23 ng/ul	207883	Chrysene-d12	21.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	83
2		Pentadecane	212	C15H32	000629-62-9	83
3		9-Methylheneicosane	310	C22H46	070475-51-3	81
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	70
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
Acq On : 17 Dec 2024 05:57
Operator : RC/JU
Sample : P5234-15
Misc : GCMS CONFIRMATION
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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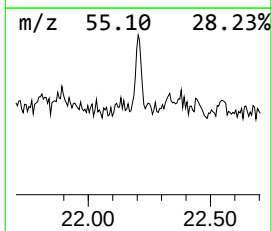
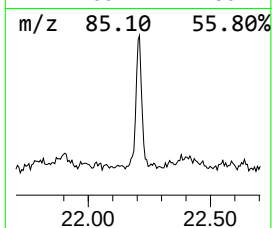
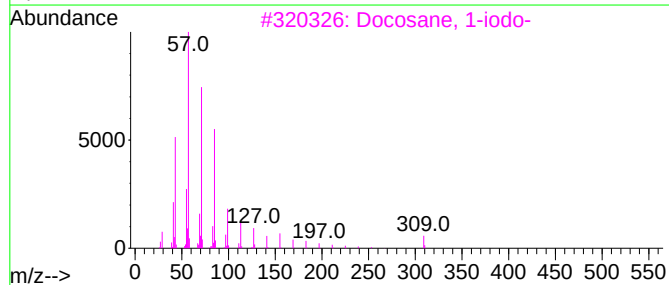
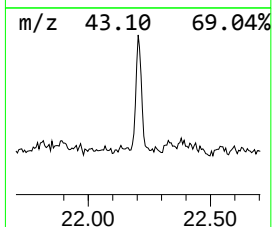
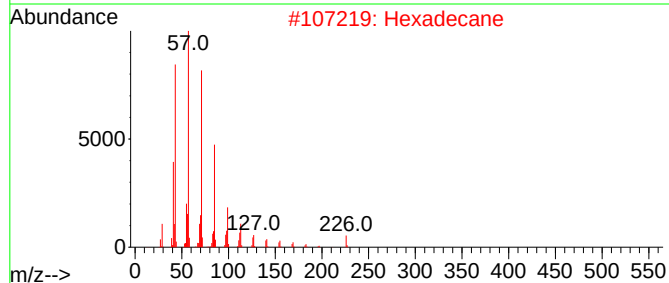
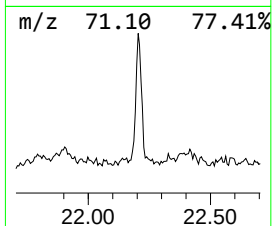
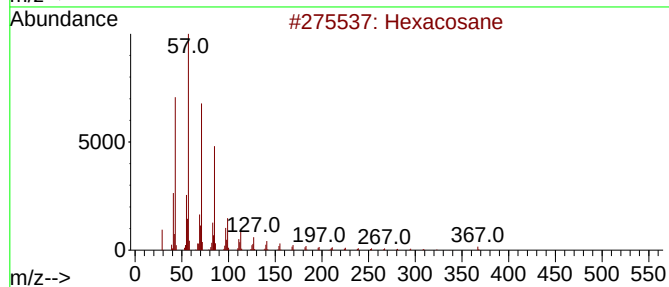
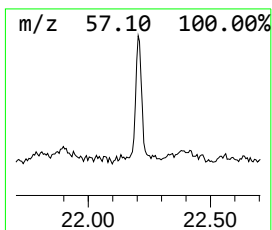
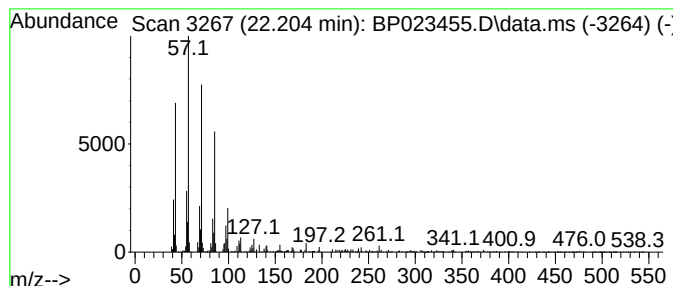
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.204	3.55 ng/ul	331167	Chrysene-d12	21.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	87
2			Hexadecane	226	C16H34	000544-76-3	87
3			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	86
4			Heptadecane	240	C17H36	000629-78-7	86
5			Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
Acq On : 17 Dec 2024 05:57
Operator : RC/JU
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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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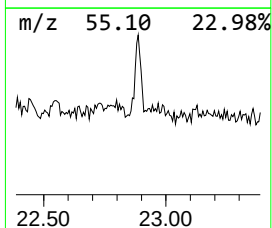
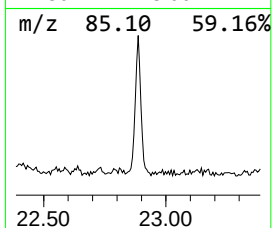
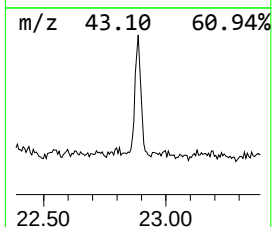
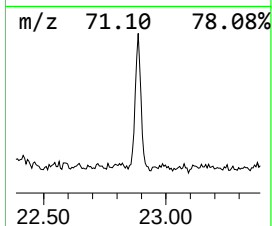
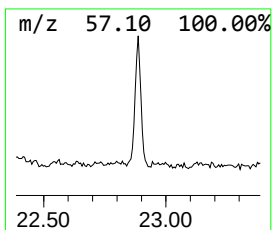
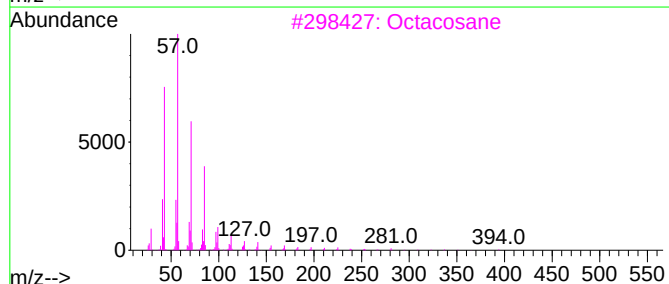
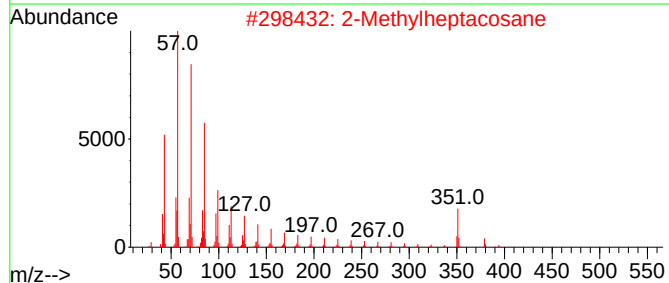
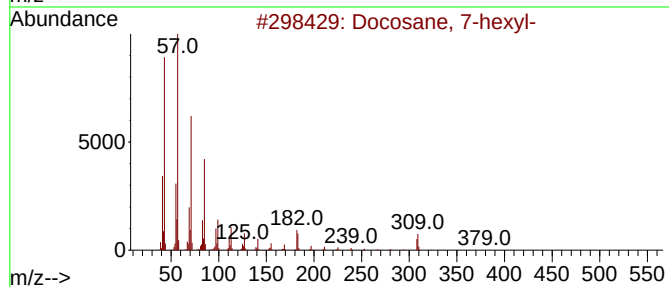
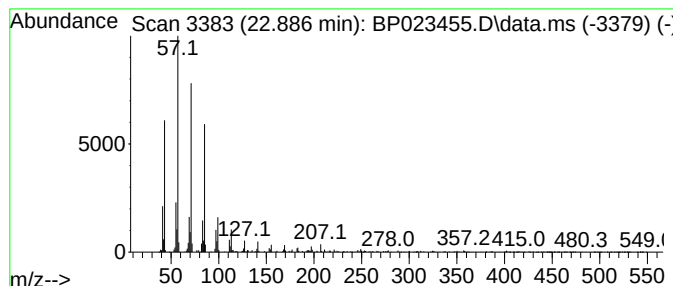
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.886	4.22 ng/ul	393650	Chrysene-d12	21.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
2		2-Methylheptacosane	394	C28H58	001561-00-8	91
3		Octacosane	394	C28H58	000630-02-4	91
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	90
5		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121624\
Data File : BP023455.D
Acq On : 17 Dec 2024 05:57
Operator : RC/JU
Sample : P5234-15
Misc : GCMS CONFIRMATION
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.017	41.9	ng/u1	2719540	1	7.522	1297930	20.0
Biphenyl, 2,4,4...	18.063	2.2	ng/u1	249640	4	16.951	2245960	20.0
(1S,4aS,4bS,7S,...	18.475	20.5	ng/u1	2304410	4	16.951	2245960	20.0
2,3,3',5,6-Pent...	18.939	2.1	ng/u1	239961	4	16.951	2245960	20.0
10,18-Bisnorabi...	19.157	2.1	ng/u1	237019	4	16.951	2245960	20.0
2,2',3,4,4'-Pen...	19.239	2.7	ng/u1	248328	5	21.386	1865090	20.0
(DEL) Alkane: S...	21.592	2.2	ng/u1	207883	5	21.386	1865090	20.0
(DEL) Alkane: S...	22.204	3.5	ng/u1	331167	5	21.386	1865090	20.0
(DEL) Alkane: S...	22.886	4.2	ng/u1	393650	5	21.386	1865090	20.0