

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
 Data File : BP021290.D
 Acq On : 01 Aug 2024 12:40
 Operator : CG/JU
 Sample : P3347-14DL 10X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AY8DL

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021290.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.181	366	373	383	rBV	21540	38912	0.81%	0.106%
2	5.311	388	395	403	rBV	13729	34878	0.72%	0.095%
3	5.669	450	456	463	rBV	13848	25754	0.53%	0.070%
4	6.893	659	664	672	rVV6	12226	34188	0.71%	0.093%
5	6.993	676	681	692	rVB3	25671	58461	1.21%	0.160%
6	7.199	706	716	724	rBV	35784	87768	1.82%	0.240%
7	7.275	724	729	737	rVV2	16488	33971	0.70%	0.093%
8	7.357	737	743	751	rVB2	16244	28631	0.59%	0.078%
9	7.628	782	789	802	rBV	492441	897986	18.60%	2.454%
10	7.981	843	849	856	rBV	146971	247485	5.13%	0.676%
11	8.157	865	879	890	rBV2	77390	167009	3.46%	0.456%
12	8.381	909	917	928	rBV2	36006	91508	1.90%	0.250%
13	8.622	951	958	965	rBV6	11045	24382	0.51%	0.067%
14	8.804	979	989	1001	rBV3	32379	115096	2.38%	0.315%
15	8.957	1007	1015	1021	rVB2	19900	33986	0.70%	0.093%
16	9.069	1024	1034	1036	rBV6	25218	60593	1.26%	0.166%
17	9.510	1101	1109	1116	rBV2	26904	61913	1.28%	0.169%
18	9.646	1122	1132	1146	rVB3	34905	110277	2.28%	0.301%
19	9.793	1146	1157	1168	rBV8	27028	97620	2.02%	0.267%
20	9.928	1168	1180	1190	rBV6	24082	74889	1.55%	0.205%
21	10.075	1199	1205	1215	rBV3	52006	118013	2.44%	0.322%
22	10.322	1241	1247	1251	rBV4	18394	38895	0.81%	0.106%
23	10.393	1252	1259	1262	rVV	800050	1304671	27.03%	3.565%
24	10.440	1262	1267	1280	rVB	3205724	4827285	100.00%	13.191%
25	10.557	1280	1287	1297	rBV2	191672	422484	8.75%	1.155%
26	10.798	1317	1328	1335	rBV5	13627	37370	0.77%	0.102%
27	10.957	1347	1355	1365	rBV2	14730	47296	0.98%	0.129%
28	11.163	1382	1390	1398	rBV6	13236	36349	0.75%	0.099%
29	11.457	1427	1440	1441	rBV5	9808	27334	0.57%	0.075%
30	11.504	1442	1448	1456	rVV5	20441	48449	1.00%	0.132%
31	11.769	1487	1493	1499	rBV6	12204	29214	0.61%	0.080%
32	11.945	1518	1523	1531	rBV	33043	63286	1.31%	0.173%
33	12.051	1534	1541	1550	rVV	303428	487935	10.11%	1.333%
34	12.210	1560	1568	1572	rVV	51918	111492	2.31%	0.305%
35	12.281	1572	1580	1589	rVB	344975	619429	12.83%	1.693%
36	12.775	1661	1664	1671	rVB4	14766	24170	0.50%	0.066%

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Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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37	13.075	1709	1715	1728	rVV2	96720	168244	3.49%	0.460%
38	13.304	1742	1754	1760	rVV4	37914	115239	2.39%	0.315%
39	13.416	1767	1773	1781	rBV3	86527	209506	4.34%	0.573%
40	13.569	1790	1799	1805	rBV2	126317	296288	6.14%	0.810%
41	13.628	1805	1809	1812	rVV	52644	100799	2.09%	0.275%
42	13.675	1812	1817	1825	rVB2	107713	226928	4.70%	0.620%
43	13.810	1831	1840	1844	rBV2	135581	272221	5.64%	0.744%
44	13.840	1844	1845	1852	rVB	53871	58890	1.22%	0.161%
45	13.910	1852	1857	1860	rBV2	19469	33142	0.69%	0.091%
46	13.957	1860	1865	1867	rBV	127362	202544	4.20%	0.553%
47	14.257	1907	1916	1922	rBV	1337927	1931539	40.01%	5.278%
48	14.328	1922	1928	1937	rVB	1858543	2924422	60.58%	7.991%
49	14.428	1938	1945	1950	rBV	100963	161098	3.34%	0.440%
50	14.569	1963	1969	1975	rVB3	82069	145305	3.01%	0.397%
51	14.634	1975	1980	1981	rBV4	16245	25573	0.53%	0.070%
52	14.675	1981	1987	1996	rVV	382544	727640	15.07%	1.988%
53	14.751	1996	2000	2007	rVB2	31652	53507	1.11%	0.146%
54	14.892	2020	2024	2027	rBV3	22532	44580	0.92%	0.122%
55	15.216	2074	2079	2081	rVV5	22453	34395	0.71%	0.094%
56	15.281	2084	2090	2095	rVV	222774	355056	7.36%	0.970%
57	15.339	2095	2100	2109	rVB	513503	832453	17.24%	2.275%
58	15.434	2112	2116	2118	rBV	37464	54209	1.12%	0.148%
59	15.463	2118	2121	2124	rVV2	82712	129881	2.69%	0.355%
60	15.498	2124	2127	2133	rVV4	93203	167506	3.47%	0.458%
61	15.551	2133	2136	2139	rVV3	28481	45255	0.94%	0.124%
62	15.592	2139	2143	2146	rVV2	49936	74984	1.55%	0.205%
63	15.645	2147	2152	2157	rVB	83368	129398	2.68%	0.354%
64	15.786	2167	2176	2184	rVV5	103786	299054	6.20%	0.817%
65	15.875	2185	2191	2199	rVB2	68209	145710	3.02%	0.398%
66	16.063	2218	2223	2232	rVB6	74391	170011	3.52%	0.465%
67	16.145	2233	2237	2243	rVB	79976	114684	2.38%	0.313%
68	16.281	2256	2260	2264	rVB2	18314	30063	0.62%	0.082%
69	16.369	2265	2275	2282	rBV5	108014	332044	6.88%	0.907%
70	16.610	2313	2316	2320	rBV4	16343	25301	0.52%	0.069%
71	16.892	2359	2364	2372	rVB	131065	226957	4.70%	0.620%
72	17.004	2380	2383	2386	rBV3	22653	29132	0.60%	0.080%
73	17.086	2391	2397	2401	rVV	1603972	2394296	49.60%	6.543%
74	17.128	2401	2404	2410	rVV	1542669	2044924	42.36%	5.588%
75	17.186	2410	2414	2417	rVV2	255215	394586	8.17%	1.078%
76	17.228	2417	2421	2429	rVV	185158	352539	7.30%	0.963%

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77	17.304	2430	2434	2440	rVB	68619	112371	2.33%	0.307%
78	17.528	2464	2472	2478	rBV	302700	486651	10.08%	1.330%
79	17.645	2489	2492	2497	rVB3	46804	66893	1.39%	0.183%
80	17.710	2500	2503	2509	rVB	50361	77228	1.60%	0.211%
81	17.780	2510	2515	2516	rBV2	42921	66795	1.38%	0.183%
82	17.986	2545	2550	2556	rBV3	144110	288830	5.98%	0.789%
83	18.045	2556	2560	2570	rVB4	126787	289683	6.00%	0.792%
84	18.151	2575	2578	2581	rVB2	36511	46308	0.96%	0.127%
85	18.198	2581	2586	2592	rBV2	153208	285333	5.91%	0.780%
86	18.322	2602	2607	2614	rBV4	78758	184667	3.83%	0.505%
87	18.422	2620	2624	2628	rBV4	38435	59552	1.23%	0.163%
88	18.469	2630	2632	2636	rVV2	35949	49501	1.03%	0.135%
89	18.522	2638	2641	2651	rVB2	47190	88573	1.83%	0.242%
90	18.875	2698	2701	2705	rVB5	29010	34655	0.72%	0.095%
91	19.227	2756	2761	2769	rVB	659103	976707	20.23%	2.669%
92	19.351	2779	2782	2787	rBV2	63832	85371	1.77%	0.233%
93	19.574	2815	2820	2823	rBV	421314	651704	13.50%	1.781%
94	19.604	2823	2825	2838	rVB	388176	600103	12.43%	1.640%
95	20.022	2893	2896	2900	rBV	65363	94141	1.95%	0.257%
96	20.127	2911	2914	2921	rVB2	111844	188184	3.90%	0.514%
97	20.239	2929	2933	2938	rBV	90054	147736	3.06%	0.404%
98	20.792	3023	3027	3038	rVB	257747	532473	11.03%	1.455%
99	21.533	3147	3153	3159	rBV	1343304	2283909	47.31%	6.241%
100	24.845	3708	3716	3732	rVB	912956	2248015	46.57%	6.143%

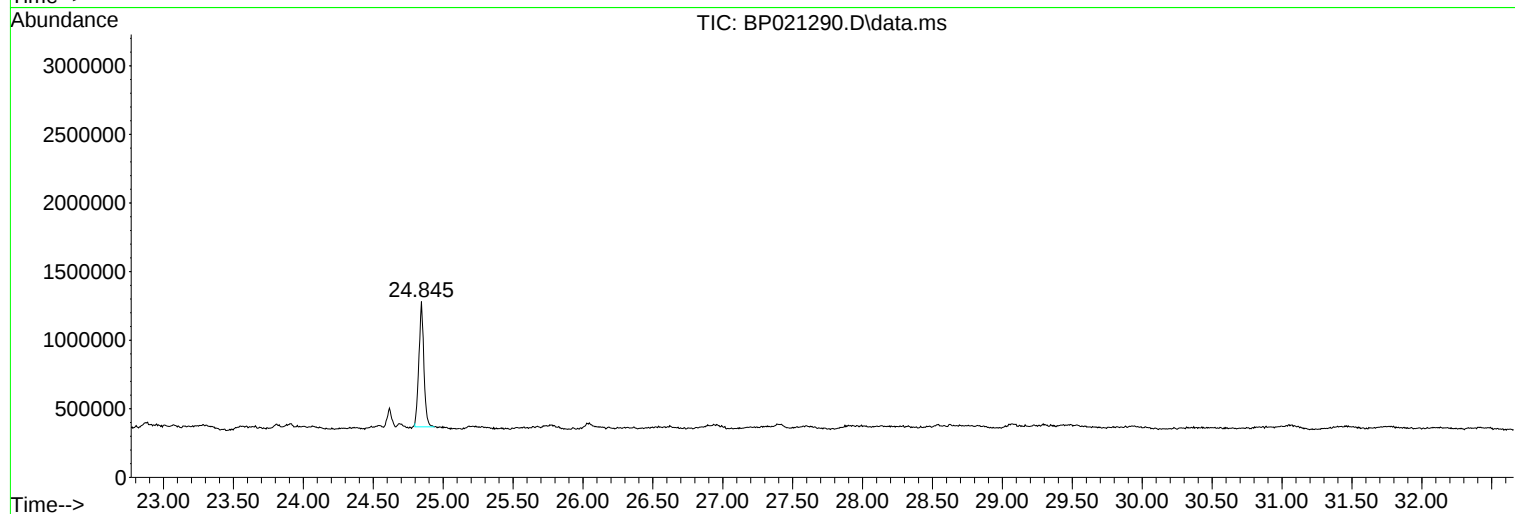
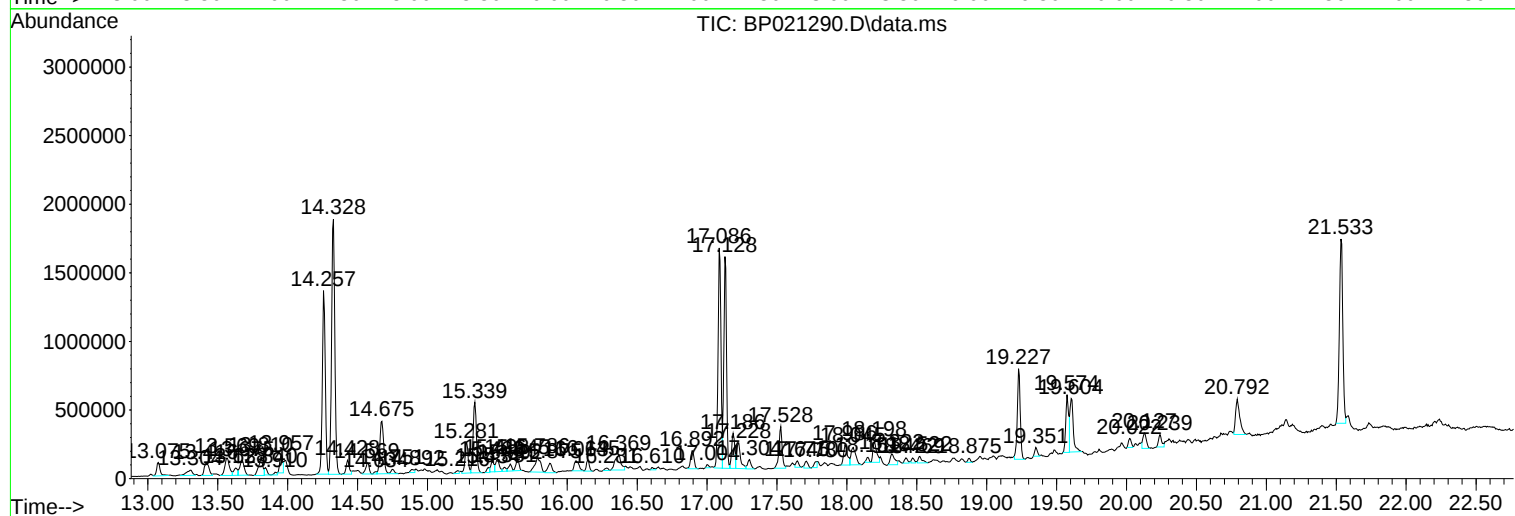
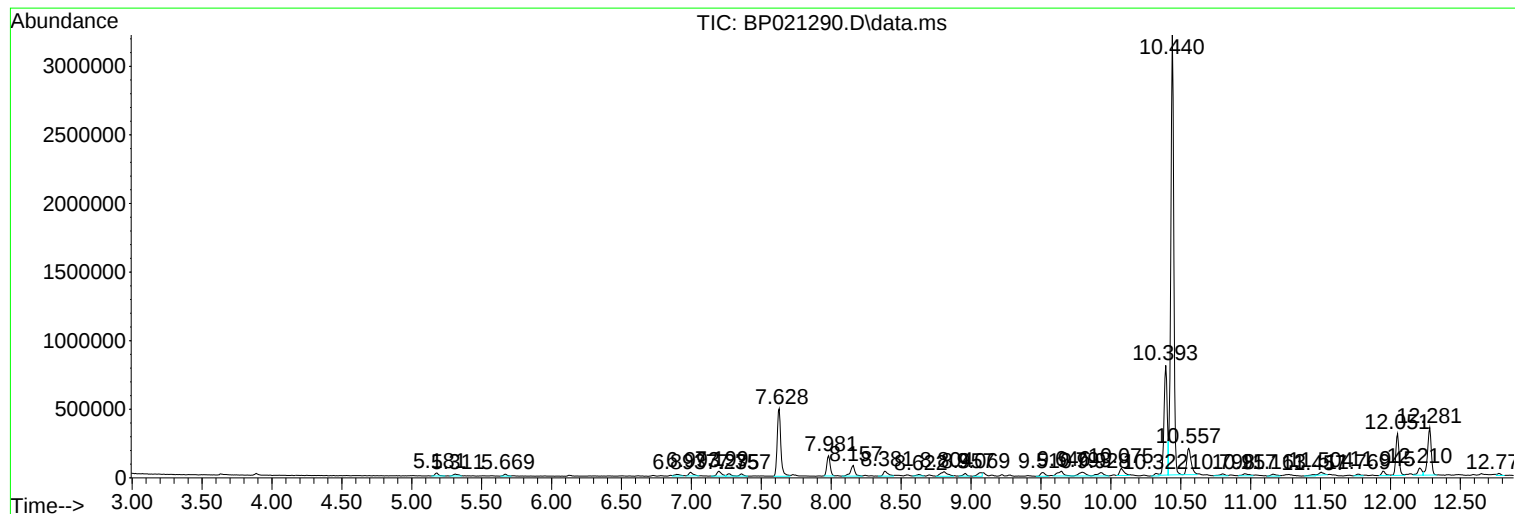
Sum of corrected areas: 36594265

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Instrument :
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C0AY8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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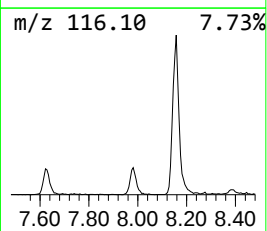
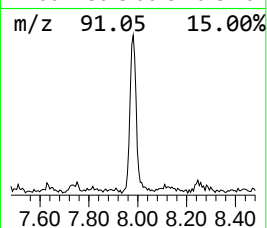
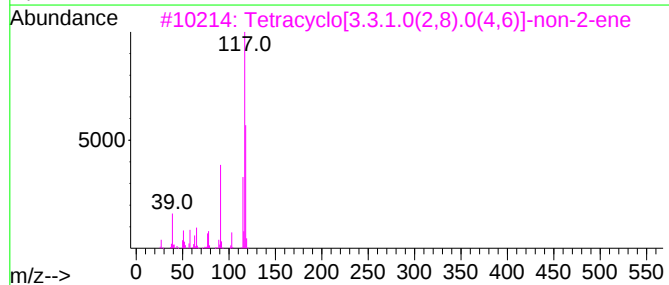
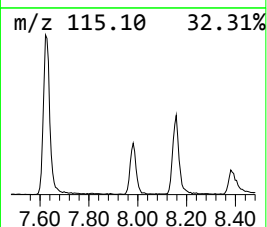
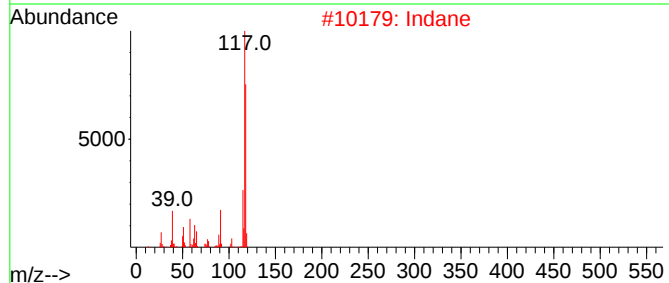
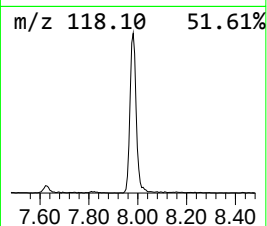
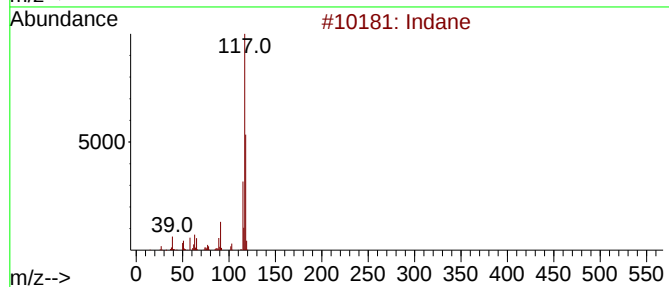
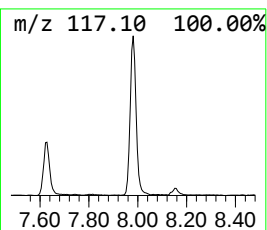
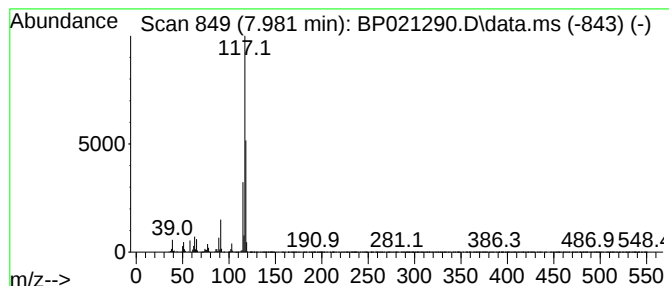
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.981	5.51 ng/ul	247485	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	91
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...			118	C9H10	1000191-13-7	87
4	Indane			118	C9H10	000496-11-7	83
5	Deltacyclene			118	C9H10	007785-10-6	74



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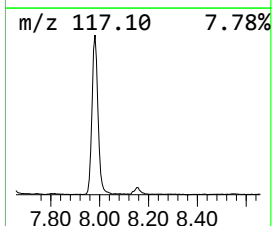
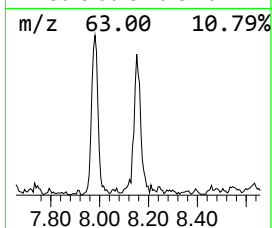
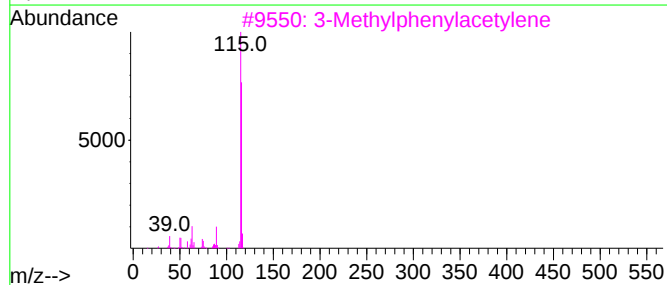
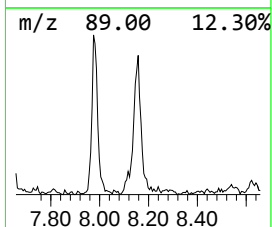
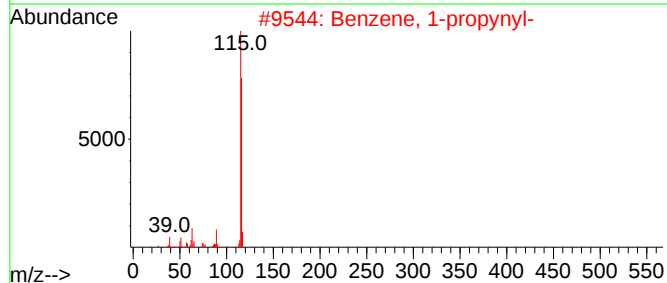
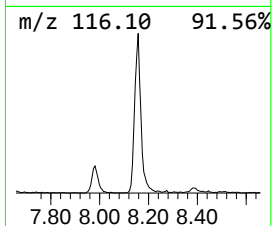
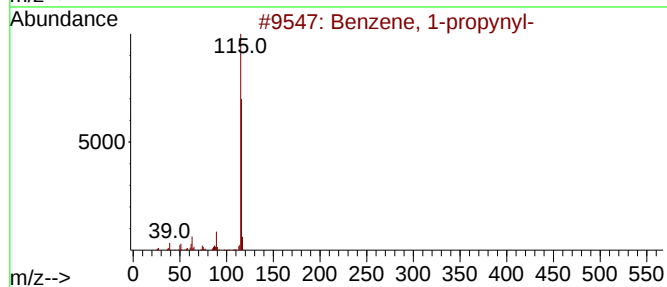
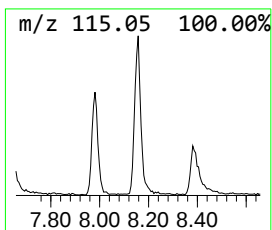
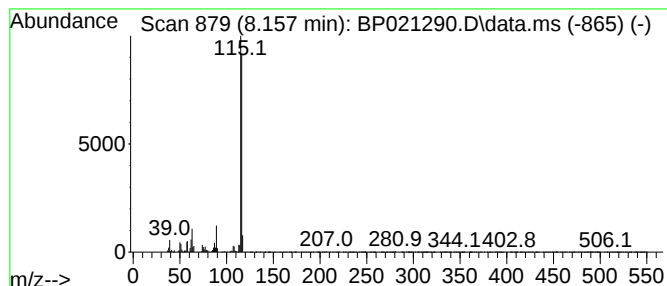
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1-propynyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.157	3.72 ng/ul	167009	1,4-Dichlorobenzene-d4	7.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3			3-Methylphenylacetylene	116	C9H8	000766-82-5	94
4			1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	91
5			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91



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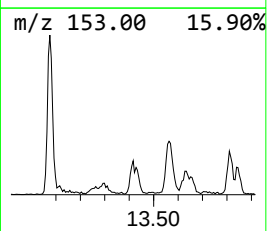
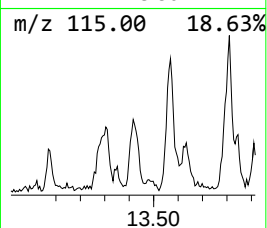
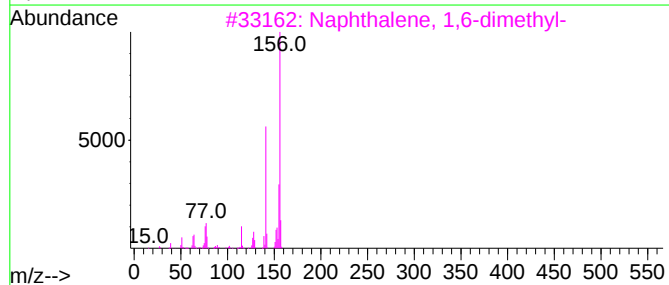
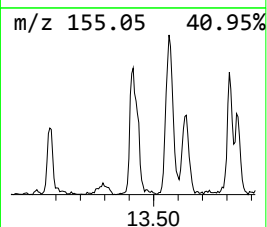
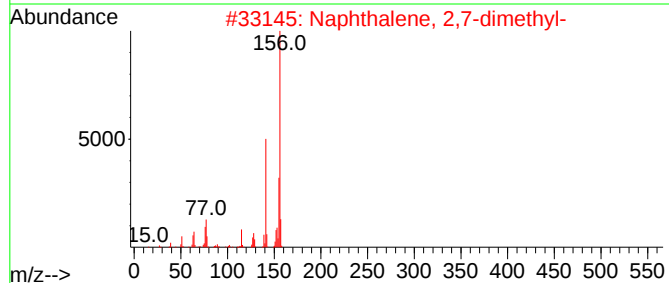
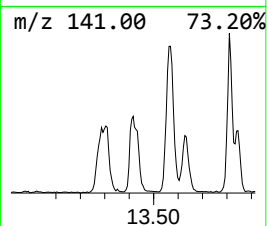
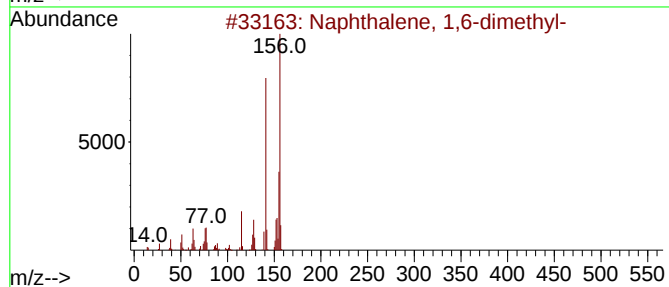
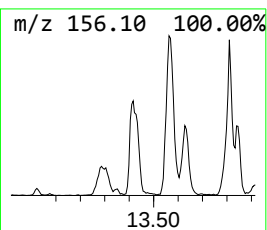
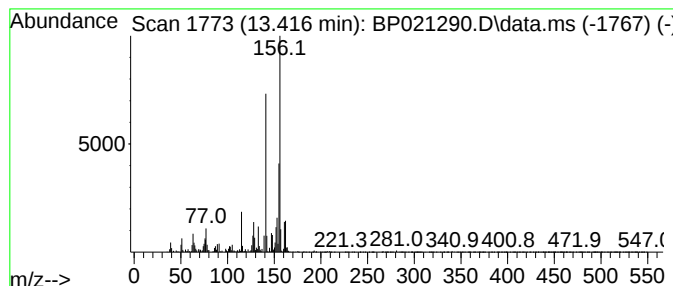
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 1,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.416	2.17 ng/ul	209506	Acenaphthene-d10	14.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
Operator : CG/JU
Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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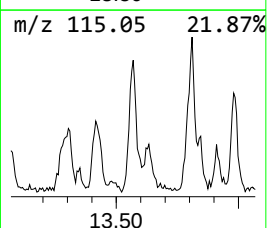
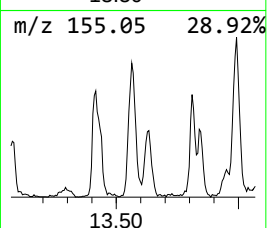
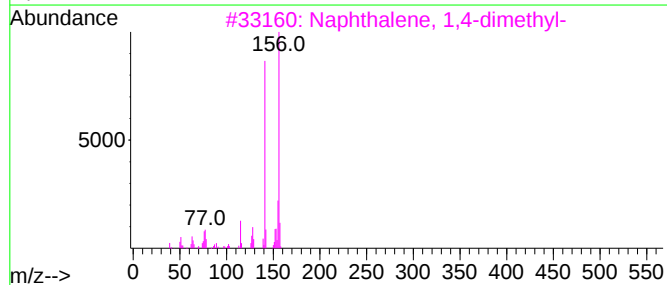
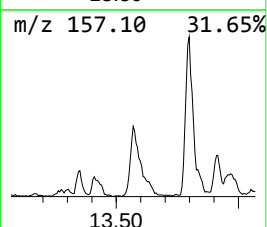
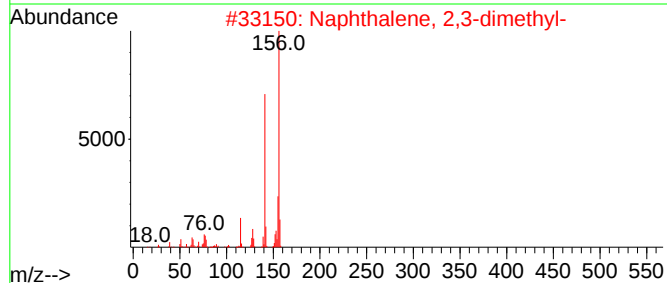
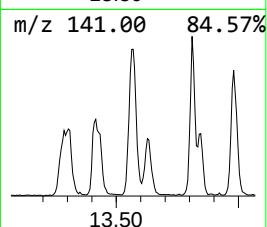
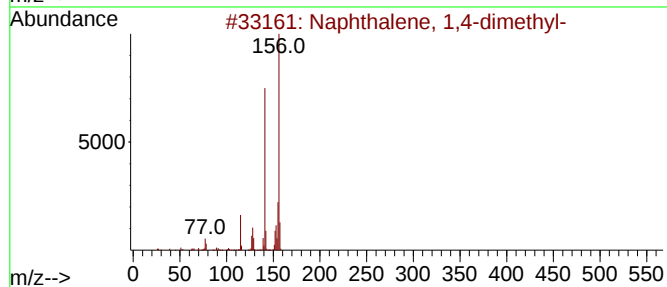
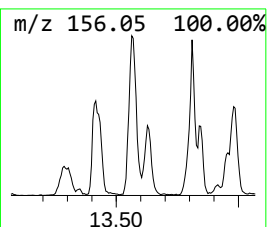
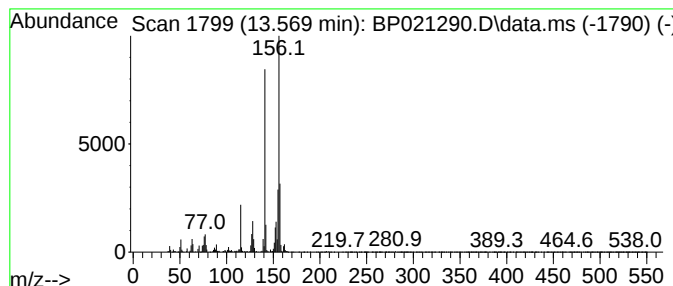
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.569	3.07 ng/ul	296288	Acenaphthene-d10	14.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	94
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
Operator : CG/JU
Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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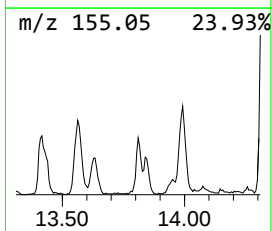
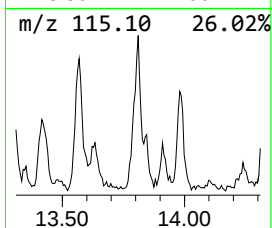
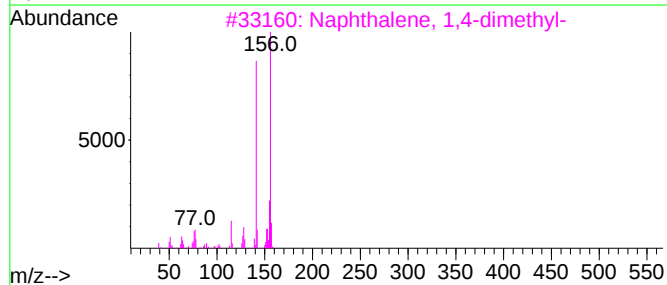
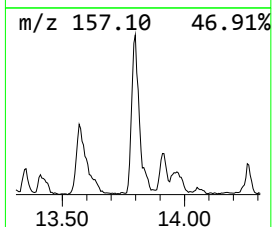
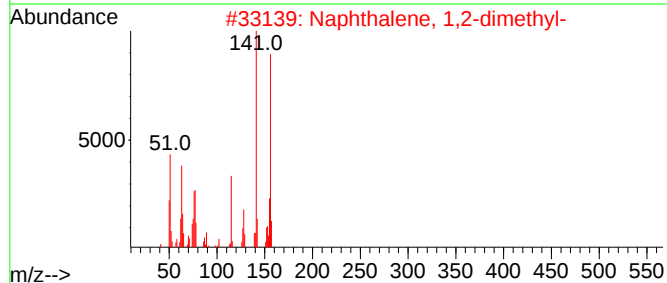
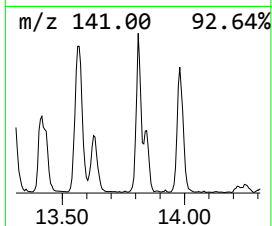
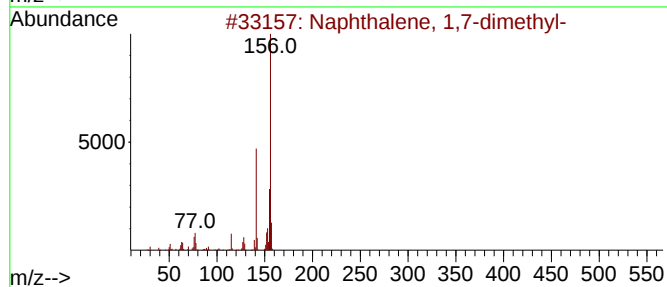
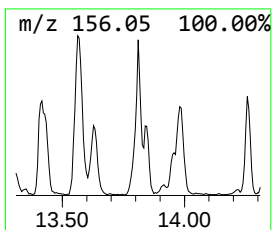
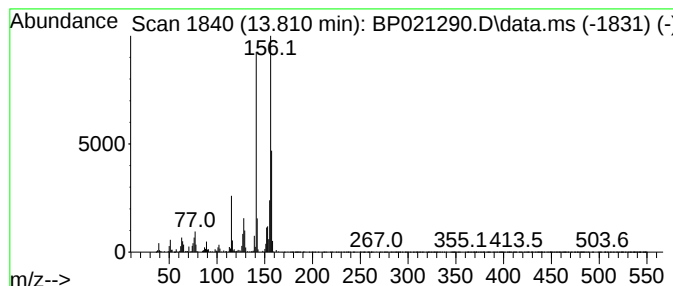
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1,7-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	2.82 ng/ul	272221	Acenaphthene-d10	14.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
2			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	94
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	94
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	94
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
Operator : CG/JU
Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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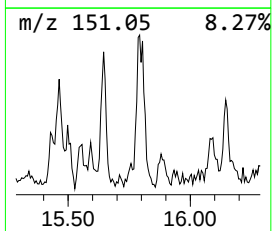
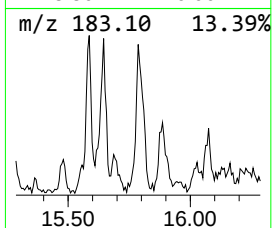
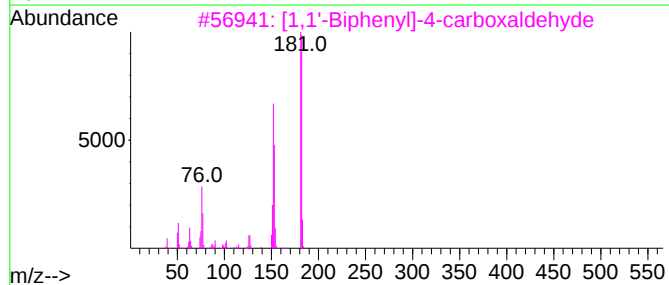
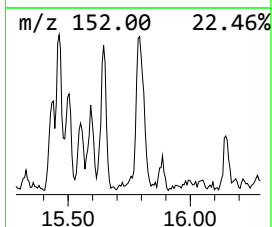
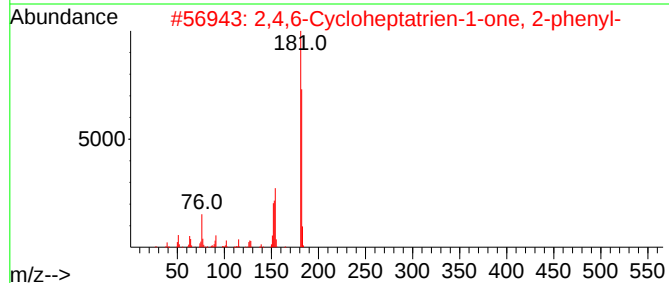
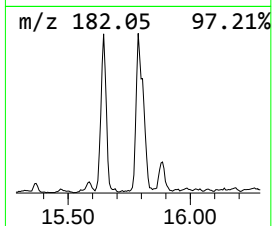
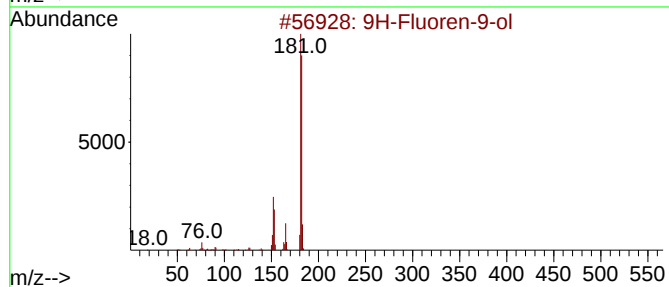
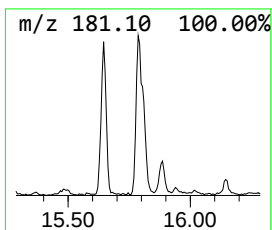
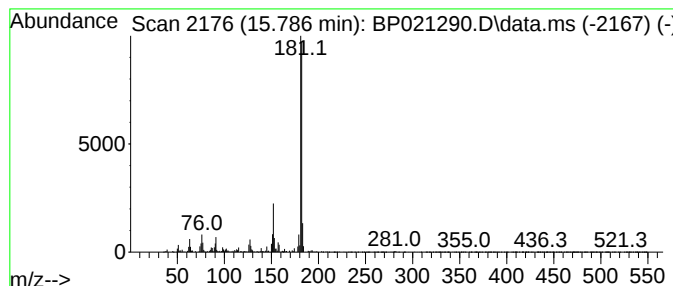
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9H-Fluoren-9-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.787	2.50 ng/ul	299054	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
2			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86
3			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86
4			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	74
5			Pyrido[2,3-b]indole, 6-methyl-	182	C12H10N2	108349-67-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
Operator : CG/JU
Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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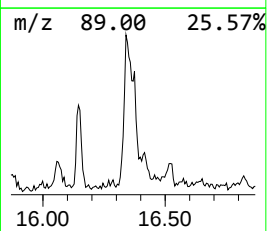
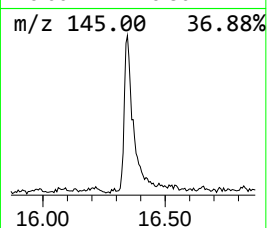
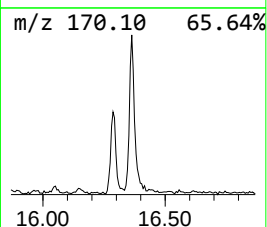
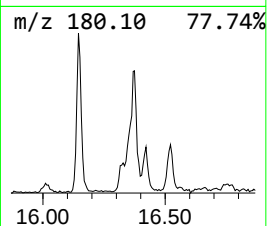
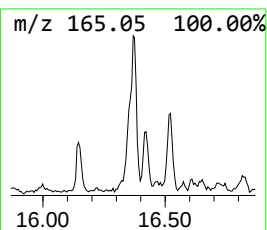
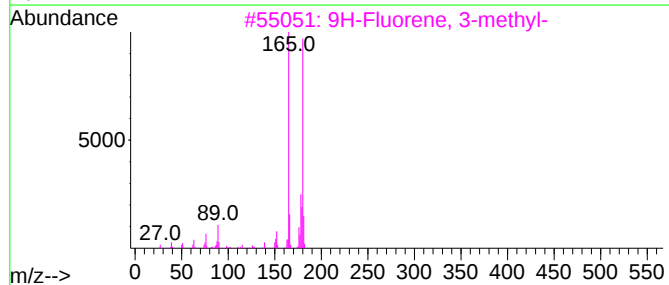
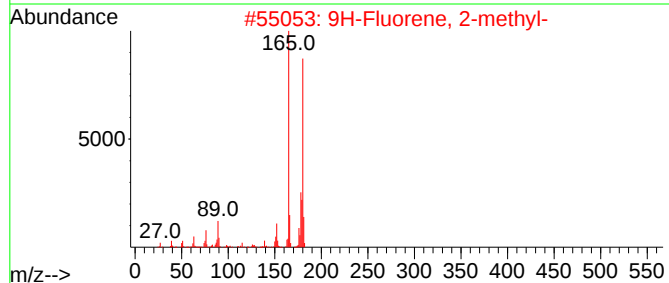
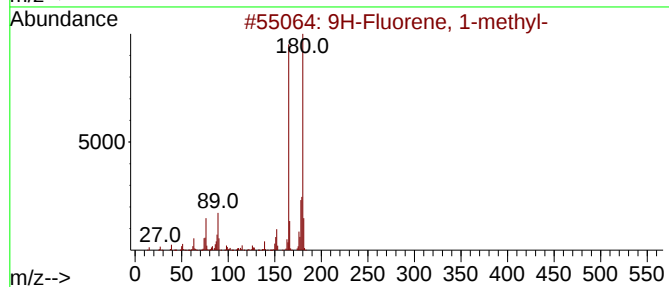
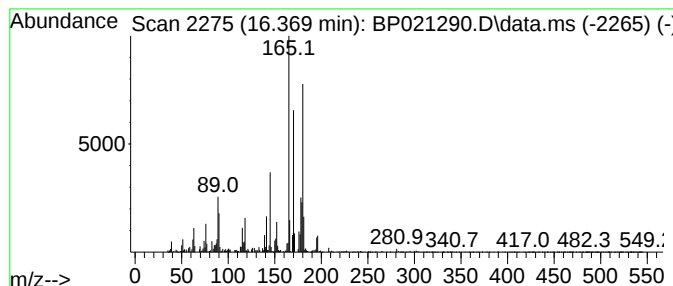
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9H-Fluorene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.369	2.77 ng/ul	332044	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 1-methyl-			180	C14H12	001730-37-6	96
2	9H-Fluorene, 2-methyl-			180	C14H12	001430-97-3	96
3	9H-Fluorene, 3-methyl-			180	C14H12	002523-39-9	93
4	9H-Fluorene, 1-methyl-			180	C14H12	001730-37-6	91
5	9H-Fluorene, 4-methyl-			180	C14H12	001556-99-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
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Sample : P3347-14DL 10X
Misc :
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Instrument :
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ClientSampleId :
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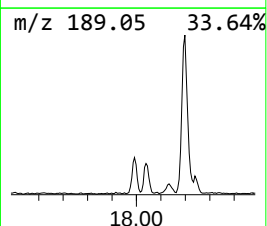
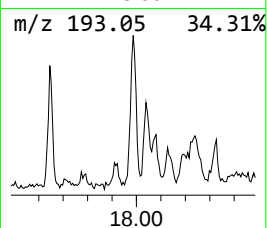
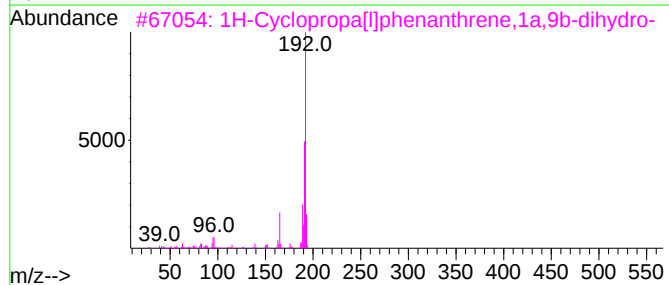
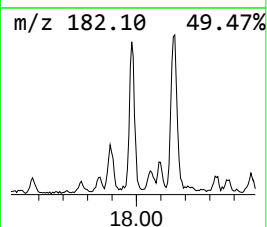
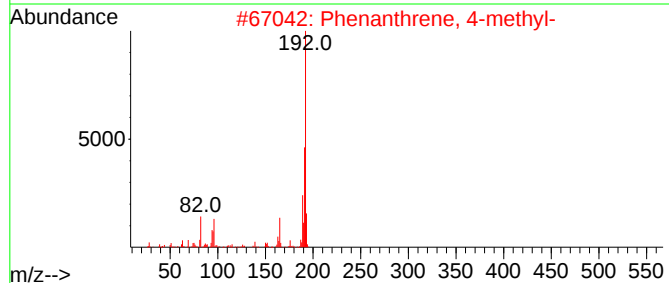
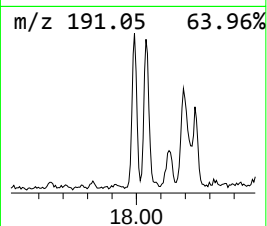
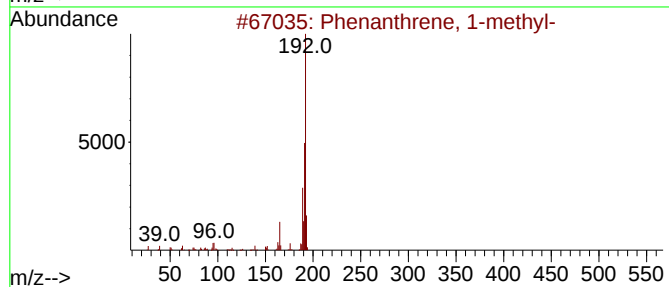
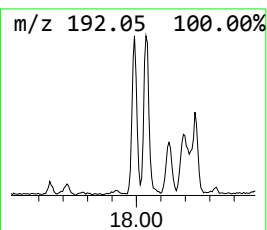
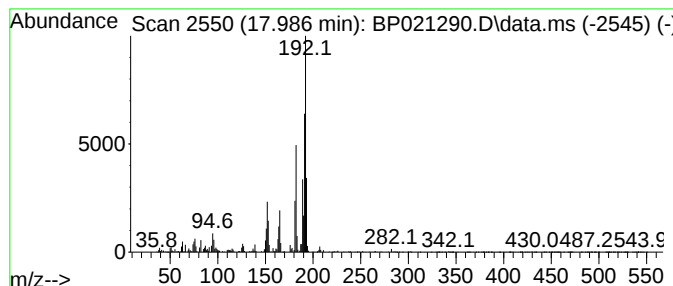
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.986	2.41 ng/ul	288830	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
2			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	78
3			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	60
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	60
5			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
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Misc :
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ClientSampleId :
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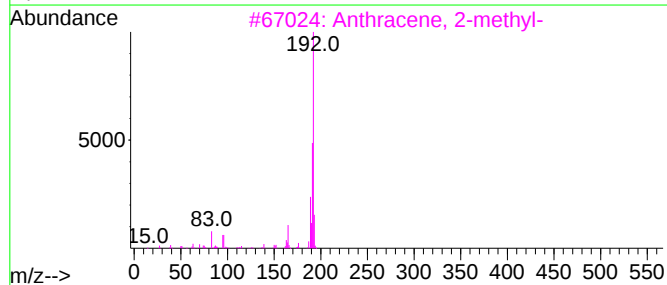
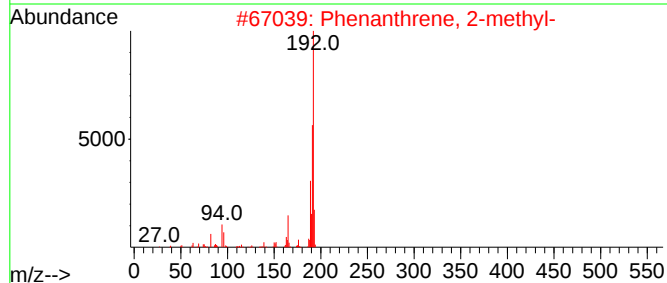
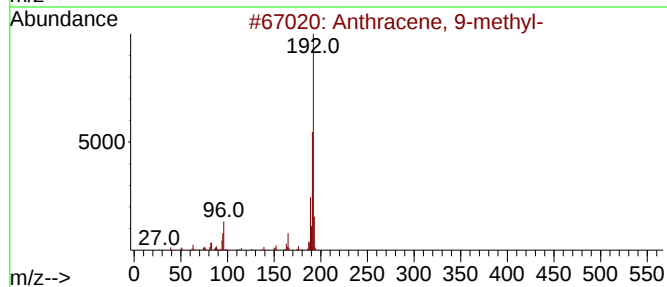
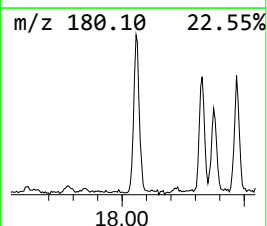
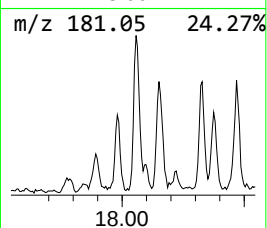
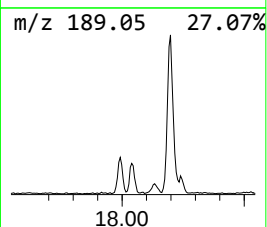
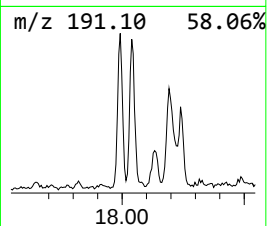
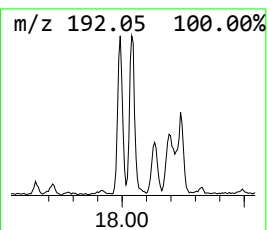
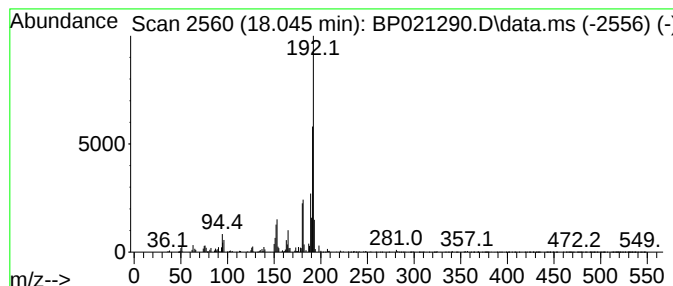
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Anthracene, 9-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.045	2.42 ng/ul	289683	Phenanthrene-d10	17.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
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Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

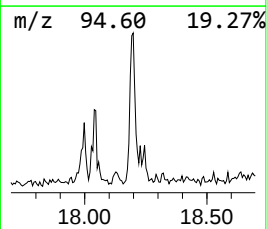
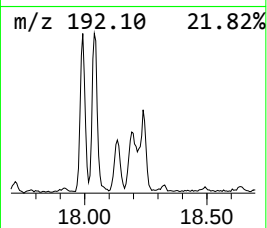
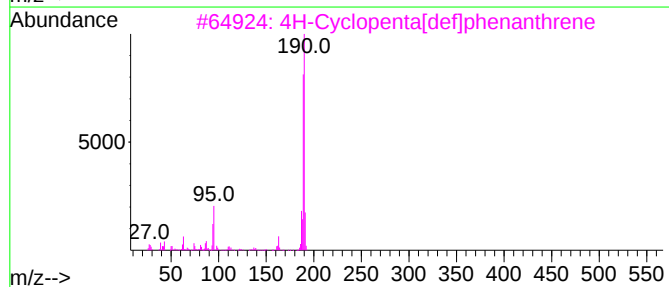
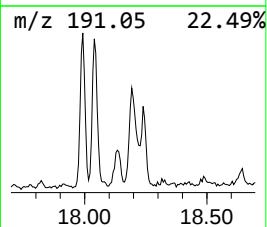
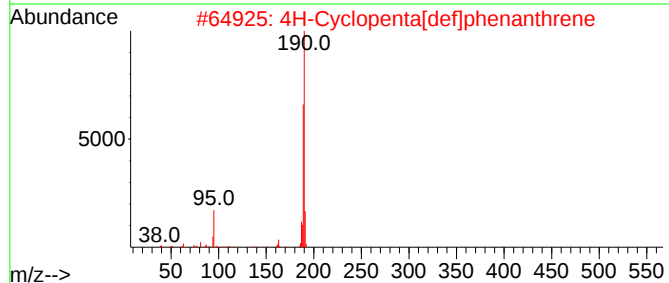
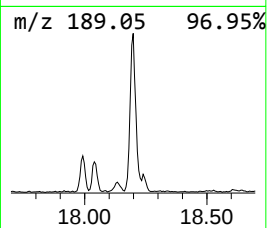
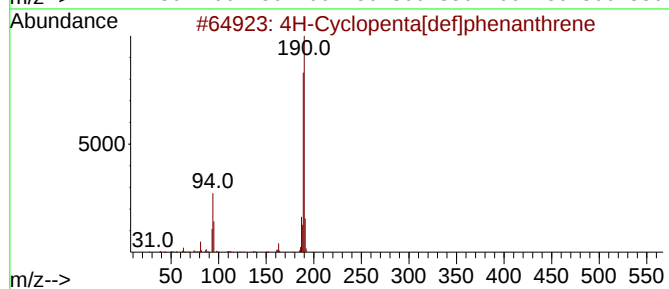
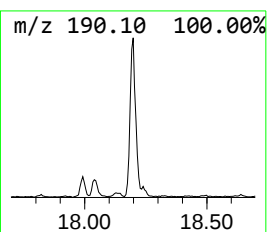
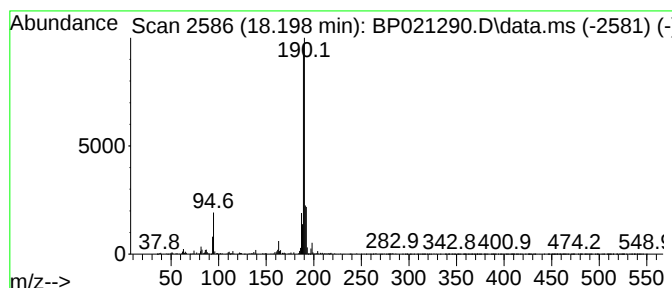
Instrument :
BNA_P
ClientSampleId :
C0AY8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.198	2.38 ng/ul	285333	Phenanthrene-d10		17.086	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	90
2	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	87
3	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	81
4	Phenol, 4-(2-thienylmethyl)-		190	C11H10O5	091680-55-6	59
5	Methyl diselenide		190	C2H6Se2	007101-31-7	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
Operator : CG/JU
Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY8DL

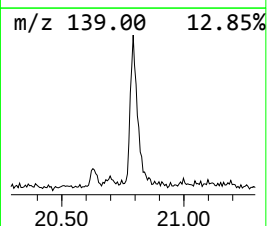
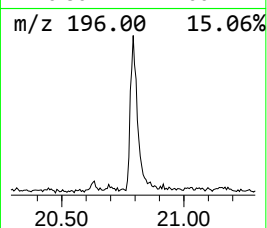
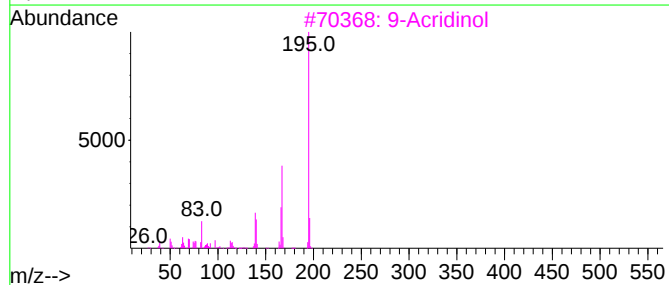
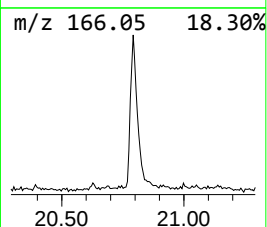
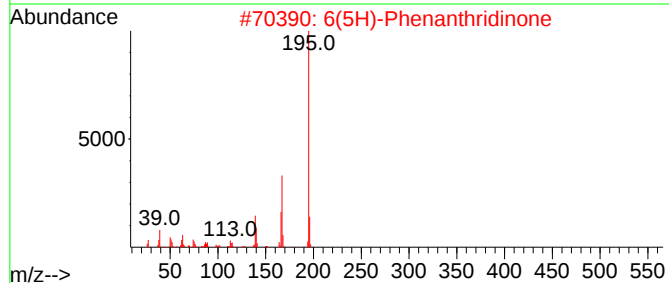
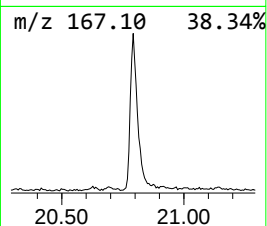
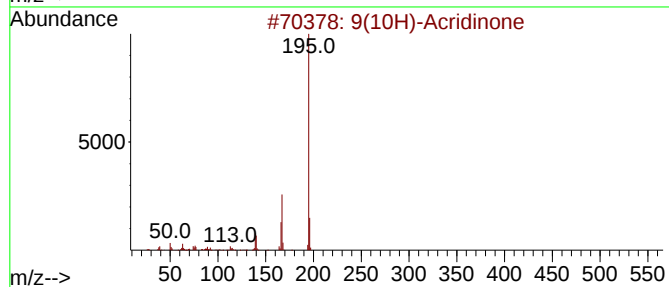
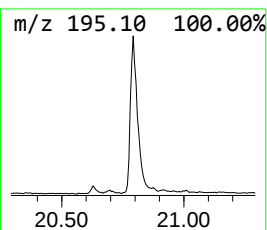
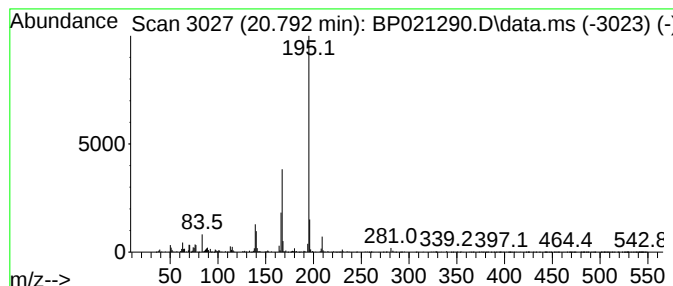
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 9(10H)-Acridinone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.792	4.66 ng/ul	532473	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9(10H)-Acridinone			195	C13H9NO	000578-95-0	96
2	6(5H)-Phenanthridinone			195	C13H9NO	001015-89-0	94
3	9-Acridinol			195	C13H9NO	000643-62-9	94
4	4-Amino-9-fluorenone			195	C13H9NO	004269-15-2	94
5	9(10H)-Acridinone			195	C13H9NO	000578-95-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021290.D
Acq On : 01 Aug 2024 12:40
Operator : CG/JU
Sample : P3347-14DL 10X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY8DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Indane	7.981	5.5	ng/ul	247485	1	7.628	897986	20.0
Benzene, 1-prop...	8.157	3.7	ng/ul	167009	1	7.628	897986	20.0
Naphthalene, 1,...	13.416	2.2	ng/ul	209506	3	14.257	1931540	20.0
Naphthalene, 1,...	13.569	3.1	ng/ul	296288	3	14.257	1931540	20.0
Naphthalene, 1,...	13.810	2.8	ng/ul	272221	3	14.257	1931540	20.0
9H-Fluoren-9-ol	15.787	2.5	ng/ul	299054	4	17.086	2394300	20.0
9H-Fluorene, 1-...	16.369	2.8	ng/ul	332044	4	17.086	2394300	20.0
Phenanthrene, 1...	17.986	2.4	ng/ul	288830	4	17.086	2394300	20.0
Anthracene, 9-m...	18.045	2.4	ng/ul	289683	4	17.086	2394300	20.0
4H-Cyclopenta[d...	18.198	2.4	ng/ul	285333	4	17.086	2394300	20.0
9(10H)-Acridinone	20.792	4.7	ng/ul	532473	5	21.533	2283910	20.0