

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
 Data File : BP006983.D
 Acq On : 15 Sep 2021 19:38
 Operator : CG/JU
 Sample : M3608-03DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBKJ9DL

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-BP090821.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006983.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.099	167	172	179	rVB	40249	58394	0.77%	0.065%
2	5.428	393	398	407	rVB	79752	118353	1.57%	0.132%
3	5.558	416	420	430	rVB	38342	82286	1.09%	0.092%
4	5.928	473	483	492	rBV	76489	128219	1.70%	0.143%
5	7.075	672	678	685	rVV2	47428	83885	1.11%	0.094%
6	7.252	702	708	715	rBV	133773	218295	2.89%	0.244%
7	7.446	735	741	750	rBV	121978	207662	2.75%	0.232%
8	7.569	756	762	769	rVB	86325	131740	1.74%	0.147%
9	7.640	769	774	781	rVB	124093	193862	2.57%	0.217%
10	7.922	814	822	830	rBV	1584235	2580851	34.18%	2.888%
11	8.034	837	841	849	rVV	57716	91355	1.21%	0.102%
12	8.293	877	885	893	rVV	238896	383051	5.07%	0.429%
13	8.458	902	913	926	rBV	626604	1042629	13.81%	1.167%
14	8.622	934	941	949	rVV2	107587	182039	2.41%	0.204%
15	9.081	1014	1019	1031	rVB2	140078	315190	4.17%	0.353%
16	9.275	1042	1052	1059	rVB	50532	87333	1.16%	0.098%
17	9.405	1059	1074	1080	rBV	98813	224387	2.97%	0.251%
18	9.799	1135	1141	1149	rBV	86111	161451	2.14%	0.181%
19	9.969	1165	1170	1181	rVB	49044	89537	1.19%	0.100%
20	10.122	1190	1196	1206	rBV4	89241	187033	2.48%	0.209%
21	10.346	1226	1234	1243	rVV2	243151	599281	7.94%	0.671%
22	10.722	1288	1298	1302	rVV	2347560	4073803	53.96%	4.558%
23	10.769	1302	1306	1314	rVV	2999610	4904130	64.95%	5.487%
24	10.887	1314	1326	1338	rVB2	368188	900560	11.93%	1.008%
25	11.081	1351	1359	1366	rBV3	32126	67298	0.89%	0.075%
26	12.105	1528	1533	1543	rVB3	32697	57740	0.76%	0.065%
27	12.275	1556	1562	1572	rVB	99211	166631	2.21%	0.186%
28	12.375	1572	1579	1586	rBV	528593	854516	11.32%	0.956%
29	12.499	1594	1600	1605	rVV3	60636	118935	1.58%	0.133%
30	12.593	1605	1616	1627	rVV	2044386	3419940	45.30%	3.827%
31	13.016	1683	1688	1690	rVV3	42460	70591	0.93%	0.079%
32	13.046	1690	1693	1698	rVV3	59885	103772	1.37%	0.116%
33	13.093	1698	1701	1714	rVB4	25464	54414	0.72%	0.061%
34	13.246	1721	1727	1736	rBV2	39077	89994	1.19%	0.101%
35	13.381	1744	1750	1758	rBV	784136	1195419	15.83%	1.338%
36	13.581	1778	1784	1796	rVB	158899	352687	4.67%	0.395%

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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-BP090821.M
 Title : SVOA CALIBRATION

37	13.722	1797	1808	1816	rBV2	314916	736893	9.76%	0.825%
38	13.869	1825	1833	1838	rBV2	408572	752977	9.97%	0.843%
39	13.922	1838	1842	1844	rBV	209904	276737	3.67%	0.310%
40	13.957	1844	1848	1857	rVB	501967	705833	9.35%	0.790%
41	14.104	1864	1873	1876	rBV2	155348	267483	3.54%	0.299%
42	14.134	1876	1878	1884	rVB	57292	74782	0.99%	0.084%
43	14.240	1890	1896	1899	rBV	472556	766605	10.15%	0.858%
44	14.546	1939	1948	1954	rVV	3981243	5964384	79.00%	6.674%
45	14.610	1954	1959	1965	rVV	3314226	4717663	62.48%	5.279%
46	14.675	1965	1970	1975	rVV	717393	1009751	13.37%	1.130%
47	14.740	1975	1981	1991	rVV5	78214	210510	2.79%	0.236%
48	14.816	1991	1994	1997	rVV	41999	57585	0.76%	0.064%
49	14.946	2009	2016	2024	rVB2	1820620	2926922	38.77%	3.275%
50	15.016	2024	2028	2034	rVB2	59357	86669	1.15%	0.097%
51	15.087	2034	2040	2044	rBV	422307	606982	8.04%	0.679%
52	15.122	2044	2046	2049	rVV	50650	63819	0.85%	0.071%
53	15.163	2049	2053	2058	rVB2	167818	253809	3.36%	0.284%
54	15.216	2058	2062	2066	rBV3	38622	53828	0.71%	0.060%
55	15.334	2075	2082	2085	rBV2	35976	75485	1.00%	0.084%
56	15.475	2100	2106	2111	rBV	159369	293283	3.88%	0.328%
57	15.534	2111	2116	2121	rVV	703751	983765	13.03%	1.101%
58	15.593	2121	2126	2131	rVB	1535545	2148377	28.45%	2.404%
59	15.657	2131	2137	2143	rBV	768333	1128840	14.95%	1.263%
60	15.716	2143	2147	2150	rVV2	85879	104335	1.38%	0.117%
61	15.751	2150	2153	2158	rVB2	60202	82653	1.09%	0.092%
62	15.810	2158	2163	2167	rBV2	119746	182741	2.42%	0.204%
63	15.887	2172	2176	2185	rVB2	92005	154189	2.04%	0.173%
64	16.028	2196	2200	2208	rVV	109839	220892	2.93%	0.247%
65	16.098	2208	2212	2222	rVB	130970	211122	2.80%	0.236%
66	16.263	2234	2240	2250	rVB3	223024	467026	6.19%	0.523%
67	16.387	2257	2261	2266	rBV	51573	67439	0.89%	0.075%
68	16.545	2284	2288	2291	rVV2	37749	59483	0.79%	0.067%
69	16.598	2294	2297	2302	rVV3	41658	76203	1.01%	0.085%
70	16.645	2302	2305	2312	rVB5	38995	65777	0.87%	0.074%
71	16.940	2346	2355	2373	rBV	2776518	3796898	50.29%	4.248%
72	17.069	2373	2377	2380	rBV2	59615	88063	1.17%	0.099%
73	17.110	2380	2384	2394	rVB	192993	277244	3.67%	0.310%
74	17.292	2409	2415	2419	rBV	4813311	6975681	92.39%	7.805%
75	17.334	2419	2422	2427	rVB	1972099	2642686	35.00%	2.957%
76	17.392	2427	2432	2437	rBV	1045265	1455702	19.28%	1.629%

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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-BP090821.M
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77	17.487	2444	2448	2458	rVB5	48615	115530	1.53%	0.129%
78	17.692	2473	2483	2493	rBV	3041998	4483853	59.39%	5.017%
79	17.787	2495	2499	2505	rVV3	59234	125166	1.66%	0.140%
80	17.845	2505	2509	2515	rVV	106132	168211	2.23%	0.188%
81	17.916	2517	2521	2524	rVV5	30247	58953	0.78%	0.066%
82	17.945	2524	2526	2530	rVV3	41827	52180	0.69%	0.058%
83	17.992	2530	2534	2537	rVB2	41039	53271	0.71%	0.060%
84	18.116	2552	2555	2559	rBV	45214	52281	0.69%	0.058%
85	18.157	2559	2562	2564	rVV	48545	61304	0.81%	0.069%
86	18.198	2564	2569	2577	rVV2	308523	489168	6.48%	0.547%
87	18.269	2578	2581	2588	rVB2	56434	75411	1.00%	0.084%
88	18.345	2588	2594	2598	rBV2	89573	157839	2.09%	0.177%
89	18.445	2609	2611	2615	rVV	72652	94575	1.25%	0.106%
90	18.487	2615	2618	2622	rVV2	93412	128518	1.70%	0.144%
91	18.581	2630	2634	2639	rVB2	104253	147884	1.96%	0.165%
92	18.634	2639	2643	2647	rVB2	34869	49386	0.65%	0.055%
93	18.681	2647	2651	2655	rBV3	48245	65835	0.87%	0.074%
94	19.298	2751	2756	2763	rVB	192007	258008	3.42%	0.289%
95	19.622	2805	2811	2823	rBV	1031807	1583899	20.98%	1.772%
96	20.016	2871	2878	2883	rBV2	138247	253824	3.36%	0.284%
97	20.075	2883	2888	2894	rBV	208940	333710	4.42%	0.373%
98	21.375	3103	3109	3118	rBV	5766364	7550101	100.00%	8.448%
99	23.633	3488	3493	3499	rVB2	525495	959985	12.71%	1.074%
100	23.792	3513	3520	3532	rVB2	3691003	7365675	97.56%	8.242%

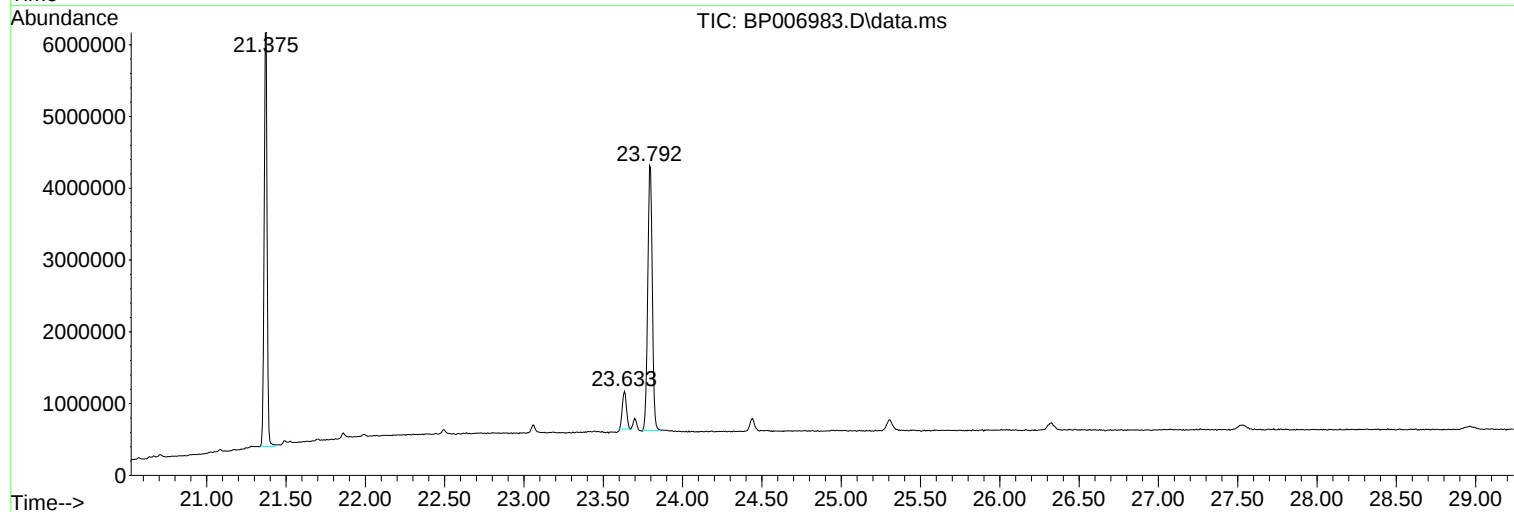
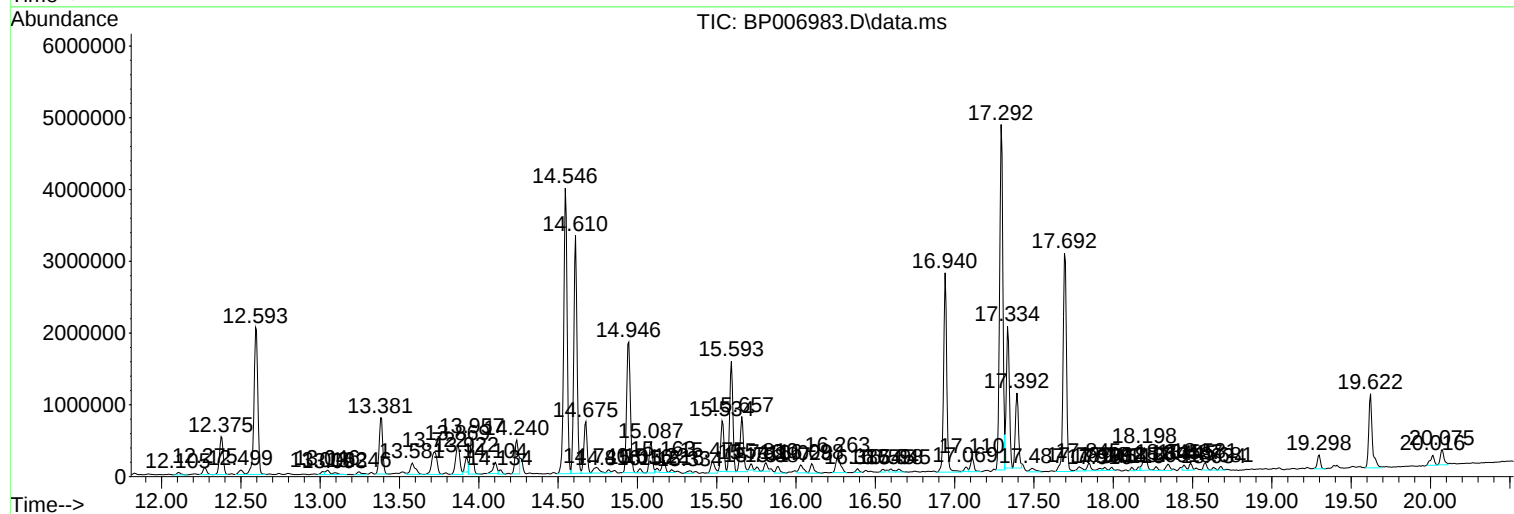
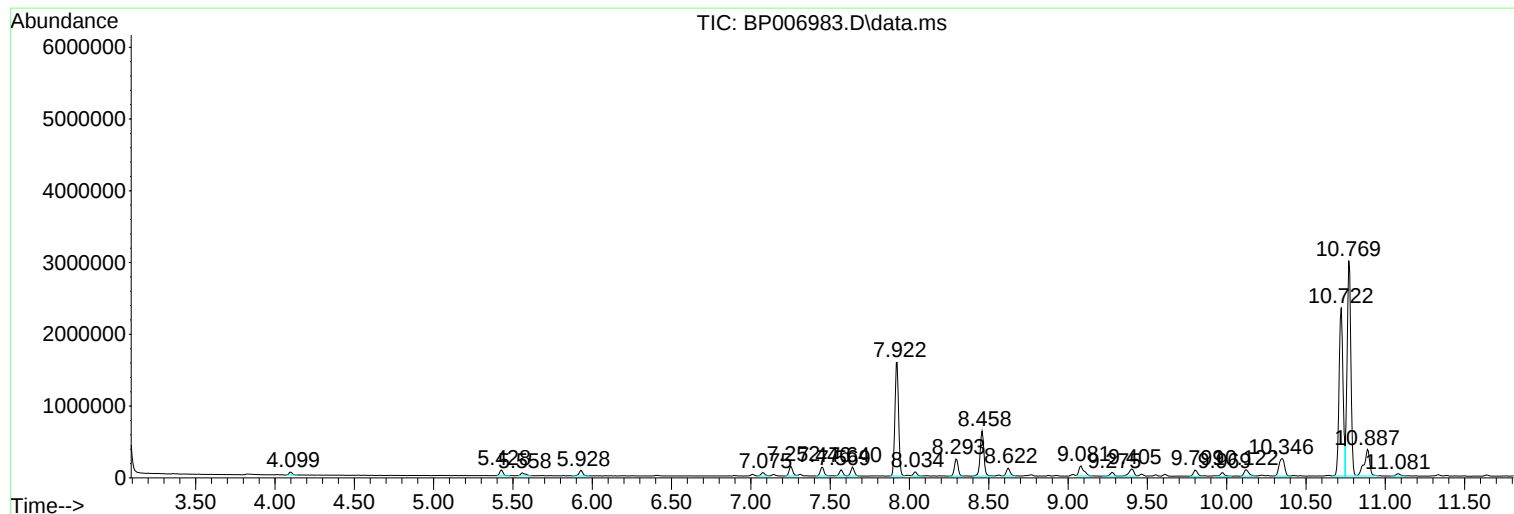
Sum of corrected areas: 89372946

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION

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



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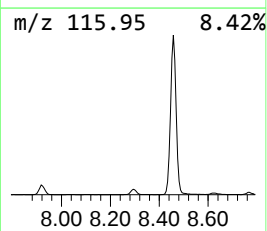
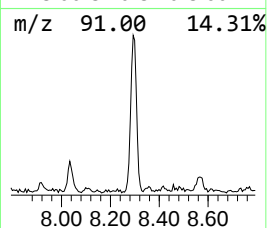
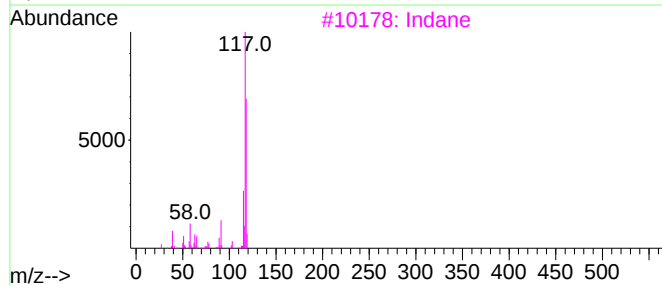
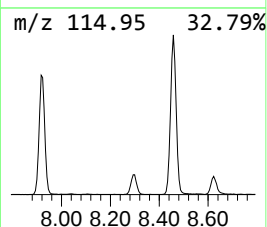
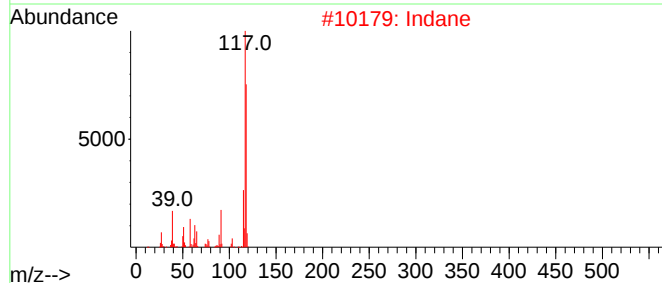
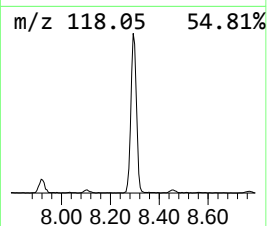
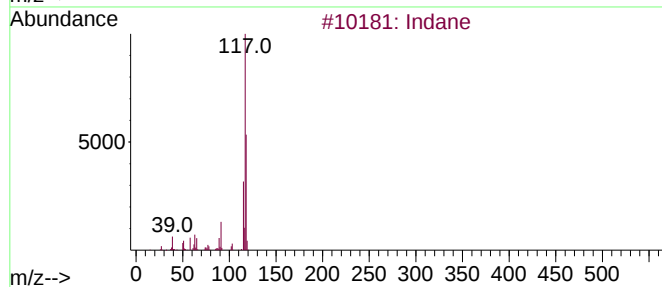
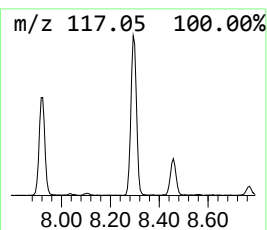
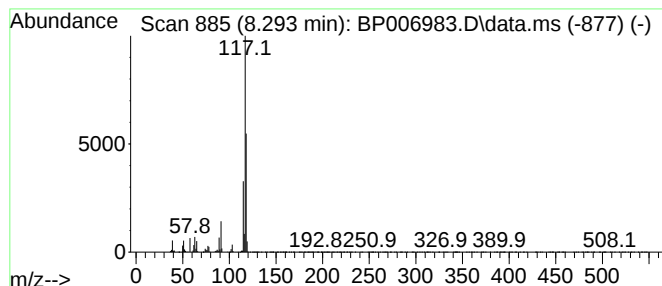
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.293	2.97 ng/ul	383051	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	95
2	Indane			118	C9H10	000496-11-7	91
3	Indane			118	C9H10	000496-11-7	87
4	Benzene, cyclopropyl-			118	C9H10	000873-49-4	87
5	Deltacyclene			118	C9H10	007785-10-6	72



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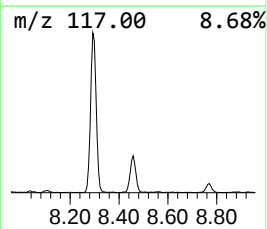
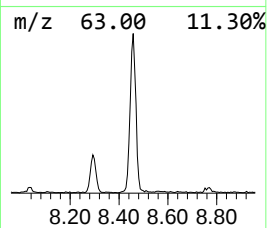
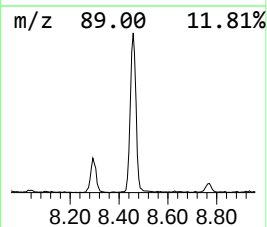
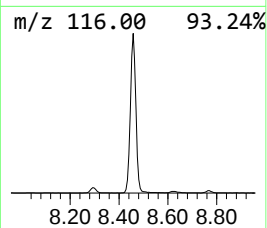
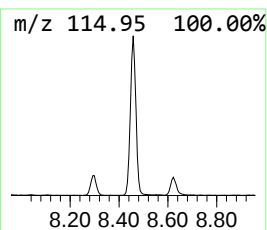
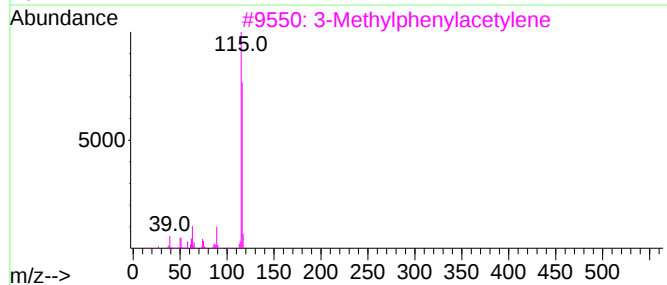
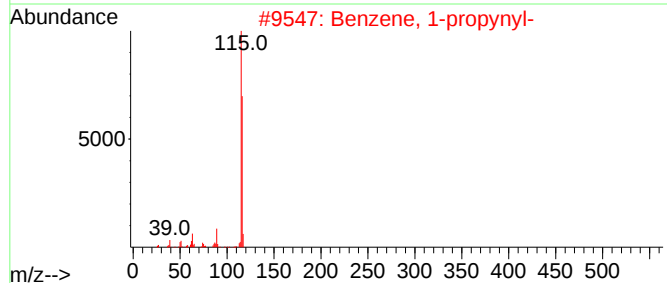
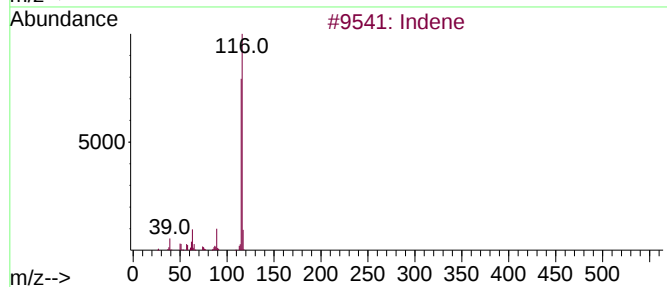
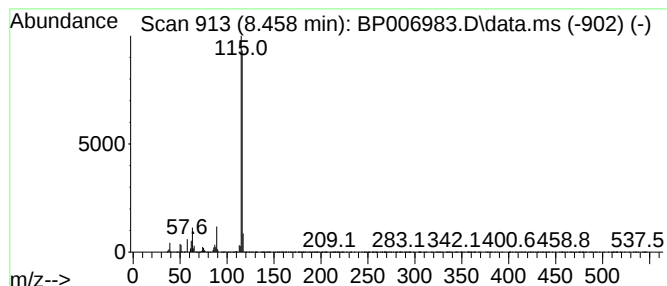
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.458	8.08 ng/ul	1042630	1,4-Dichlorobenzene-d4	7.922

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
3	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91



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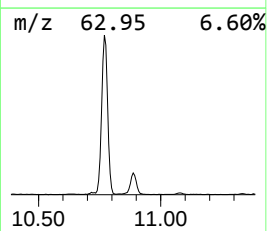
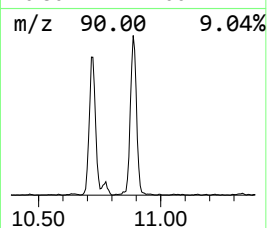
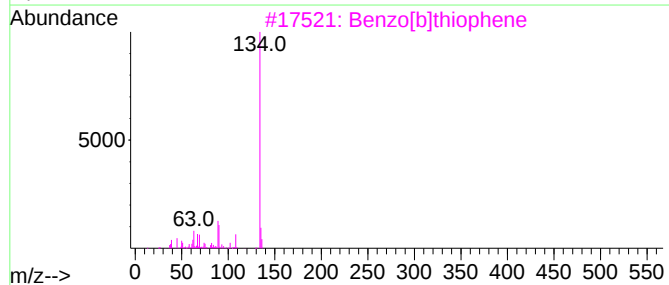
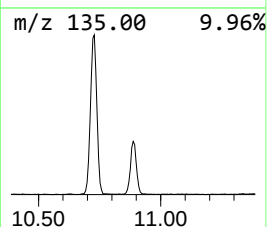
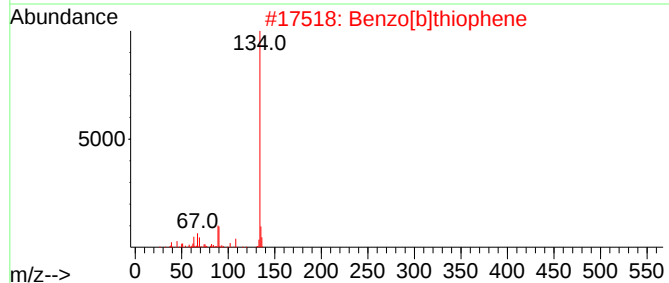
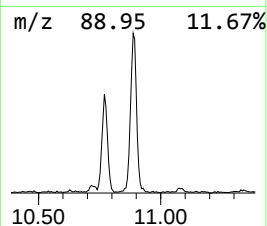
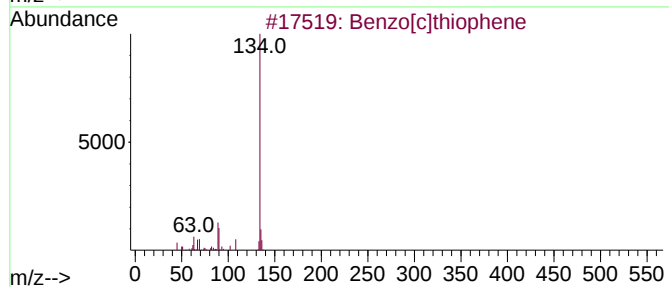
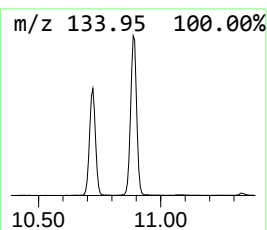
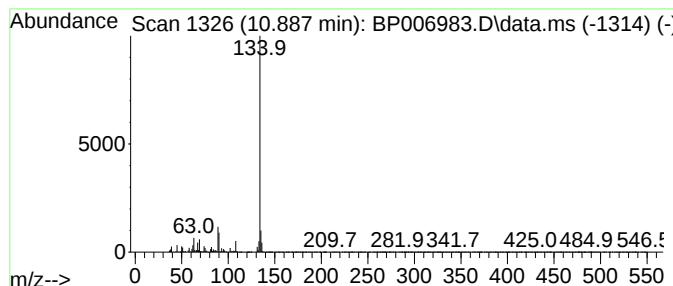
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzo[c]thiophene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.887	4.42 ng/ul	900560	Naphthalene-d8	10.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2			Benzo[b]thiophene	134	C8H6S	000095-15-8	95
3			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
4			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
5			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
Data File : BP006983.D
Acq On : 15 Sep 2021 19:38
Operator : CG/JU
Sample : M3608-03DL 10X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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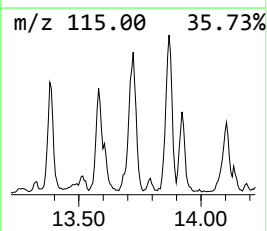
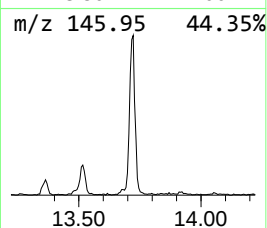
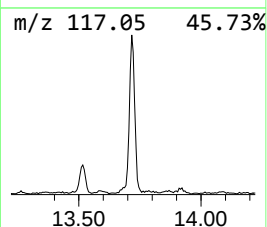
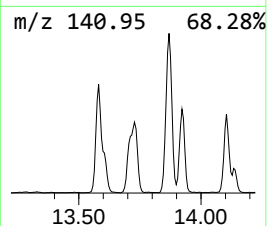
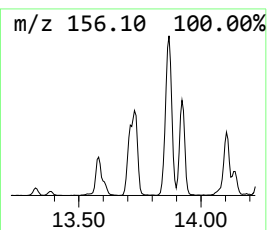
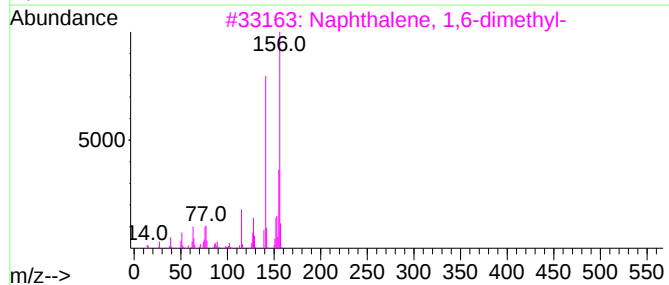
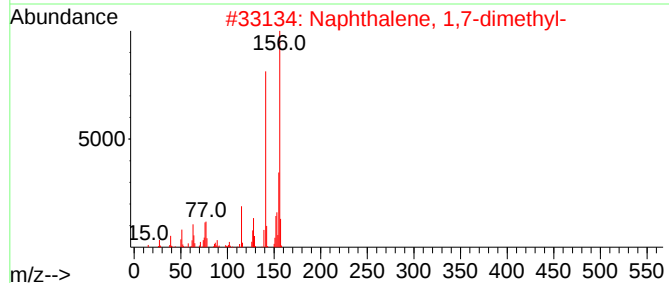
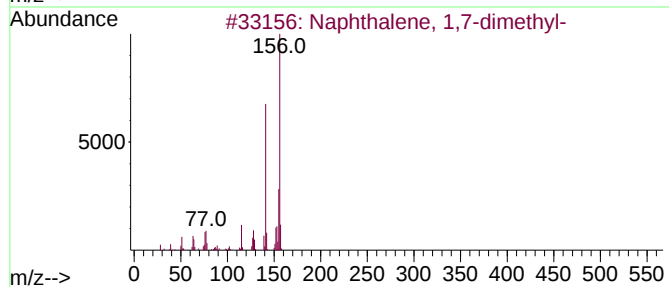
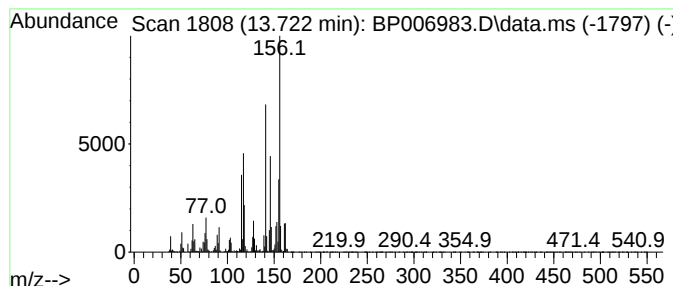
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 1,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.722	2.47 ng/ul	736893	Acenaphthene-d10	14.546

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
Data File : BP006983.D
Acq On : 15 Sep 2021 19:38
Operator : CG/JU
Sample : M3608-03DL 10X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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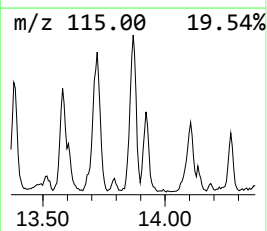
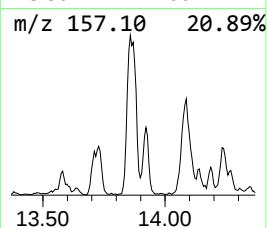
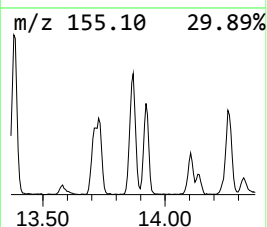
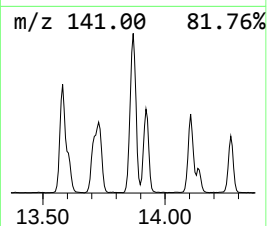
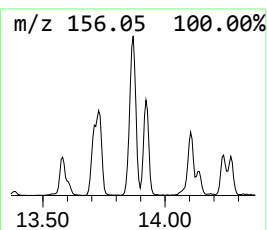
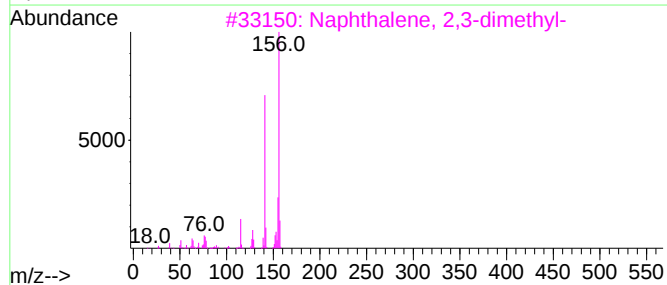
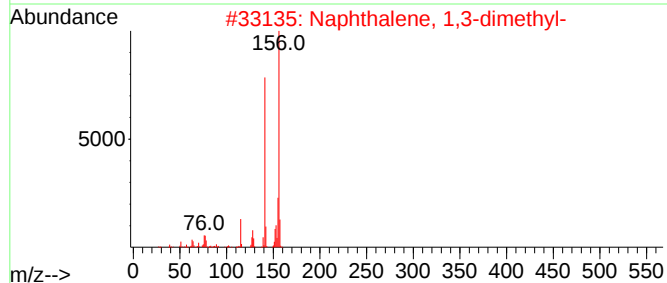
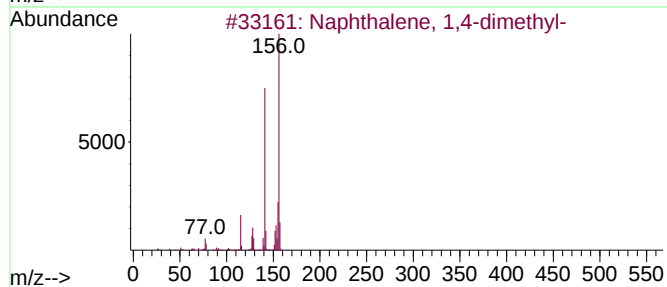
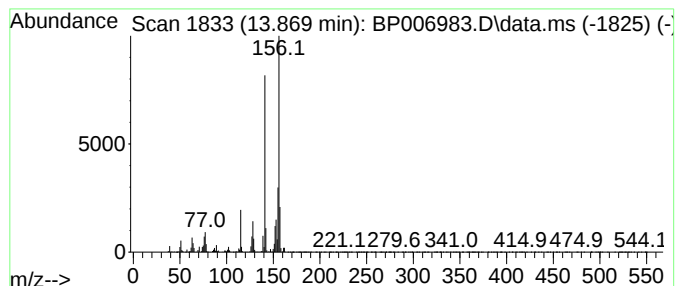
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.869	2.52 ng/ul	752977	Acenaphthene-d10	14.546

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
Data File : BP006983.D
Acq On : 15 Sep 2021 19:38
Operator : CG/JU
Sample : M3608-03DL 10X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBKJ9DL

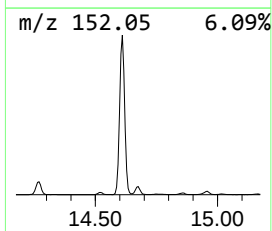
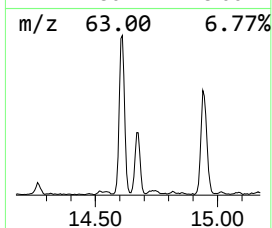
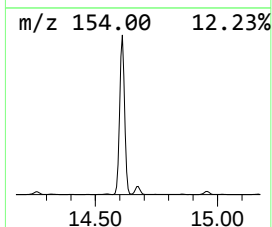
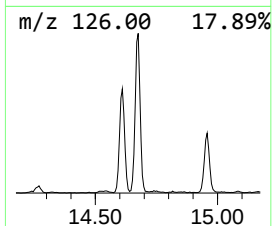
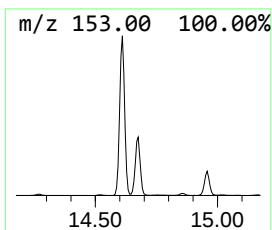
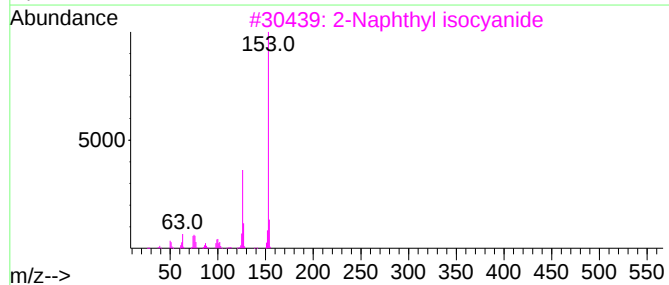
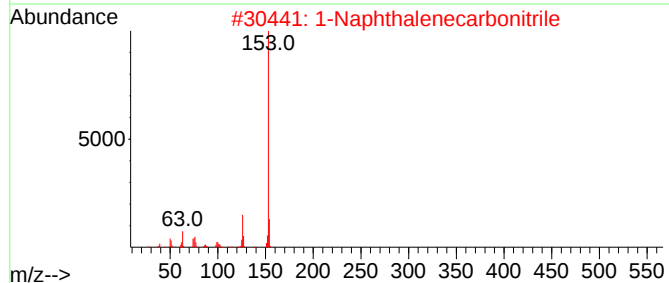
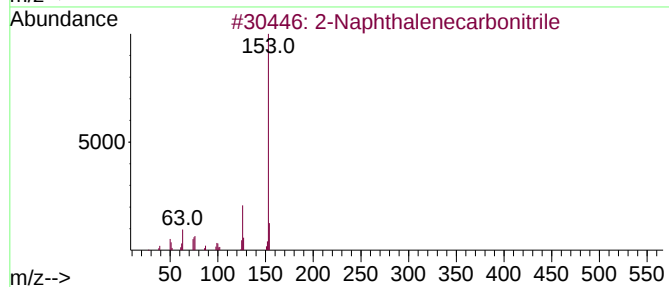
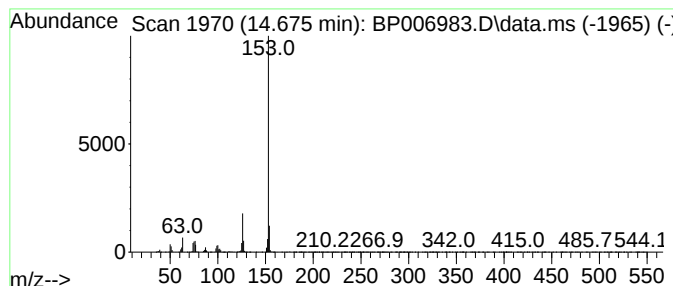
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Naphthalenecarbonitrile Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.675	3.39 ng/ul	1009750	Acenaphthene-d10	14.546

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile			153	C11H7N	000613-46-7	97
2	1-Naphthalenecarbonitrile			153	C11H7N	000086-53-3	97
3	2-Naphthyl isocyanide			153	C11H7N	010124-78-4	95
4	1-Naphthalenecarbonitrile			153	C11H7N	000086-53-3	94
5	2-Naphthalenecarbonitrile			153	C11H7N	000613-46-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
Data File : BP006983.D
Acq On : 15 Sep 2021 19:38
Operator : CG/JU
Sample : M3608-03DL 10X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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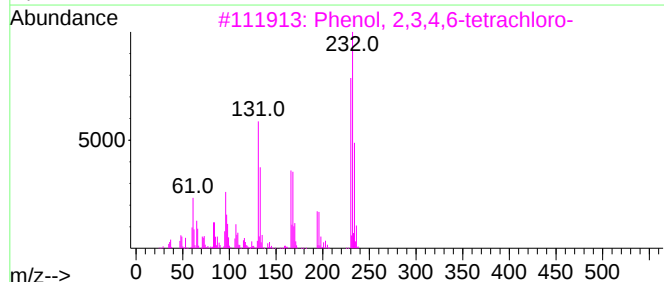
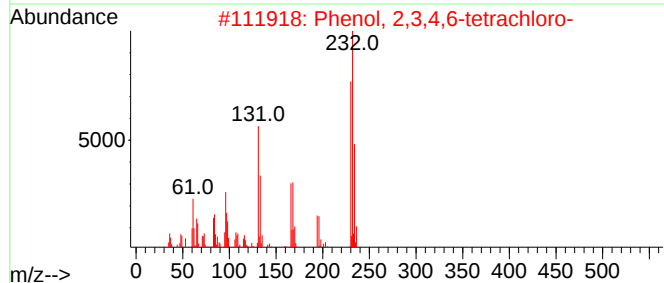
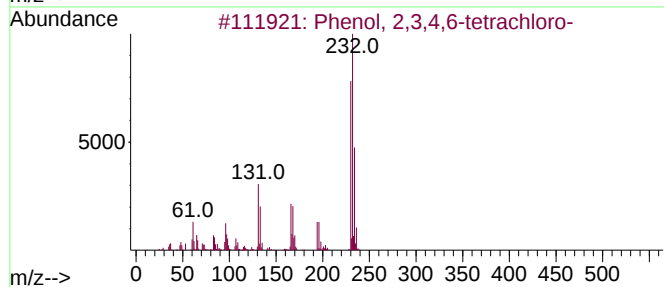
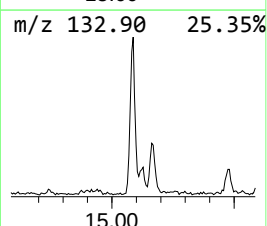
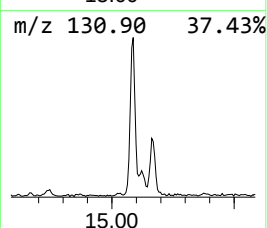
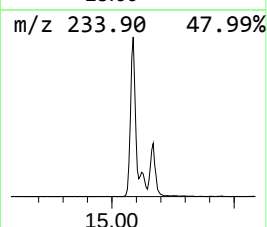
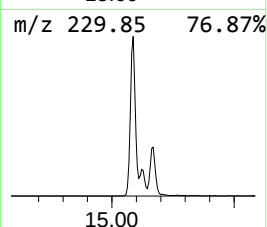
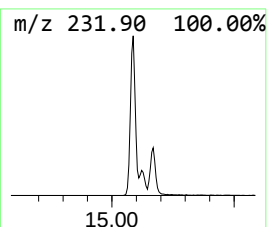
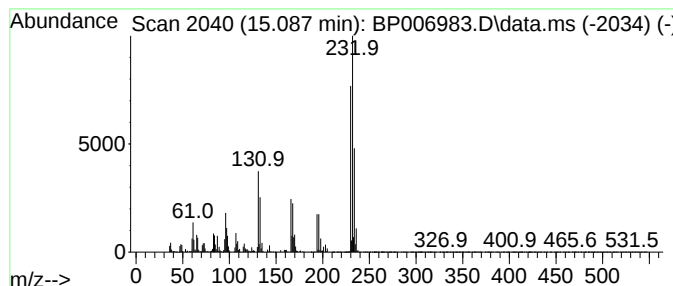
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.087	2.04 ng/ul	606982	Acenaphthene-d10	14.546

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
3			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
4			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
5			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091521\
Data File : BP006983.D
Acq On : 15 Sep 2021 19:38
Operator : CG/JU
Sample : M3608-03DL 10X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBKJ9DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.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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Indene	8.458	8.1	ng/ul	1042630	1	7.922	2580850	20.0
Benzo[c]thiophene	10.887	4.4	ng/ul	900560	2	10.722	4073800	20.0
Naphthalene, 1,...	13.722	2.5	ng/ul	736893	3	14.546	5964380	20.0
Naphthalene, 1,...	13.869	2.5	ng/ul	752977	3	14.546	5964380	20.0
2-Naphthaleneca...	14.675	3.4	ng/ul	1009750	3	14.546	5964380	20.0
Phenol, 2,3,5,6...	15.087	2.0	ng/ul	606982	3	14.546	5964380	20.0