

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
 Data File : BN015108.D
 Acq On : 26 Jun 2021 09:50
 Operator : CG/JU
 Sample : M2680-02DL 5X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGD72DL

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_N\METHODS\SFAM-EPA-BN061821MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.999	150	155	163	rBV	21122	29777	0.64%	0.064%
2	5.410	390	395	402	rBV	22984	41595	0.89%	0.089%
3	6.893	639	647	656	rVB	66619	116543	2.50%	0.250%
4	7.063	671	676	685	rVB	55228	86897	1.86%	0.187%
5	7.257	704	709	716	rBV	70139	108735	2.33%	0.234%
6	7.722	781	788	798	rVB	661414	1052157	22.55%	2.261%
7	8.416	900	906	913	rVB	75844	117611	2.52%	0.253%
8	8.869	977	983	993	rVB	57259	93363	2.00%	0.201%
9	9.581	1099	1104	1115	rVB	35109	55870	1.20%	0.120%
10	10.116	1188	1195	1201	rVB	67932	110002	2.36%	0.236%
11	10.498	1252	1260	1266	rBV	890324	1488473	31.91%	3.198%
12	10.628	1277	1282	1288	rBV	55541	91552	1.96%	0.197%
13	13.757	1810	1814	1820	rVB	123692	169063	3.62%	0.363%
14	14.039	1854	1862	1872	rBV	153907	249893	5.36%	0.537%
15	14.351	1908	1915	1921	rBV	1283622	1886233	40.43%	4.053%
16	14.410	1921	1925	1932	rVB	33310	51613	1.11%	0.111%
17	14.539	1942	1947	1957	rBV3	20841	37429	0.80%	0.080%
18	14.751	1978	1983	1991	rBV	35903	56506	1.21%	0.121%
19	15.345	2075	2084	2089	rVV	204879	296029	6.35%	0.636%
20	15.398	2089	2093	2100	rVV3	53830	75761	1.62%	0.163%
21	15.775	2152	2157	2163	rBV	32154	44943	0.96%	0.097%
22	15.845	2163	2169	2176	rVB2	20767	39663	0.85%	0.085%
23	16.410	2253	2265	2269	rBV6	19602	40722	0.87%	0.087%
24	16.751	2319	2323	2329	rVV3	26177	44366	0.95%	0.095%
25	16.833	2331	2337	2342	rVV3	18164	36387	0.78%	0.078%
26	16.916	2347	2351	2361	rVB	42248	64490	1.38%	0.139%
27	17.098	2376	2382	2386	rBV	1527537	2246405	48.15%	4.826%
28	17.139	2386	2389	2394	rVV	885960	1134545	24.32%	2.438%
29	17.192	2394	2398	2401	rVV	237088	347585	7.45%	0.747%
30	17.228	2401	2404	2410	rVB	258419	351934	7.54%	0.756%
31	17.286	2410	2414	2420	rVB	35492	55005	1.18%	0.118%
32	17.498	2441	2450	2455	rBV2	110252	188020	4.03%	0.404%
33	17.975	2525	2531	2534	rVV	108354	156291	3.35%	0.336%
34	18.022	2534	2539	2547	rVB	191832	293914	6.30%	0.631%
35	18.098	2547	2552	2557	rBV	92889	141426	3.03%	0.304%
36	18.169	2558	2564	2568	rBV3	149635	308265	6.61%	0.662%

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Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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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_N\METHODS\SFAM-EPA-BN061821MA.M
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37	18.204	2568	2570	2575	rVB2	102432	118191	2.53%	0.254%
38	18.316	2585	2589	2594	rBV4	24108	42565	0.91%	0.091%
39	18.422	2602	2607	2611	rBV2	43958	72702	1.56%	0.156%
40	18.480	2612	2617	2620	rVV	88812	129496	2.78%	0.278%
41	18.516	2620	2623	2626	rVB2	27796	34391	0.74%	0.074%
42	18.727	2655	2659	2662	rBV3	47558	65736	1.41%	0.141%
43	18.775	2664	2667	2670	rVB2	26457	32602	0.70%	0.070%
44	18.816	2670	2674	2678	rVB2	41554	59452	1.27%	0.128%
45	18.898	2682	2688	2692	rBV	84616	161892	3.47%	0.348%
46	18.963	2693	2699	2702	rBV4	80404	138631	2.97%	0.298%
47	19.033	2708	2711	2719	rVB2	63588	92430	1.98%	0.199%
48	19.110	2719	2724	2727	rBV6	34748	57929	1.24%	0.124%
49	19.163	2727	2733	2739	rVB	1829742	2532412	54.29%	5.441%
50	19.257	2746	2749	2753	rVB3	58232	74795	1.60%	0.161%
51	19.404	2771	2774	2778	rVB2	43215	51033	1.09%	0.110%
52	19.498	2784	2790	2791	rBV2	593805	776322	16.64%	1.668%
53	19.522	2791	2794	2803	rVV	1511245	2228727	47.78%	4.788%
54	19.598	2803	2807	2814	rVB2	108514	196792	4.22%	0.423%
55	19.704	2821	2825	2831	rBV	125516	170337	3.65%	0.366%
56	19.763	2831	2835	2841	rVB	101290	156932	3.36%	0.337%
57	19.851	2845	2850	2857	rBV4	130045	334641	7.17%	0.719%
58	19.986	2869	2873	2876	rBV3	85029	145072	3.11%	0.312%
59	20.021	2876	2879	2883	rVB	216094	276489	5.93%	0.594%
60	20.063	2883	2886	2890	rBV	104399	109969	2.36%	0.236%
61	20.121	2893	2896	2900	rBV	114636	142922	3.06%	0.307%
62	20.180	2900	2906	2911	rVB2	194876	325797	6.98%	0.700%
63	20.263	2917	2920	2926	rBV2	46492	65543	1.41%	0.141%
64	20.321	2926	2930	2934	rBV2	91461	122779	2.63%	0.264%
65	20.369	2934	2938	2943	rBV2	97208	171575	3.68%	0.369%
66	20.516	2959	2963	2968	rVB	892767	944832	20.25%	2.030%
67	20.563	2968	2971	2976	rBV2	163767	185760	3.98%	0.399%
68	20.739	2997	3001	3003	rBV5	35539	58785	1.26%	0.126%
69	20.774	3003	3007	3012	rVV	137952	199530	4.28%	0.429%
70	20.939	3029	3035	3039	rBV3	147935	277490	5.95%	0.596%
71	20.980	3039	3042	3045	rVV	196267	247048	5.30%	0.531%
72	21.027	3045	3050	3054	rVV3	447010	649131	13.92%	1.395%
73	21.068	3054	3057	3064	rVB2	133751	211155	4.53%	0.454%
74	21.221	3081	3083	3085	rBV2	40703	30922	0.66%	0.066%
75	21.245	3085	3087	3089	rVV	44618	51079	1.09%	0.110%
76	21.292	3089	3095	3099	rVV3	2359227	4664963	100.00%	10.023%

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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

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
 Title : SVOA CALIBRATION

77	21.333	3099	3102	3111	rVB	1321531	1718285	36.83%	3.692%
78	21.463	3118	3124	3128	rBV2	372240	584001	12.52%	1.255%
79	21.610	3144	3149	3155	rBV2	200932	328274	7.04%	0.705%
80	21.792	3177	3180	3183	rBV	63387	95065	2.04%	0.204%
81	21.845	3186	3189	3194	rVV	250758	368591	7.90%	0.792%
82	21.904	3195	3199	3203	rVB	443071	543802	11.66%	1.168%
83	22.045	3219	3223	3226	rBV	117551	168458	3.61%	0.362%
84	22.080	3226	3229	3235	rVB3	109643	163538	3.51%	0.351%
85	22.151	3237	3241	3250	rVB4	97655	201950	4.33%	0.434%
86	22.368	3274	3278	3286	rVB	332134	491516	10.54%	1.056%
87	22.874	3357	3364	3369	rBV2	1265880	2966443	63.59%	6.373%
88	23.045	3388	3393	3400	rVB	249776	429878	9.22%	0.924%
89	23.174	3412	3415	3417	rBV2	56542	78561	1.68%	0.169%
90	23.351	3439	3445	3450	rBV	569074	1056633	22.65%	2.270%
91	23.398	3450	3453	3456	rVV2	160354	268244	5.75%	0.576%
92	23.445	3456	3461	3469	rVB	1215104	2228466	47.77%	4.788%
93	23.545	3471	3478	3483	rBV2	1273563	2475273	53.06%	5.318%
94	23.592	3483	3486	3495	rVB2	327306	582928	12.50%	1.252%
95	24.104	3567	3573	3584	rVB2	224337	455191	9.76%	0.978%
96	24.857	3695	3701	3707	rBV2	141382	289311	6.20%	0.622%
97	25.733	3843	3850	3853	rBV3	95189	244642	5.24%	0.526%
98	25.798	3854	3861	3874	rVB2	514096	1500401	32.16%	3.224%
99	26.062	3902	3906	3914	rVB	103385	229224	4.91%	0.492%
100	26.486	3970	3978	3995	rVB2	281338	957551	20.53%	2.057%

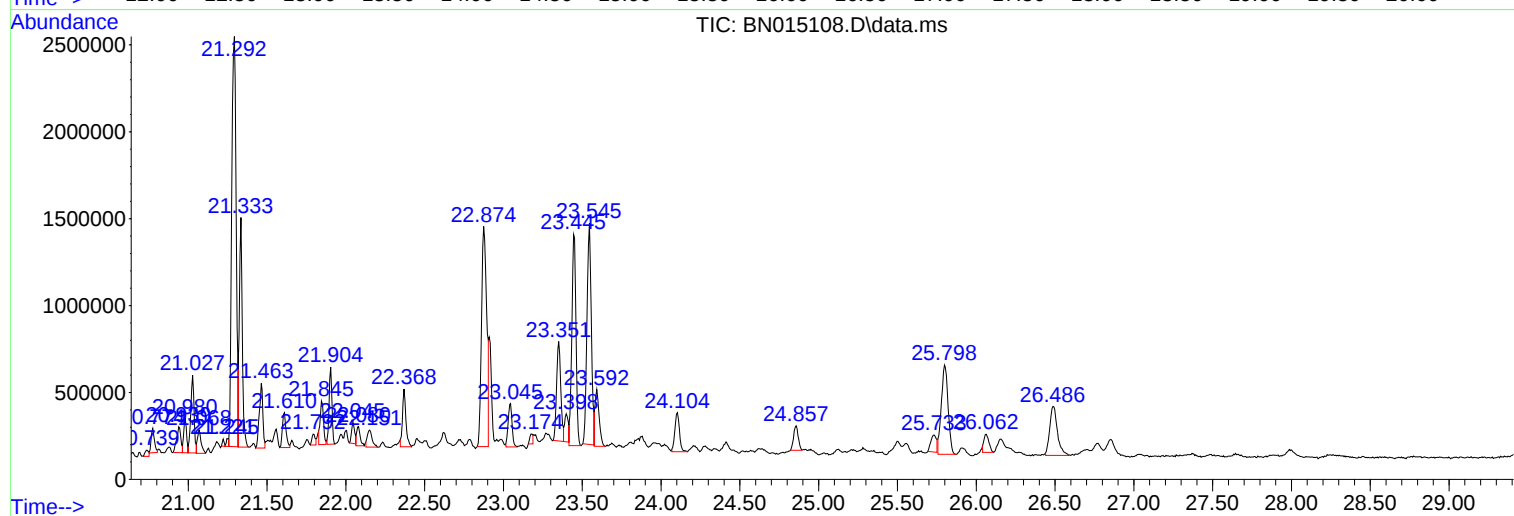
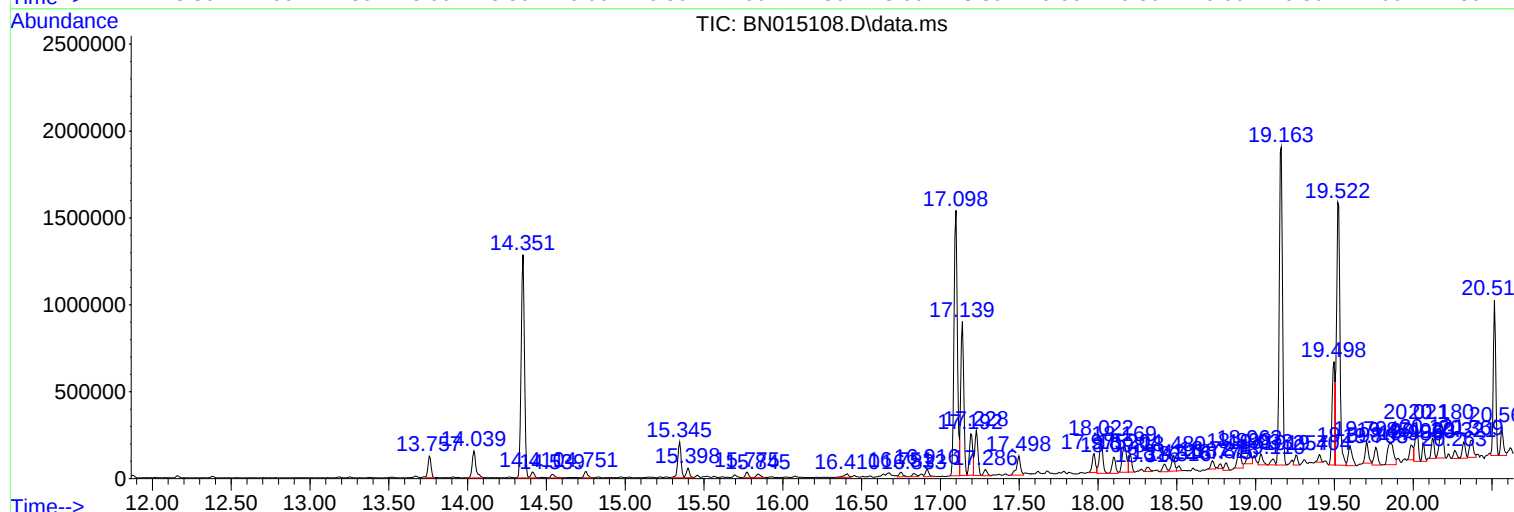
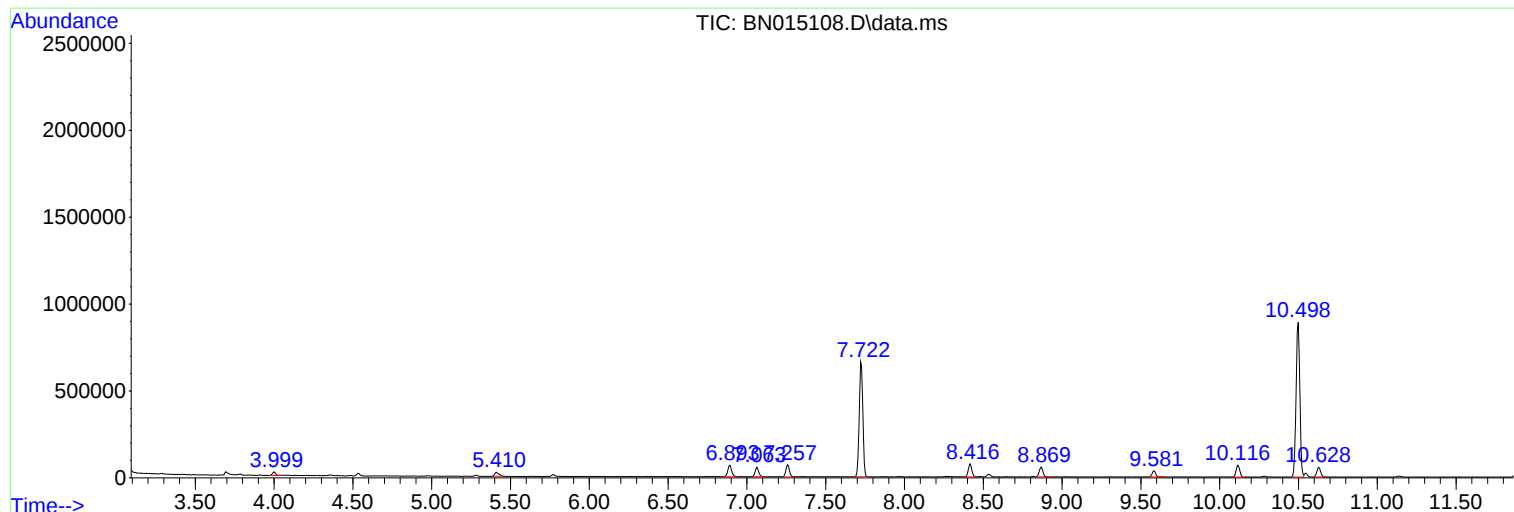
Sum of corrected areas: 46544138

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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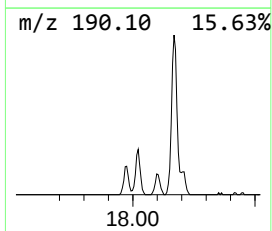
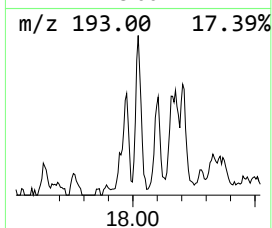
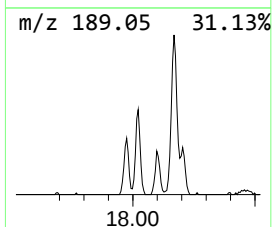
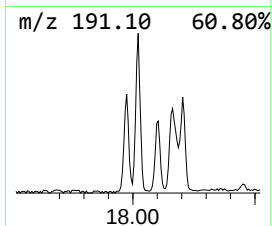
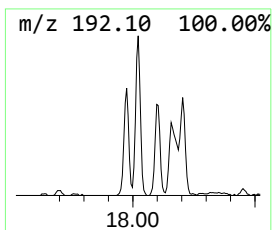
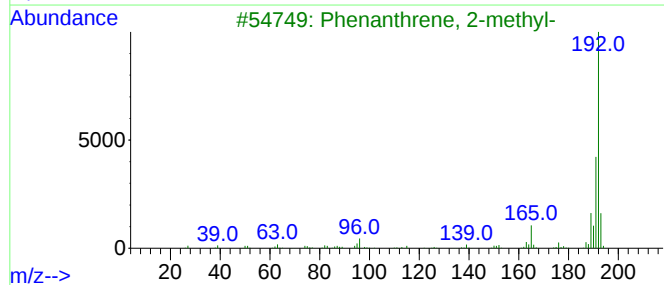
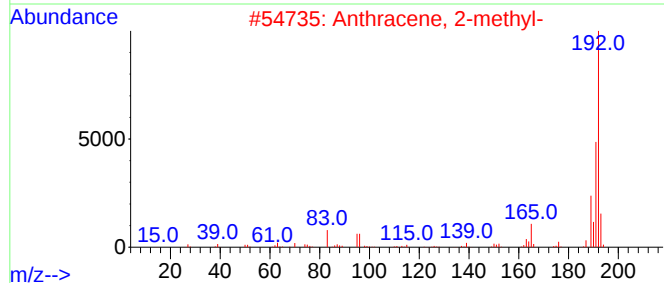
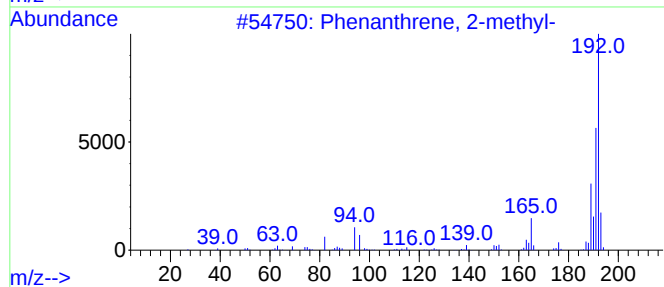
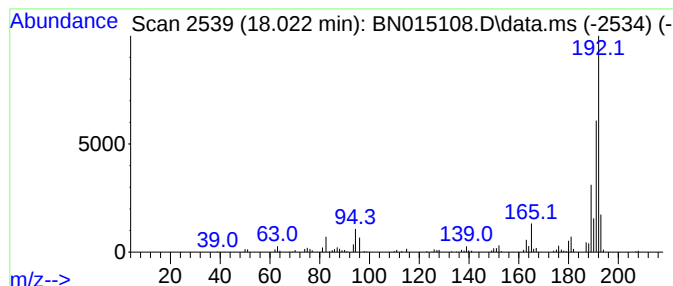
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	2.62 ng/ul	293914	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	96
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



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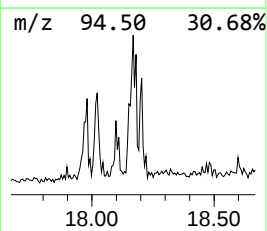
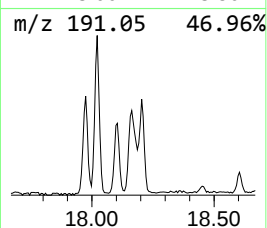
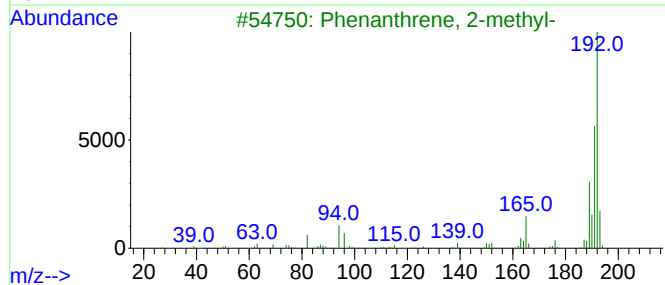
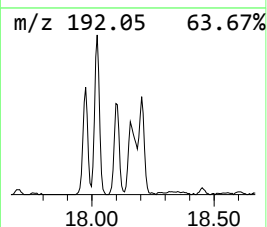
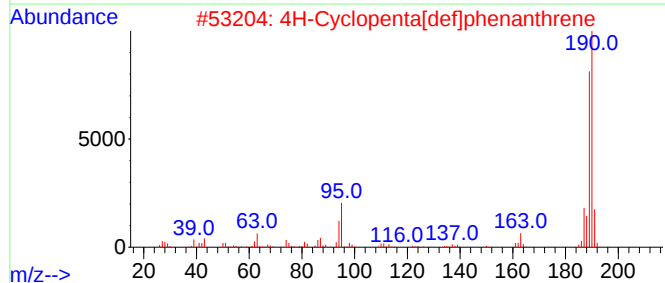
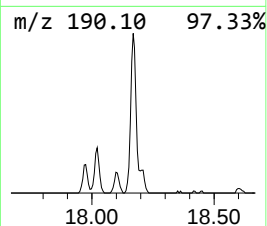
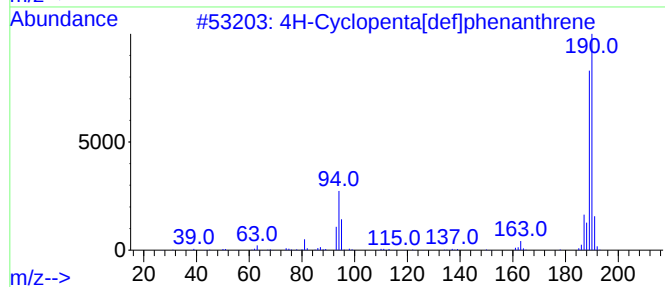
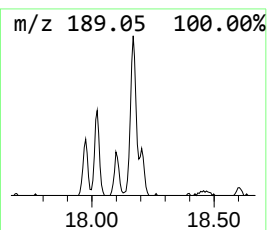
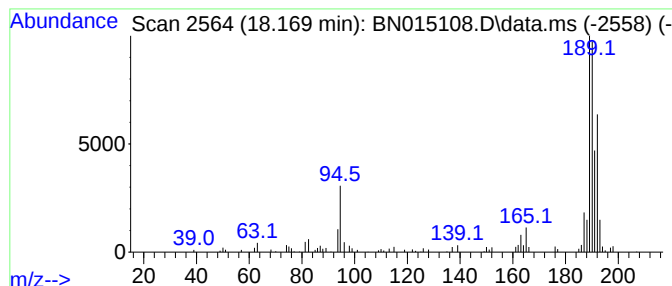
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.169	2.74 ng/ul	308265	Phenanthrene-d10	17.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	42
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	38
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38



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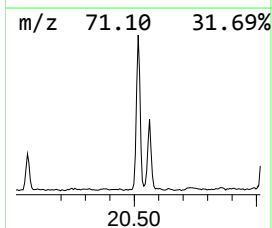
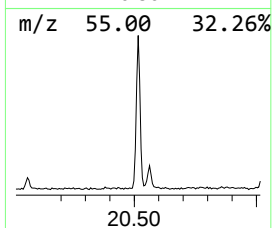
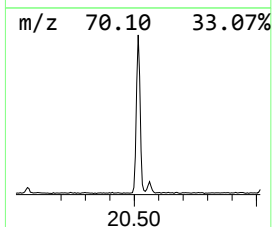
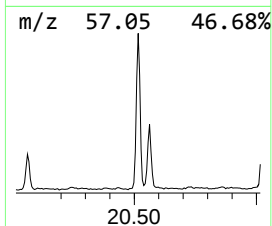
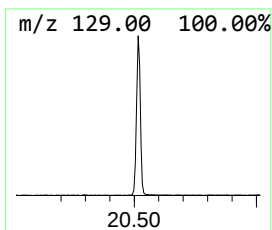
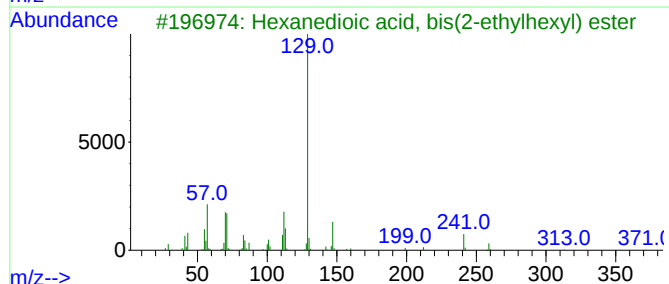
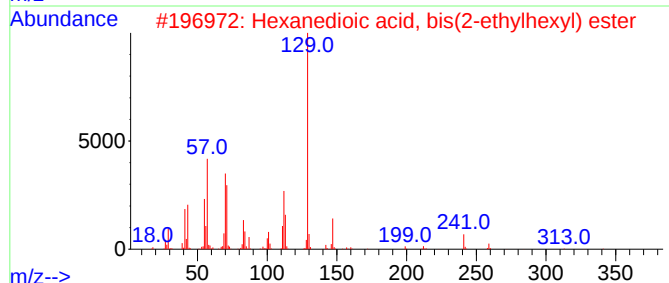
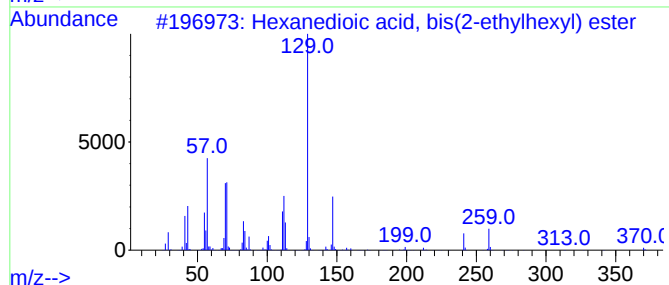
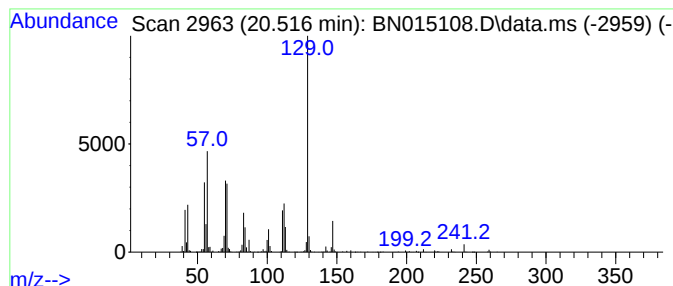
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexanedioic acid, bis(2-eth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.516	4.05 ng/ul	944832	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
4			Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	83
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
Operator : CG/JU
Sample : M2680-02DL 5X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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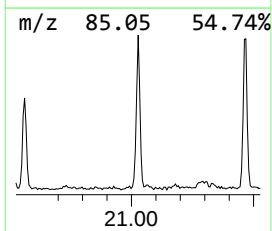
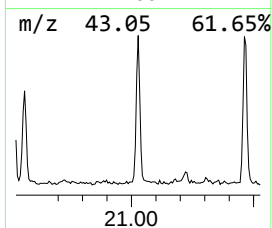
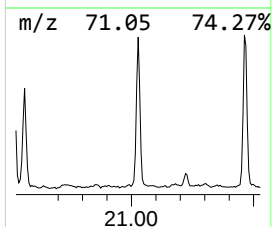
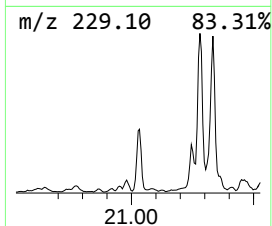
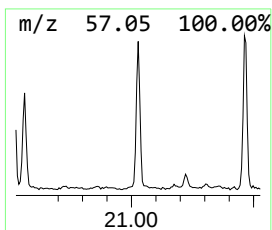
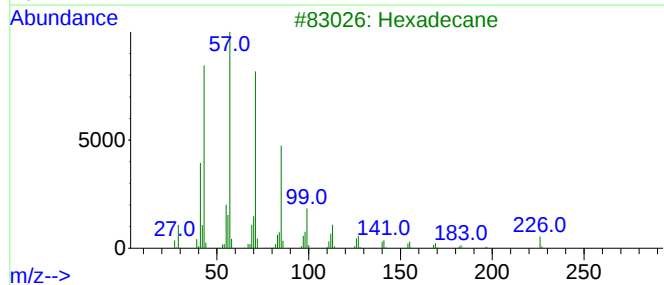
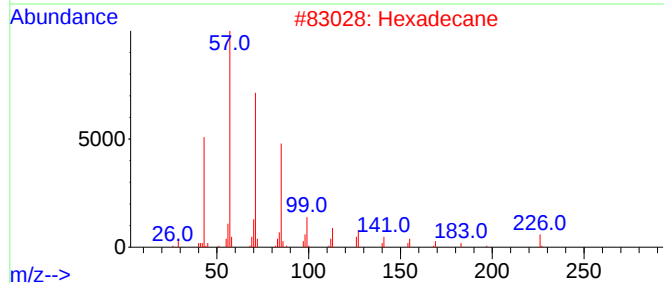
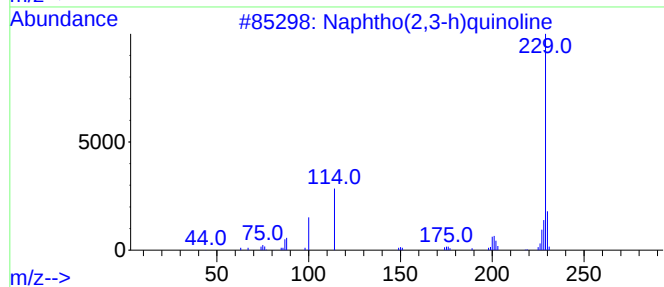
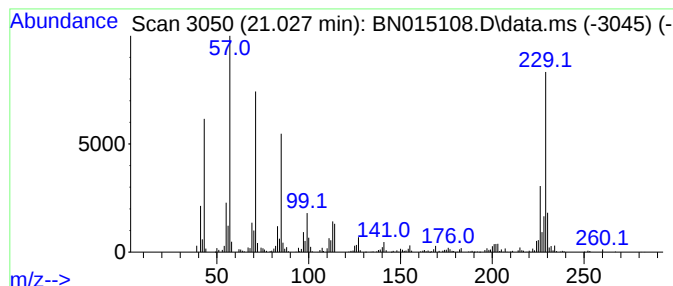
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphtho(2,3-h)quinoline Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.027	2.78 ng/ul	649131	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho(2,3-h)quinoline	229	C17H11N	000084-56-0	60
2			Hexadecane	226	C16H34	000544-76-3	49
3			Hexadecane	226	C16H34	000544-76-3	49
4			Hexadecane	226	C16H34	000544-76-3	49
5			Heneicosane	296	C21H44	000629-94-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
Operator : CG/JU
Sample : M2680-02DL 5X
Misc :
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Instrument :
BNA_N
ClientSampleId :
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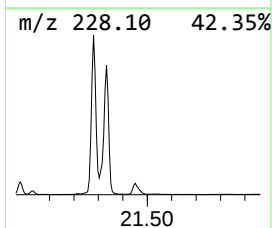
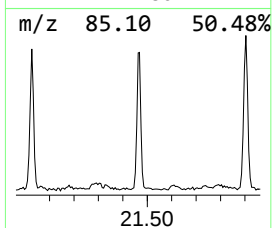
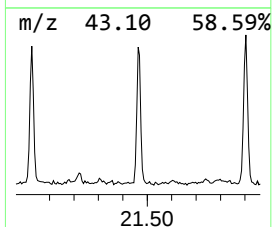
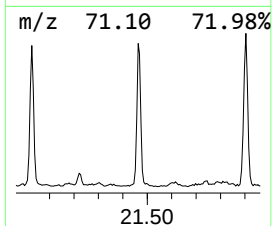
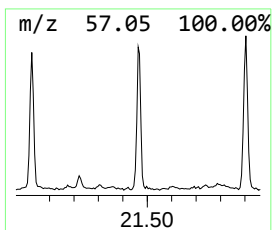
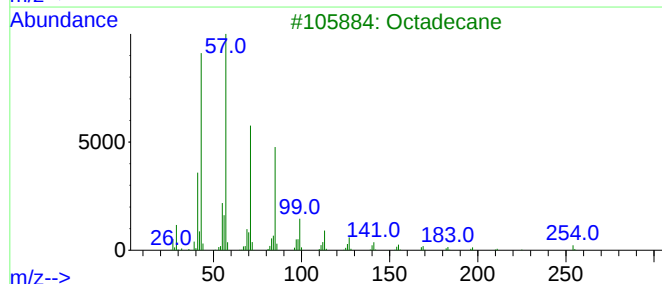
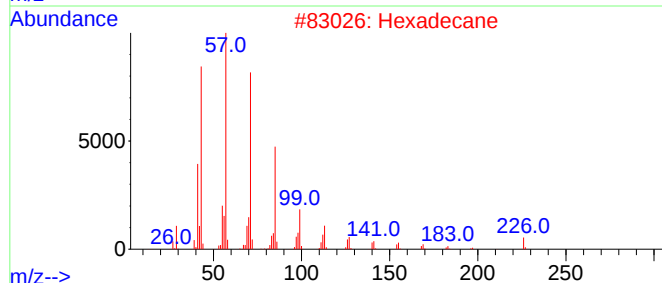
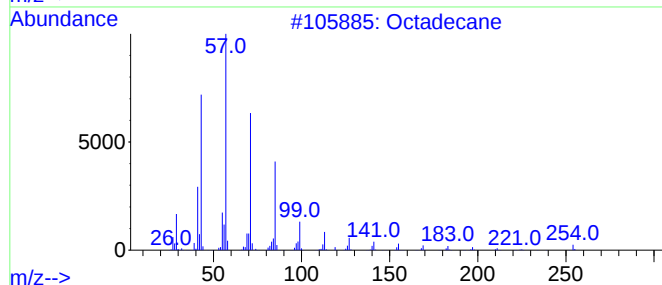
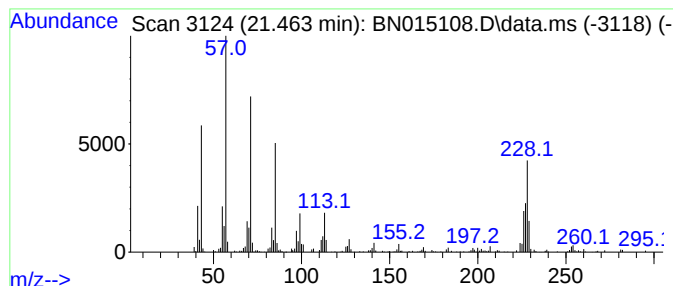
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.463	2.50 ng/ul	584001	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	95
2			Hexadecane	226	C16H34	000544-76-3	87
3			Octadecane	254	C18H38	000593-45-3	64
4			Heneicosane	296	C21H44	000629-94-7	50
5			Decane, 1-iodo-	268	C10H21I	002050-77-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
Operator : CG/JU
Sample : M2680-02DL 5X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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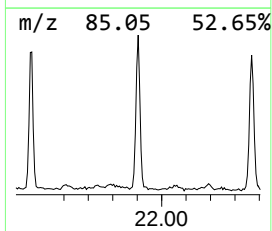
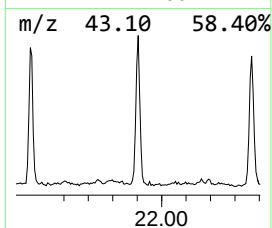
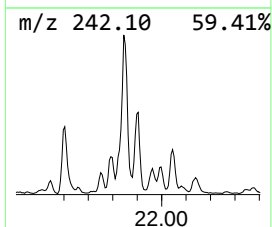
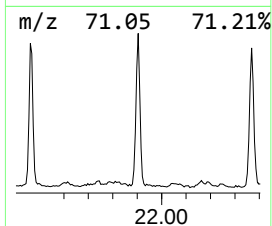
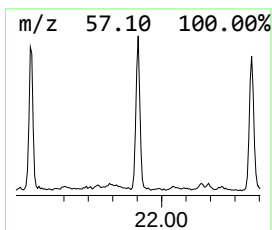
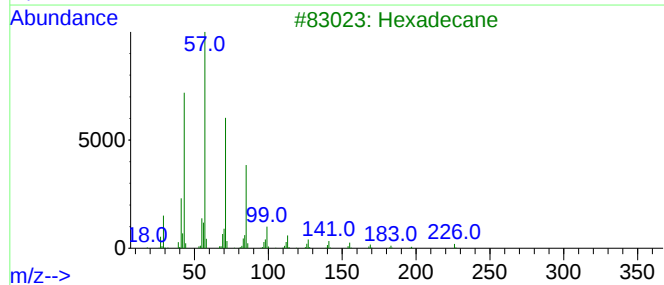
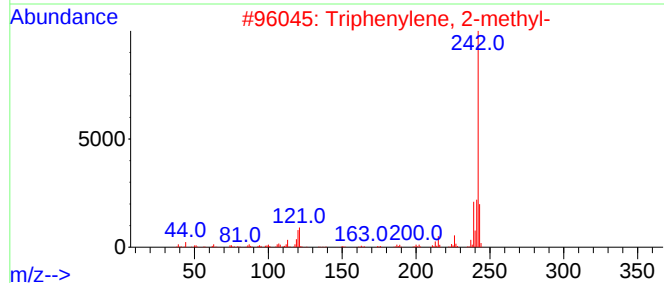
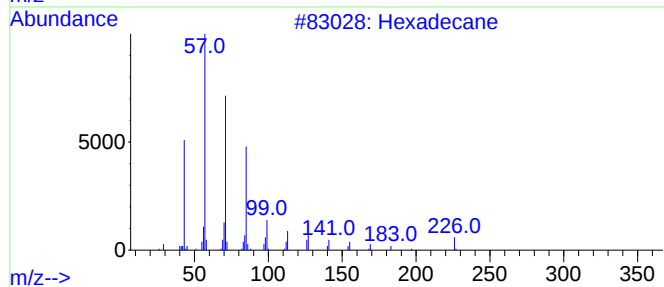
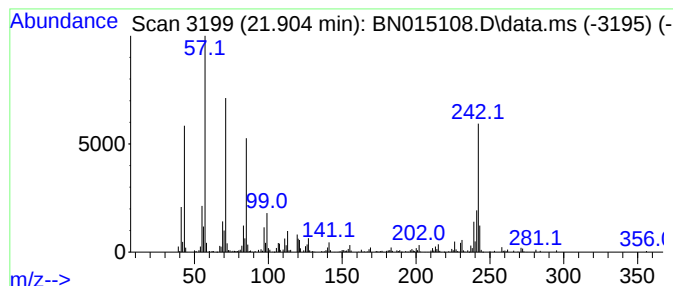
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.904	2.33 ng/ul	543802	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	93
2			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	92
3			Hexadecane	226	C16H34	000544-76-3	89
4			Hexadecane	226	C16H34	000544-76-3	89
5			Heptadecane	240	C17H36	000629-78-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
Operator : CG/JU
Sample : M2680-02DL 5X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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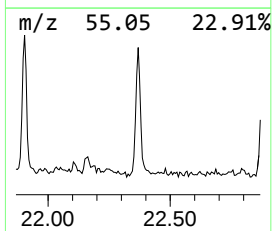
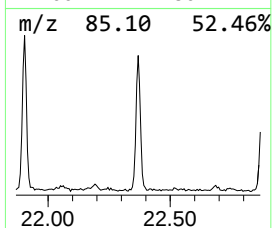
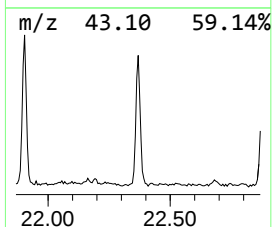
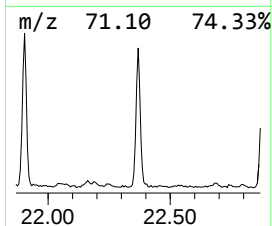
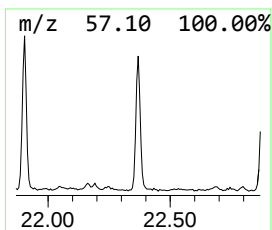
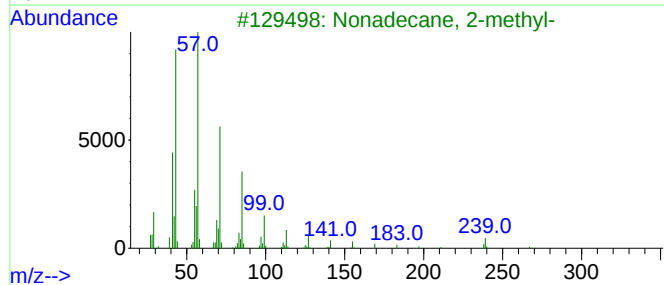
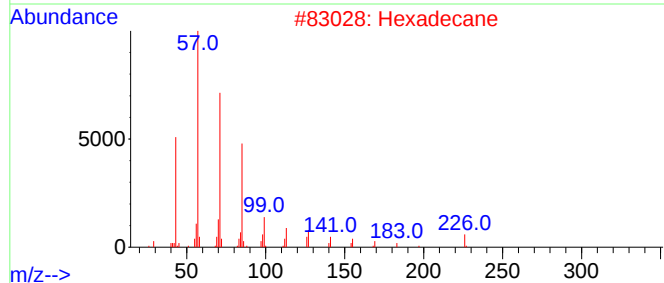
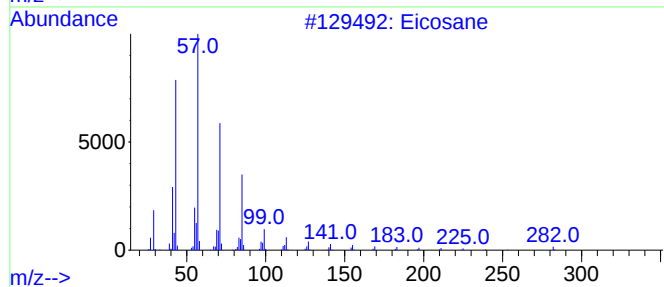
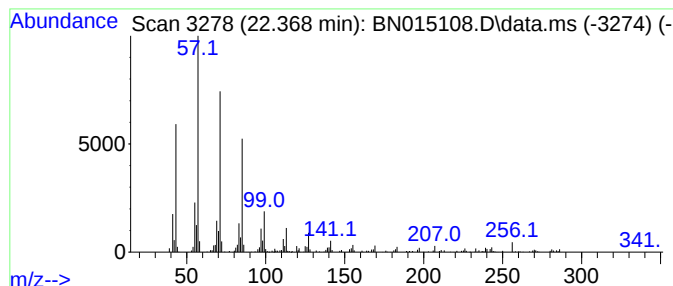
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.368	2.11 ng/ul	491516	Chrysene-d12	21.298

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	95
2		Hexadecane	226	C16H34	000544-76-3	91
3		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
Operator : CG/JU
Sample : M2680-02DL 5X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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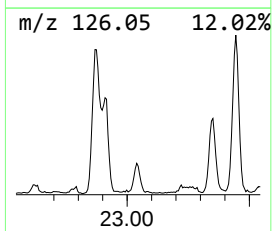
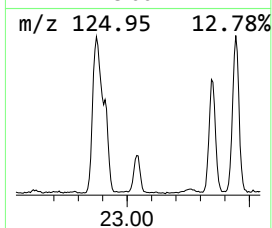
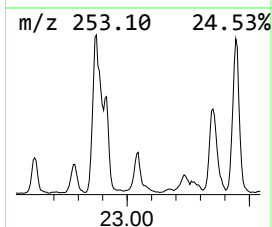
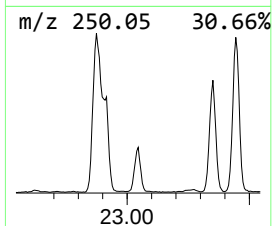
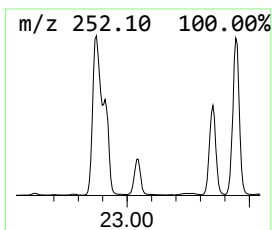
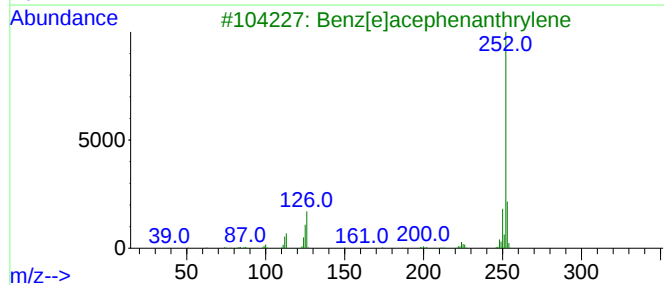
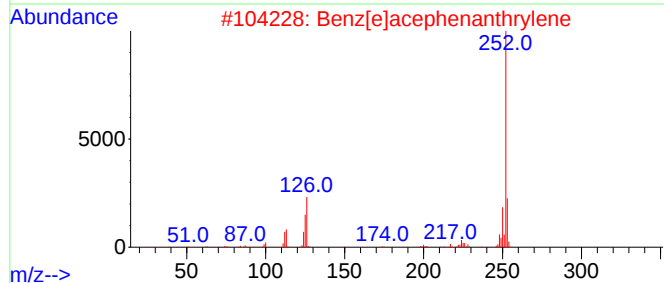
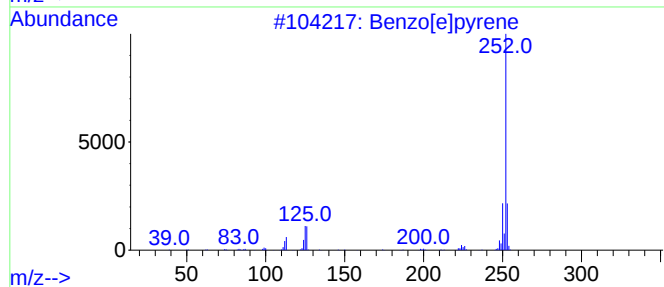
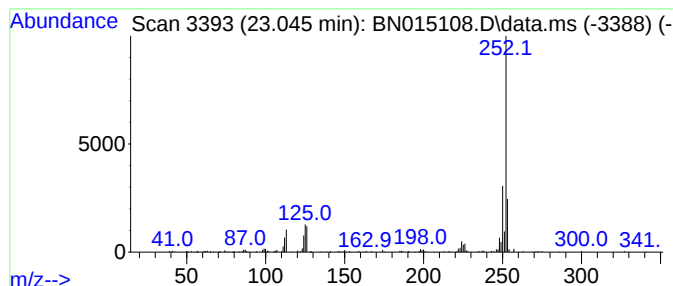
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.045	3.47 ng/ul	429878	Perylene-d12	23.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	94
2			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
3			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
5			Benzo[a]pyrene	252	C20H12	000050-32-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
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Instrument :
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ClientSampleId :
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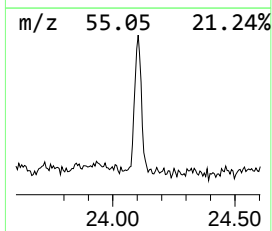
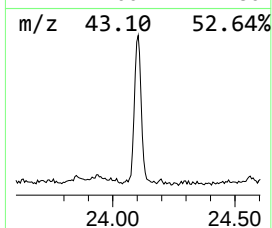
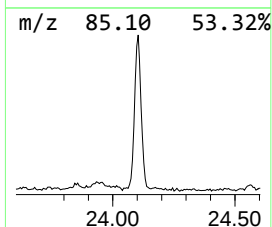
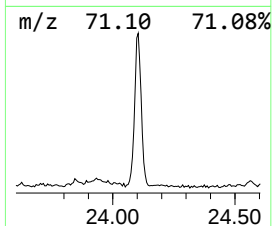
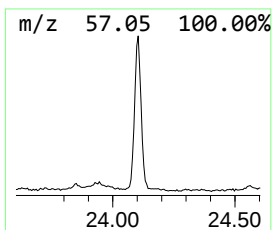
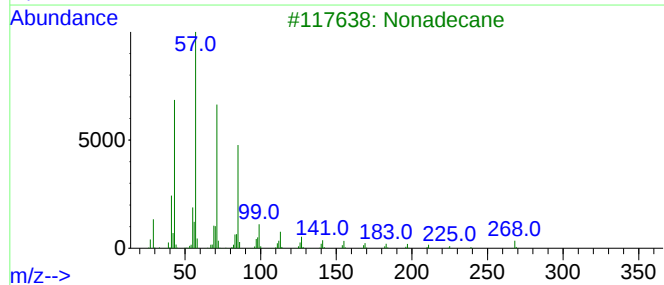
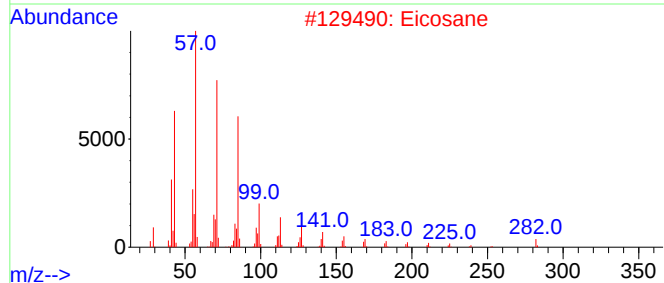
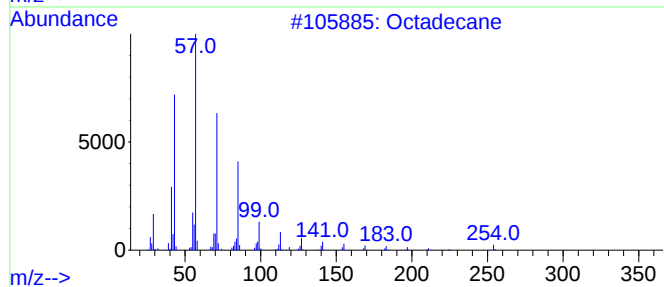
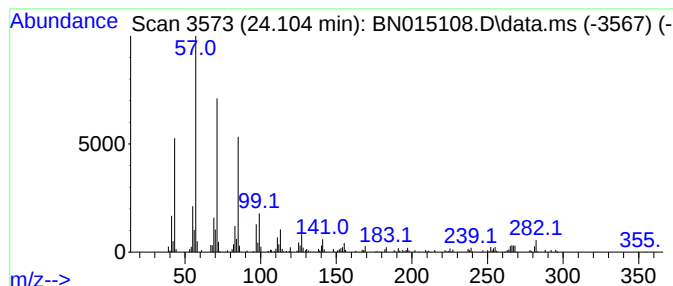
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.104	3.68 ng/ul	455191	Perylene-d12	23.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	97
2			Eicosane	282	C20H42	000112-95-8	97
3			Nonadecane	268	C19H40	000629-92-5	96
4			Eicosane	282	C20H42	000112-95-8	93
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
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ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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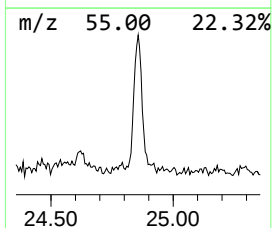
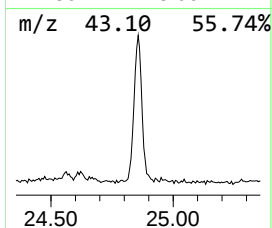
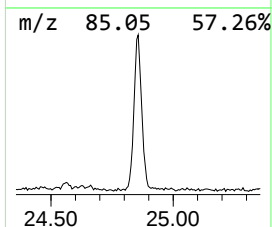
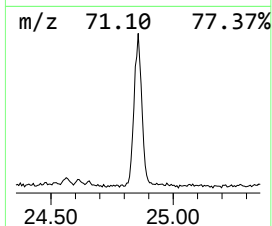
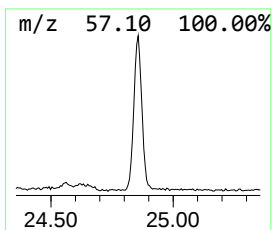
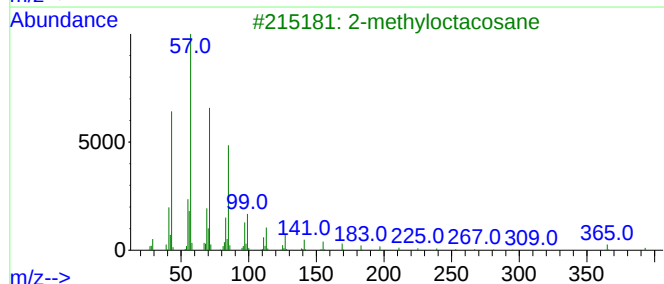
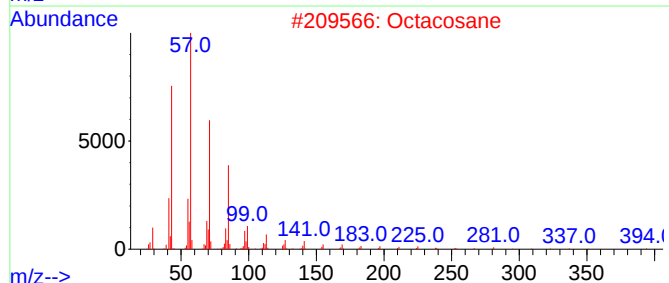
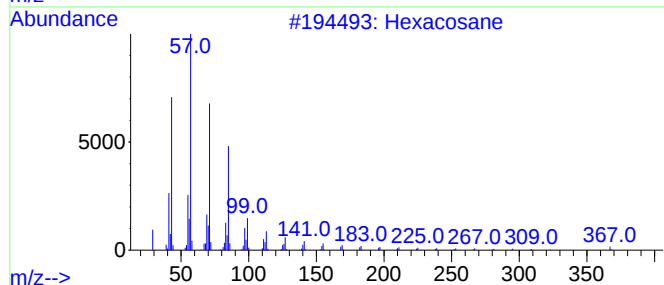
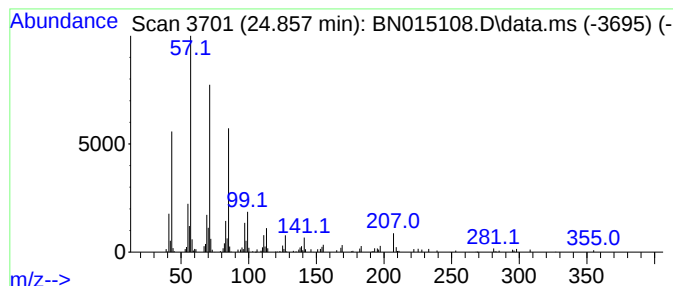
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.857	2.34 ng/ul	289311	Perylene-d12	23.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Octacosane	394	C28H58	000630-02-4	91
3			2-methyloctacosane	408	C29H60	1000376-72-8	91
4			Triacontane	422	C30H62	000638-68-6	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015108.D
Acq On : 26 Jun 2021 09:50
Operator : CG/JU
Sample : M2680-02DL 5X
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGD72DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 2...	18.022	2.6	ng/u1	293914	4	17.098	2246410	20.0
4H-Cyclopenta[d...	18.169	2.7	ng/u1	308265	4	17.098	2246410	20.0
Hexanedioic aci...	20.516	4.0	ng/u1	944832	5	21.298	4664960	20.0
Naphtho(2,3-h)q...	21.027	2.8	ng/u1	649131	5	21.298	4664960	20.0
(DEL) Alkane: S...	21.463	2.5	ng/u1	584001	5	21.298	4664960	20.0
(DEL) Alkane: S...	21.904	2.3	ng/u1	543802	5	21.298	4664960	20.0
(DEL) Alkane: S...	22.368	2.1	ng/u1	491516	5	21.298	4664960	20.0
Benzo[e]pyrene	23.045	3.5	ng/u1	429878	6	23.545	2475270	20.0
(DEL) Alkane: S...	24.104	3.7	ng/u1	455191	6	23.545	2475270	20.0
(DEL) Alkane: S...	24.857	2.3	ng/u1	289311	6	23.545	2475270	20.0