

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
 Data File : BP022823.D
 Acq On : 04 Nov 2024 12:17
 Operator : RC/JU
 Sample : P4568-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4EA4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022823.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.176	29	32	42	rVB	9153	14076	0.11%	0.009%
2	4.423	237	244	255	rVB5	12195	23922	0.19%	0.014%
3	6.852	647	657	672	rBV2	264968	672591	5.22%	0.406%
4	6.975	672	678	684	rVV	113648	169387	1.31%	0.102%
5	7.064	686	693	705	rVV	1012500	1591249	12.35%	0.961%
6	7.205	705	717	735	rVB2	666479	1301593	10.10%	0.786%
7	7.664	790	795	804	rVB	1860397	2977478	23.11%	1.798%
8	7.975	839	848	859	rBV	1847688	2989180	23.20%	1.805%
9	8.346	904	911	918	rBV	178285	303051	2.35%	0.183%
10	8.428	918	925	946	rVB	1359370	2109362	16.37%	1.274%
11	8.681	960	968	978	rVB	1017992	1741873	13.52%	1.052%
12	8.775	978	984	987	rBV	167362	283603	2.20%	0.171%
13	8.811	987	990	1002	rVB	294789	464652	3.61%	0.281%
14	9.169	1045	1051	1058	rVB4	9323	17288	0.13%	0.010%
15	9.334	1071	1079	1095	rBV	1779764	2855934	22.17%	1.725%
16	9.487	1098	1105	1106	rVV	158757	246029	1.91%	0.149%
17	9.510	1106	1109	1121	rVV	326105	549670	4.27%	0.332%
18	9.640	1124	1131	1138	rVB	63912	107304	0.83%	0.065%
19	9.805	1152	1159	1172	rVB	1475333	2190283	17.00%	1.323%
20	10.046	1189	1200	1222	rBV2	1070033	2359707	18.31%	1.425%
21	10.381	1244	1257	1264	rBV	355928	624456	4.85%	0.377%
22	10.528	1275	1282	1303	rBV	126432	275878	2.14%	0.167%
23	10.710	1304	1313	1330	rVB	2262987	3558233	27.62%	2.149%
24	11.681	1464	1478	1493	rBV	408455	780781	6.06%	0.471%
25	11.934	1515	1521	1532	rVB3	17264	37055	0.29%	0.022%
26	12.046	1532	1540	1547	rBV	944566	1565856	12.15%	0.946%
27	12.475	1608	1613	1621	rBV	21652	30620	0.24%	0.018%
28	12.663	1636	1645	1653	rBV	2740476	4037771	31.34%	2.438%
29	12.746	1653	1659	1675	rVB	444770	883038	6.85%	0.533%
30	13.651	1807	1813	1824	rVB	476233	693619	5.38%	0.419%
31	13.840	1836	1845	1853	rBV	2662389	4145326	32.17%	2.503%
32	13.940	1855	1862	1865	rVV2	549036	943892	7.33%	0.570%
33	13.969	1865	1867	1878	rVB	425638	468174	3.63%	0.283%
34	14.257	1905	1916	1920	rBV	637634	945204	7.34%	0.571%
35	14.316	1920	1926	1932	rVV	3153143	4241078	32.92%	2.561%
36	14.375	1932	1936	1946	rVB3	21235	34573	0.27%	0.021%

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37	14.504	1950	1958	1968	rBV3	656270	1499446	11.64%	0.905%
38	14.581	1968	1971	1975	rVV2	85396	133428	1.04%	0.081%
39	14.640	1975	1981	1998	rVB	2442475	2928811	22.73%	1.769%
40	14.910	2021	2027	2035	rVB2	35331	59763	0.46%	0.036%
41	15.145	2057	2067	2075	rVB2	8495	18985	0.15%	0.011%
42	15.269	2080	2088	2091	rBV	710443	1118209	8.68%	0.675%
43	15.322	2091	2097	2105	rVV2	5919913	9226852	71.61%	5.572%
44	15.404	2105	2111	2120	rVV	152148	272015	2.11%	0.164%
45	15.575	2133	2140	2146	rBV	52246	96989	0.75%	0.059%
46	15.634	2146	2150	2156	rVB	207726	271784	2.11%	0.164%
47	15.928	2196	2200	2208	rVV6	9314	14519	0.11%	0.009%
48	16.016	2208	2215	2226	rVB4	23856	55387	0.43%	0.033%
49	16.222	2240	2250	2259	rBV	945999	1420918	11.03%	0.858%
50	16.351	2265	2272	2277	rBV	7853907	9769820	75.82%	5.900%
51	16.710	2323	2333	2347	rBV	4294307	6513787	50.55%	3.934%
52	17.057	2385	2392	2395	rBV	916749	1377914	10.69%	0.832%
53	17.098	2395	2399	2403	rVV	1640284	2161814	16.78%	1.305%
54	17.151	2403	2408	2410	rVV	990237	1292725	10.03%	0.781%
55	17.187	2410	2414	2422	rVB	1519958	2061476	16.00%	1.245%
56	17.351	2434	2442	2445	rBV8	8144	18766	0.15%	0.011%
57	17.475	2458	2463	2475	rVB5	15524	36004	0.28%	0.022%
58	17.687	2494	2499	2505	rBV5	43489	76272	0.59%	0.046%
59	17.745	2505	2509	2514	rVB	64967	75075	0.58%	0.045%
60	17.928	2535	2540	2544	rBV8	11069	19783	0.15%	0.012%
61	17.981	2544	2549	2555	rBV	175070	246785	1.92%	0.149%
62	18.063	2556	2563	2568	rVV	4854761	5989021	46.48%	3.617%
63	18.163	2577	2580	2584	rVB6	13853	18286	0.14%	0.011%
64	18.275	2595	2599	2610	rVB6	15203	35623	0.28%	0.022%
65	18.386	2612	2618	2625	rBV6	19347	59181	0.46%	0.036%
66	18.451	2626	2629	2635	rVV6	16078	28777	0.22%	0.017%
67	18.528	2638	2642	2649	rVB	162955	205692	1.60%	0.124%
68	18.763	2680	2682	2686	rVB4	10014	14167	0.11%	0.009%
69	18.857	2692	2698	2702	rBV2	46153	91280	0.71%	0.055%
70	18.969	2715	2717	2721	rVB5	15758	19089	0.15%	0.012%
71	19.069	2730	2734	2737	rBV4	13048	21705	0.17%	0.013%
72	19.104	2737	2740	2745	rVV3	23987	32804	0.25%	0.020%
73	19.186	2747	2754	2763	rVV	5935529	8412135	65.29%	5.080%
74	19.322	2773	2777	2783	rVB8	36097	61125	0.47%	0.037%
75	19.434	2791	2796	2802	rVB	39151	68452	0.53%	0.041%
76	19.533	2807	2813	2815	rBV	1061850	1649527	12.80%	0.996%

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77	19.563	2815	2818	2830	rVB	2485486	3616259	28.07%	2.184%
78	20.098	2905	2909	2916	rBV5	44520	94281	0.73%	0.057%
79	20.257	2932	2936	2942	rBV3	93778	122042	0.95%	0.074%
80	20.510	2974	2979	2984	rVB	1377896	1493505	11.59%	0.902%
81	21.163	3087	3090	3099	rVB4	113828	233954	1.82%	0.141%
82	21.304	3111	3114	3122	rVB5	77810	112147	0.87%	0.068%
83	21.392	3122	3129	3134	rBV	4942681	7108039	55.17%	4.292%
84	21.463	3134	3141	3147	rVV2	5699020	9943379	77.17%	6.005%
85	21.528	3147	3152	3162	rVB	8446094	12884779	100.00%	7.781%
86	23.716	3512	3524	3536	rVB	5292965	12264700	95.19%	7.406%
87	24.510	3650	3659	3665	rBV	663212	1840026	14.28%	1.111%
88	24.580	3665	3671	3683	rVB	987813	2480209	19.25%	1.498%
89	24.727	3687	3696	3707	rVB	692458	1739193	13.50%	1.050%
90	28.486	4314	4335	4354	rVB	1535146	7099242	55.10%	4.287%
91	29.545	4506	4515	4519	rBV	296718	856171	6.64%	0.517%

Sum of corrected areas: 165595903

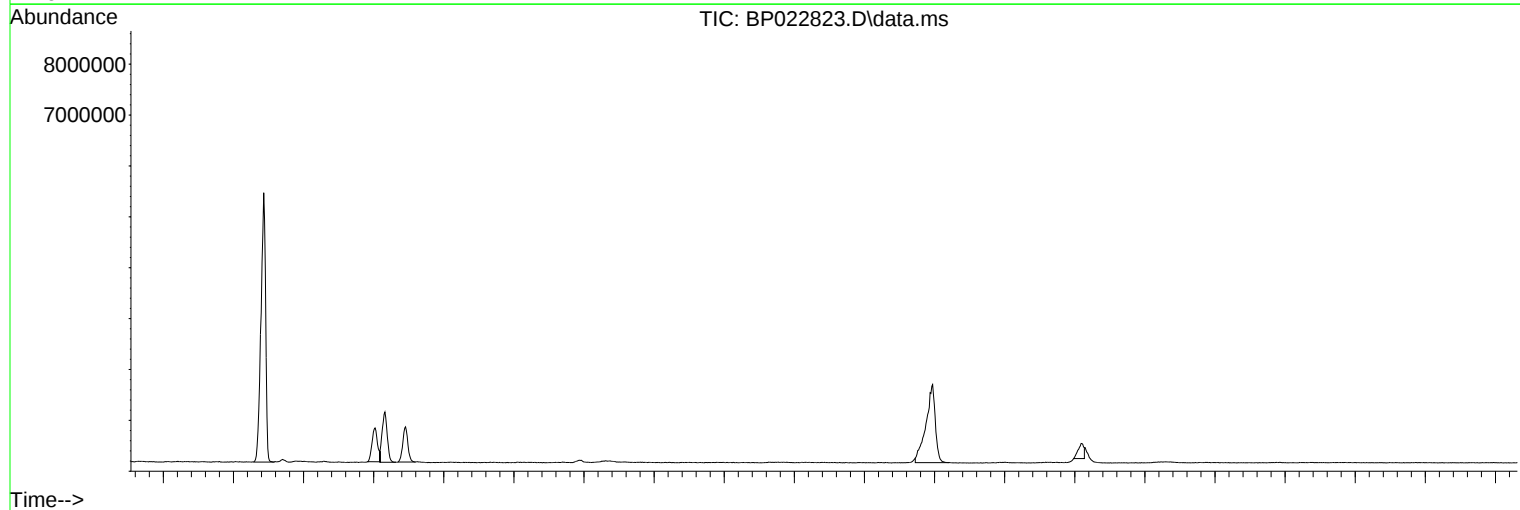
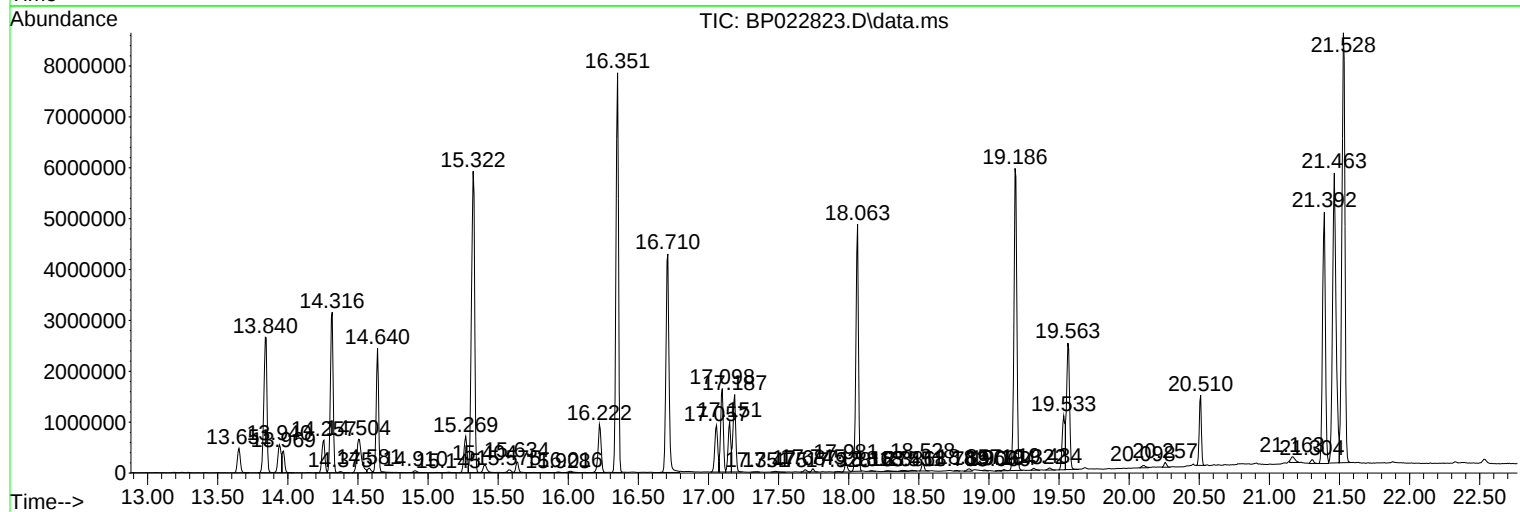
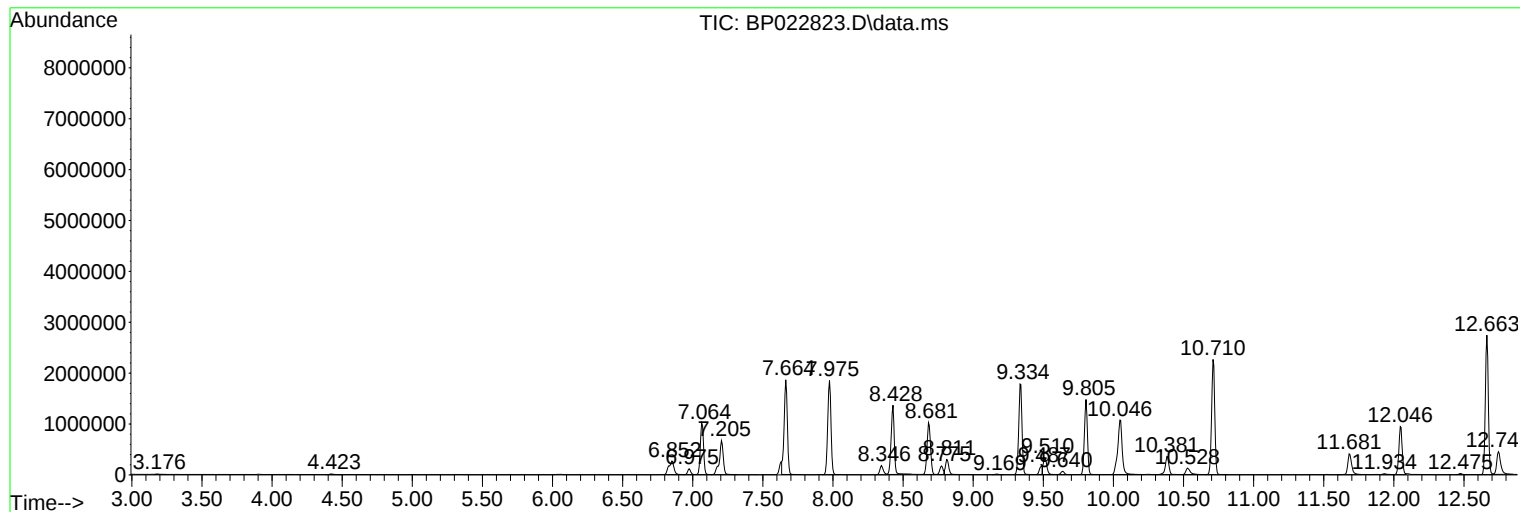
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P



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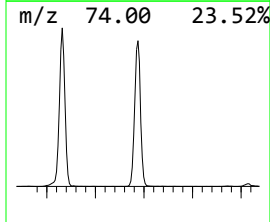
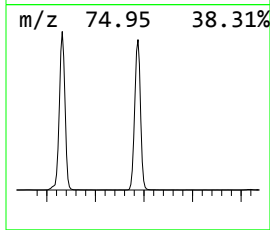
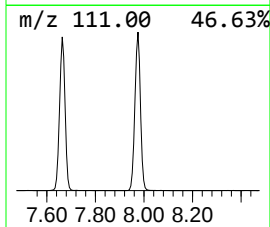
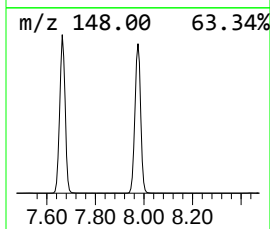
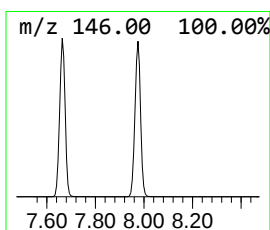
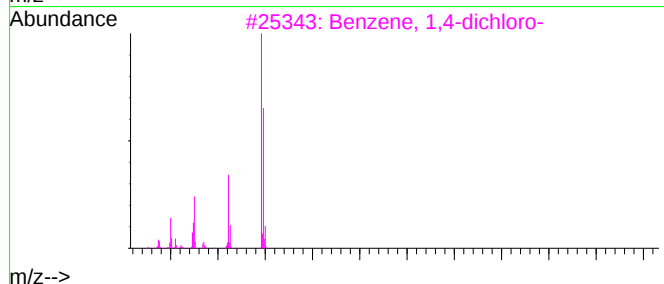
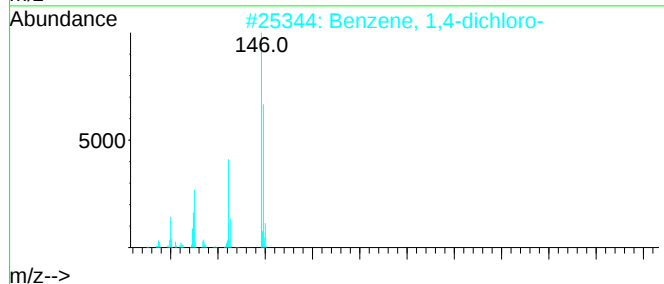
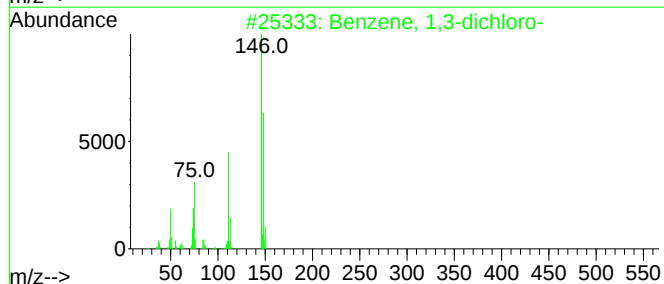
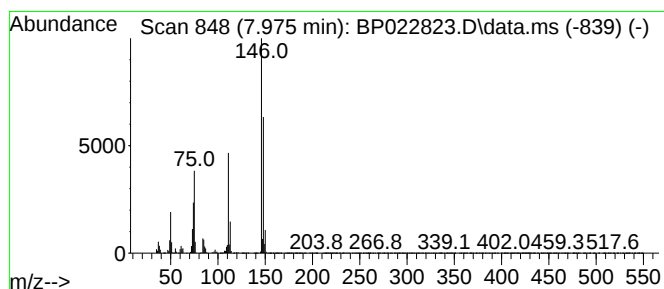
TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.975	20.08 ng/ul	2989180	1,4-Dichlorobenzene-d4	7.628

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
4	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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Quant Title : SVOA CALIBRATION

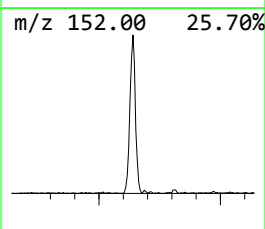
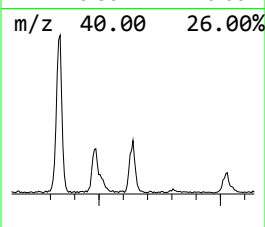
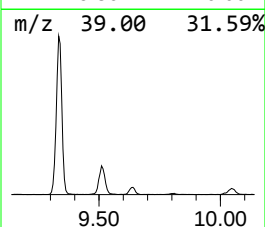
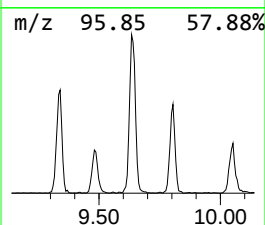
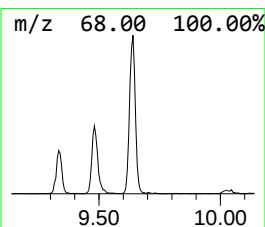
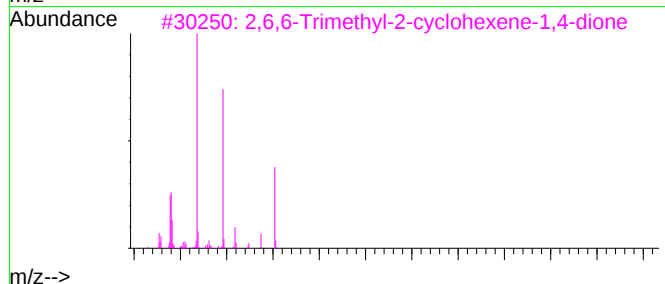
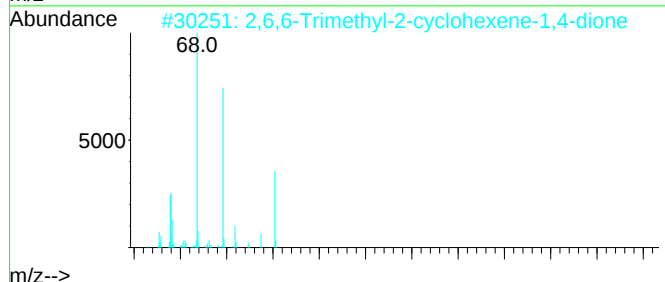
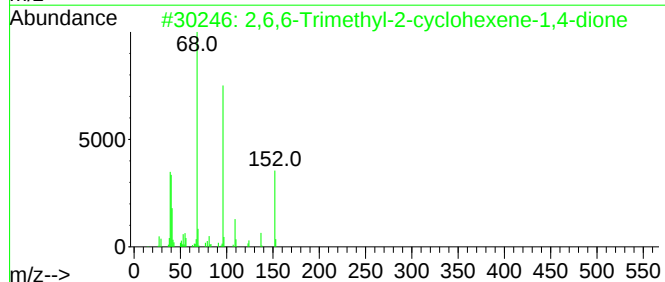
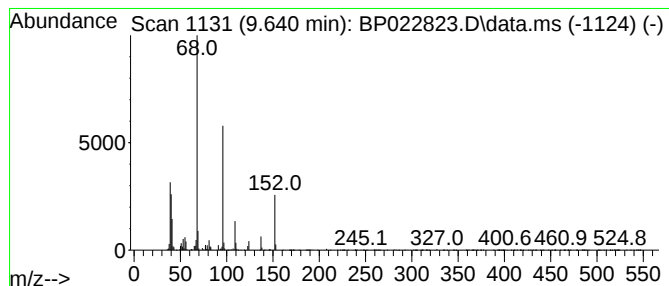
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Peak Number 2 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.640	3.44 ng/ul	107304	Naphthalene-d8	10.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	96
2	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	94
3	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	94
4	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	62
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	53



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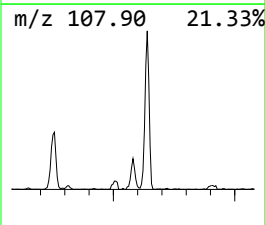
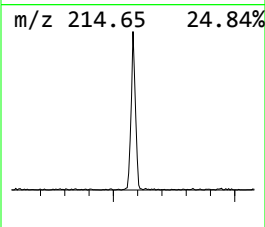
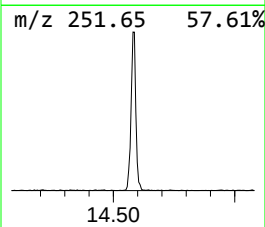
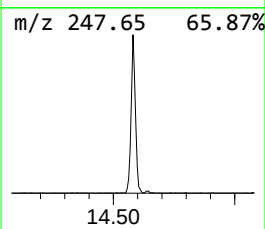
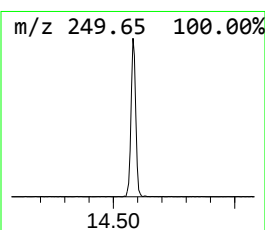
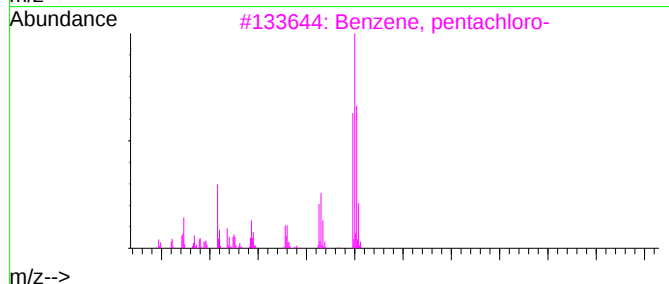
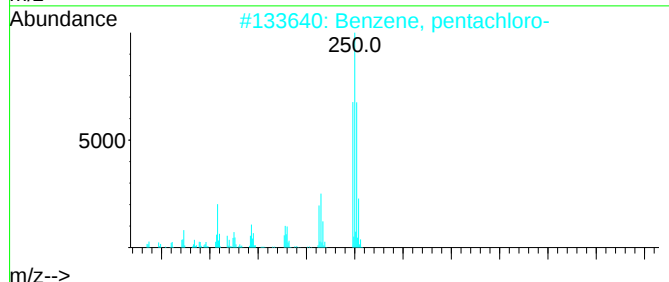
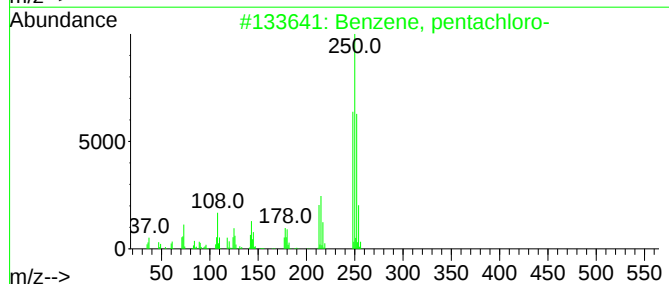
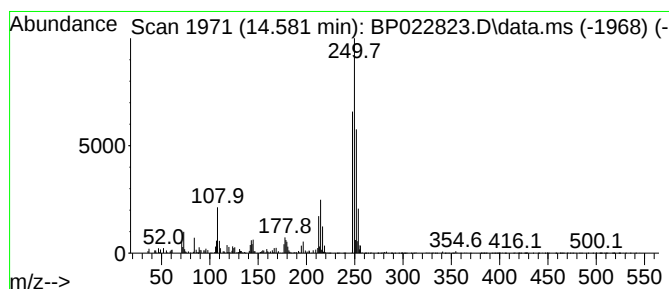
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, pentachloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.581	2.82 ng/ul	133428	Acenaphthene-d10		14.257	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, pentachloro-		248	C6HC15	000608-93-5	95
2	Benzene, pentachloro-		248	C6HC15	000608-93-5	94
3	Benzene, pentachloro-		248	C6HC15	000608-93-5	94
4	Benzene, pentachloro-		248	C6HC15	000608-93-5	93
5	Benzene, pentachloro-		248	C6HC15	000608-93-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022823.D
Acq On : 04 Nov 2024 12:17
Operator : RC/JU
Sample : P4568-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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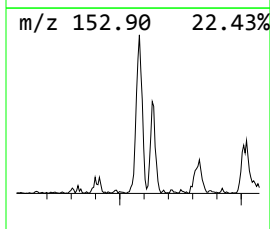
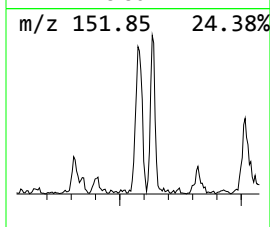
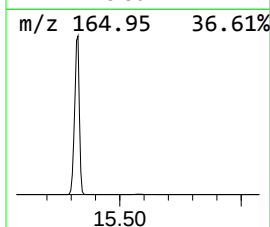
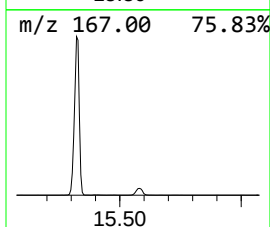
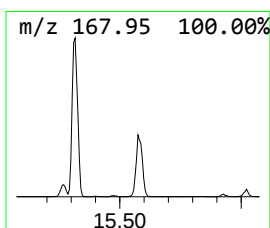
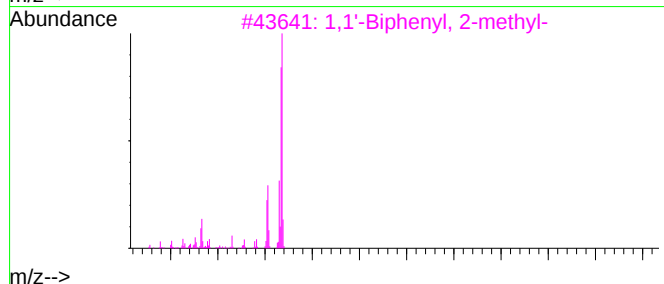
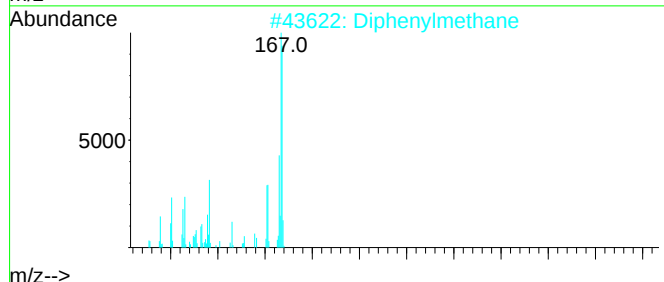
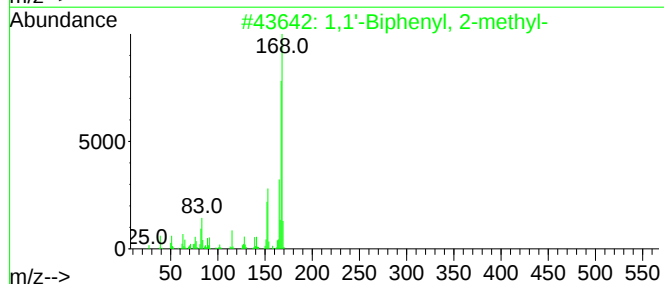
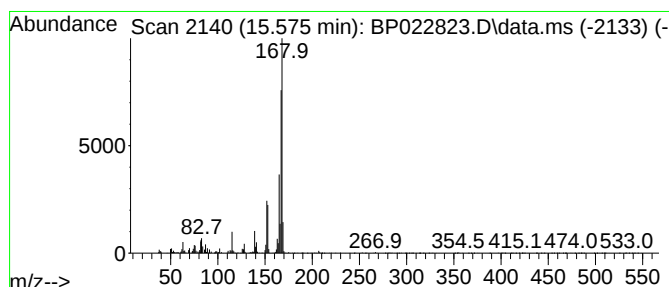
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 1,1'-Biphenyl, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.575	2.05 ng/ul	96989	Acenaphthene-d10	14.257	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	94
2	Diphenylmethane	168	C13H12	000101-81-5	93
3	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	91
4	Diphenylmethane	168	C13H12	000101-81-5	90
5	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022823.D
Acq On : 04 Nov 2024 12:17
Operator : RC/JU
Sample : P4568-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EA4

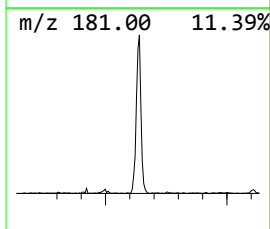
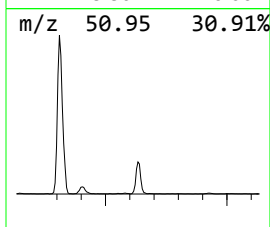
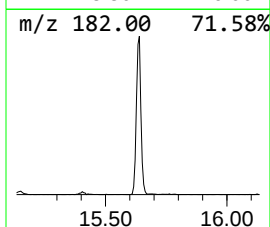
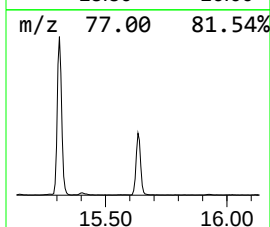
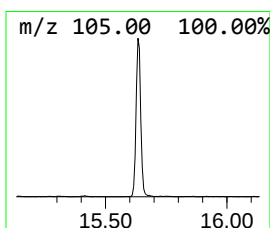
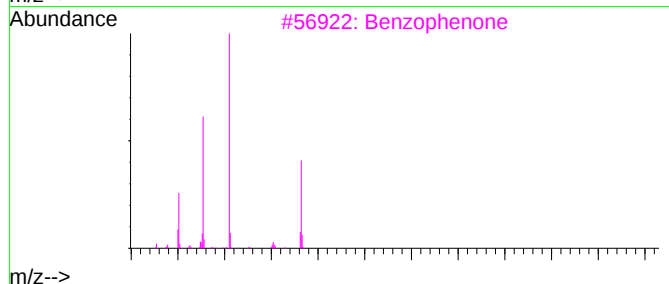
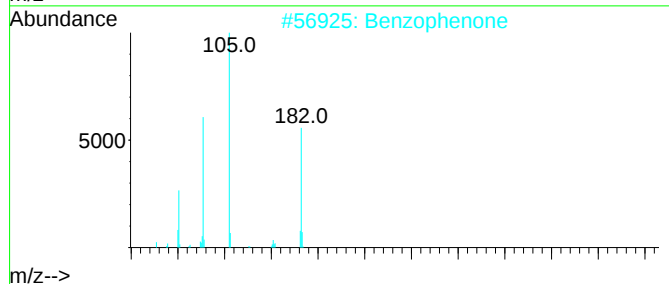
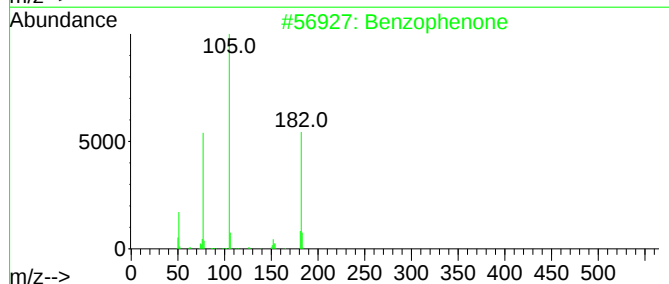
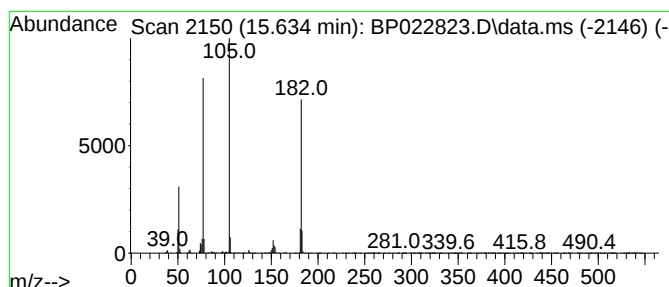
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.634	5.75 ng/ul	271784	Acenaphthene-d10	14.257

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	93
5	Benzophenone	182	C13H10O	000119-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022823.D
Acq On : 04 Nov 2024 12:17
Operator : RC/JU
Sample : P4568-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EA4

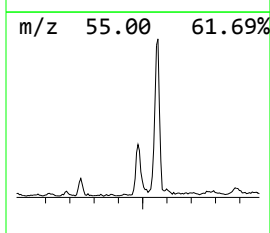
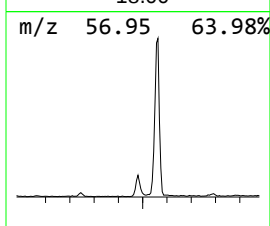
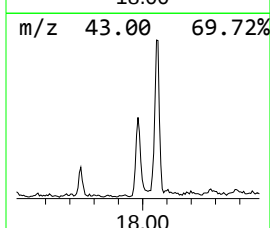
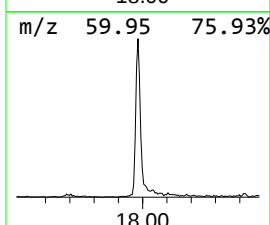
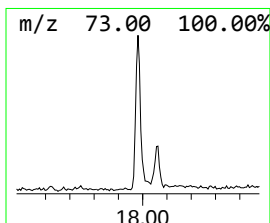
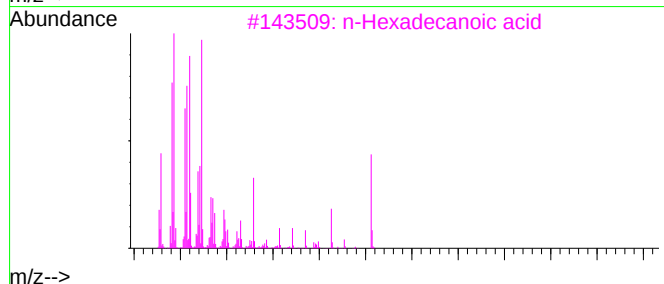
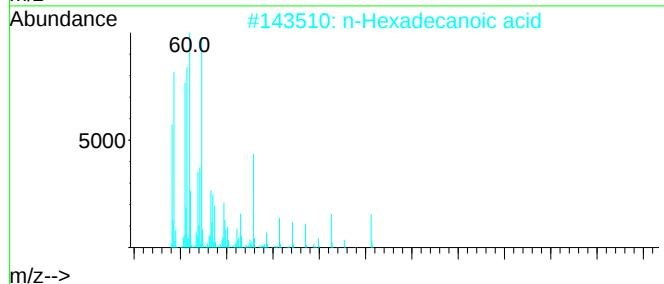
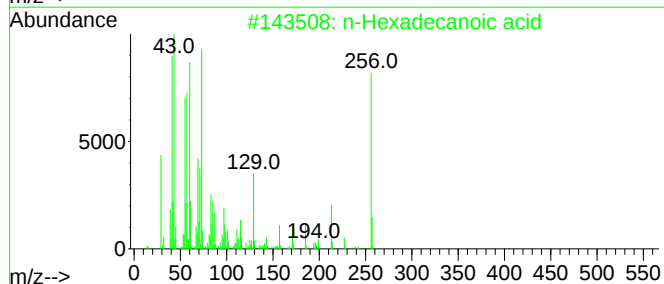
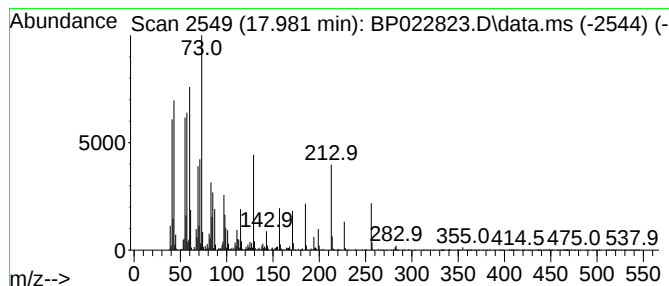
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.981	3.58 ng/ul	246785	Phenanthrene-d10	17.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022823.D
Acq On : 04 Nov 2024 12:17
Operator : RC/JU
Sample : P4568-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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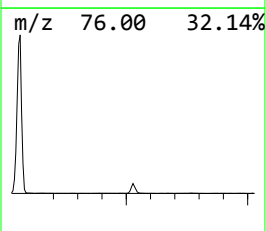
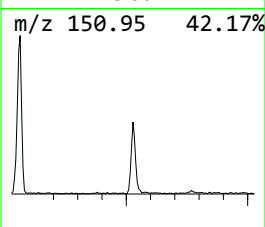
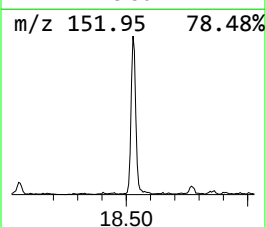
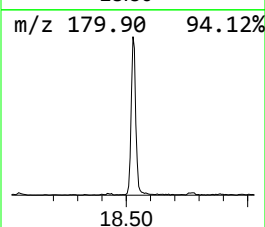
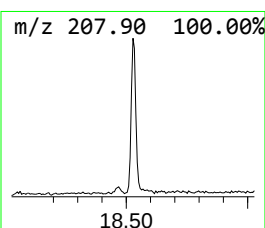
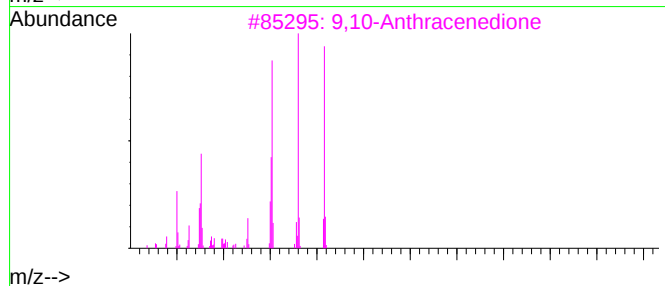
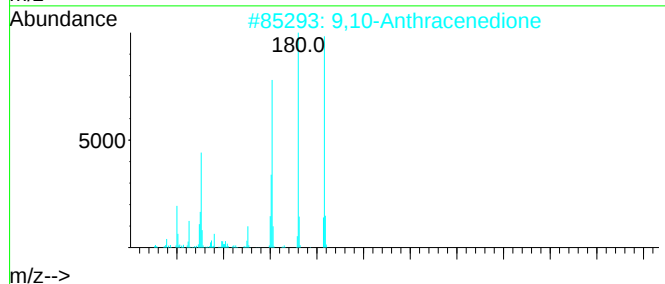
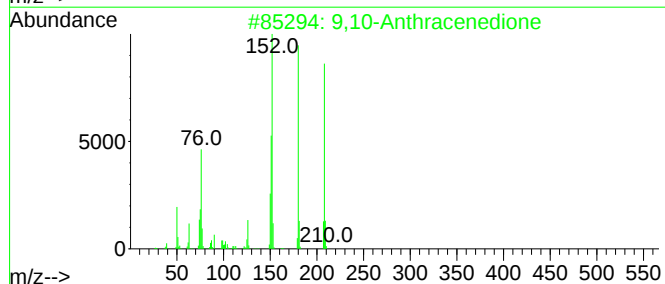
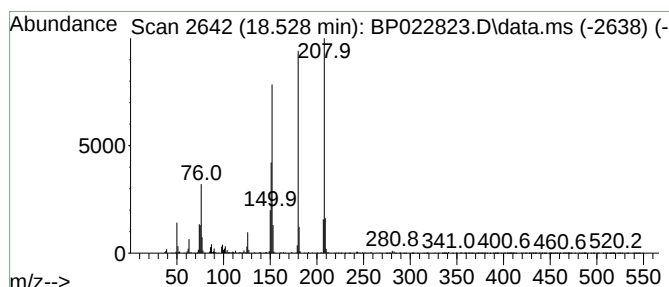
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.528	2.99 ng/ul	205692	Phenanthrene-d10	17.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	91
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	89
5	1H-Phenalen-1-one	180	C13H8O	000548-39-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022823.D
Acq On : 04 Nov 2024 12:17
Operator : RC/JU
Sample : P4568-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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
Benzene, 1,3-di...	7.975	20.1	ng/u1	2989180	1	7.628	2977480	20.0
2,6,6-Trimethyl...	9.640	3.4	ng/u1	107304	2	10.381	624456	20.0
Benzene, pentac...	14.581	2.8	ng/u1	133428	3	14.257	945204	20.0
1,1'-Biphenyl, ...	15.575	2.0	ng/u1	96989	3	14.257	945204	20.0
Benzophenone	15.634	5.8	ng/u1	271784	3	14.257	945204	20.0
n-Hexadecanoic ...	17.981	3.6	ng/u1	246785	4	17.057	1377910	20.0
9,10-Anthracene...	18.528	3.0	ng/u1	205692	4	17.057	1377910	20.0