

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053124\
 Data File : BG061406.D
 Acq On : 1 Jun 2024 00:21
 Operator : MA/JU
 Sample : P2588-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH5S5

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_G\Methods\SFAM-EPA-BG052124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061406.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.082	10	16	35	rVB	118989	237088	13.64%	2.144%
2	3.334	55	59	68	rVB4	3971	6424	0.37%	0.058%
3	3.605	100	105	115	rBV	48870	80753	4.65%	0.730%
4	3.740	123	128	141	rVB	46803	85495	4.92%	0.773%
5	4.733	294	297	304	rBV3	3178	6438	0.37%	0.058%
6	7.060	687	693	701	rBV	70257	120985	6.96%	1.094%
7	7.201	710	717	724	rVB	44826	70017	4.03%	0.633%
8	7.412	747	753	758	rBV	66767	110164	6.34%	0.996%
9	7.459	758	761	766	rVB	18907	28152	1.62%	0.255%
10	7.870	823	831	838	rBV	109976	186318	10.72%	1.685%
11	8.587	946	953	964	rBV2	74053	128760	7.41%	1.164%
12	9.028	1020	1028	1036	rBB	65726	114801	6.61%	1.038%
13	9.586	1117	1123	1128	rBV	9549	14977	0.86%	0.135%
14	9.750	1144	1151	1158	rVB	63720	103841	5.98%	0.939%
15	10.303	1239	1245	1252	rBV	90084	154843	8.91%	1.400%
16	10.438	1261	1268	1277	rBV2	14598	27918	1.61%	0.252%
17	10.667	1299	1307	1314	rBV	145205	255615	14.71%	2.311%
18	10.802	1324	1330	1341	rVB2	71549	127549	7.34%	1.153%
19	11.196	1392	1397	1402	rVB3	3717	5155	0.30%	0.047%
20	11.407	1426	1433	1446	rVB2	36001	64707	3.72%	0.585%
21	13.910	1851	1859	1866	rVB	175497	252062	14.51%	2.279%
22	14.198	1901	1908	1917	rVB	180843	288332	16.59%	2.607%
23	14.510	1950	1961	1967	rVB	240060	377785	21.74%	3.416%
24	14.733	1989	1999	2012	rBV	752610	1737573	100.00%	15.710%
25	14.815	2012	2013	2021	rVV4	2839	4465	0.26%	0.040%
26	15.503	2124	2130	2136	rBV	262694	391701	22.54%	3.542%
27	15.638	2148	2153	2160	rBV	89843	126625	7.29%	1.145%
28	16.237	2249	2255	2259	rBV5	3334	5670	0.33%	0.051%
29	16.913	2365	2370	2374	rBV3	4917	7571	0.44%	0.068%
30	17.259	2422	2429	2433	rVV	344429	506373	29.14%	4.578%
31	17.300	2433	2436	2440	rVV	78368	108154	6.22%	0.978%
32	17.353	2440	2445	2456	rVB2	333483	502407	28.91%	4.542%
33	17.453	2457	2462	2467	rVB4	3444	4669	0.27%	0.042%
34	17.659	2492	2497	2508	rVB5	7011	13221	0.76%	0.120%
35	17.841	2523	2528	2532	rBV5	3046	4945	0.28%	0.045%
36	17.929	2538	2543	2546	rBV2	5750	7609	0.44%	0.069%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
 Title : SVOA CALIBRATION

37	18.135	2573	2578	2580	rBV2	16509	24407	1.40%	0.221%
38	18.182	2583	2586	2592	rVV	23745	34236	1.97%	0.310%
39	18.264	2597	2600	2605	rBV3	3742	5533	0.32%	0.050%
40	18.329	2605	2611	2615	rBV4	20360	35016	2.02%	0.317%
41	18.364	2615	2617	2623	rVB3	10730	13642	0.79%	0.123%
42	18.446	2627	2631	2635	rBV5	4443	7106	0.41%	0.064%
43	18.528	2640	2645	2650	rVB6	3200	4925	0.28%	0.045%
44	18.634	2659	2663	2666	rBV	12313	16176	0.93%	0.146%
45	18.675	2666	2670	2678	rVB2	12947	20892	1.20%	0.189%
46	18.881	2702	2705	2711	rVB5	5048	6707	0.39%	0.061%
47	18.934	2711	2714	2716	rBV2	7159	7447	0.43%	0.067%
48	18.969	2718	2720	2724	rVB3	5050	5046	0.29%	0.046%
49	19.016	2724	2728	2731	rBV5	11936	22270	1.28%	0.201%
50	19.051	2731	2734	2739	rVB3	12622	16487	0.95%	0.149%
51	19.098	2739	2742	2748	rVB6	13662	17126	0.99%	0.155%
52	19.139	2748	2749	2752	rVB3	5300	4390	0.25%	0.040%
53	19.192	2754	2758	2759	rBV3	12047	12112	0.70%	0.110%
54	19.316	2772	2779	2786	rVB	184870	269693	15.52%	2.438%
55	19.380	2786	2790	2793	rBV5	5836	9547	0.55%	0.086%
56	19.463	2802	2804	2809	rVB3	4857	5374	0.31%	0.049%
57	19.510	2809	2812	2814	rBV4	5339	7083	0.41%	0.064%
58	19.557	2816	2820	2829	rVB3	14492	20944	1.21%	0.189%
59	19.651	2830	2836	2839	rBV	448150	686246	39.49%	6.205%
60	19.751	2849	2853	2860	rVB2	11565	18478	1.06%	0.167%
61	19.856	2867	2871	2876	rVB	8962	14062	0.81%	0.127%
62	19.915	2876	2881	2886	rBV5	11415	16160	0.93%	0.146%
63	20.003	2889	2896	2898	rBV7	16619	30672	1.77%	0.277%
64	20.168	2912	2924	2930	rBV3	32348	78241	4.50%	0.707%
65	20.273	2937	2942	2946	rBV3	16249	22839	1.31%	0.206%
66	20.332	2947	2952	2955	rVV2	21656	33873	1.95%	0.306%
67	20.409	2963	2965	2970	rVB6	5776	7203	0.41%	0.065%
68	20.473	2970	2976	2978	rBV2	16746	26627	1.53%	0.241%
69	20.514	2980	2983	2988	rVB2	19371	26646	1.53%	0.241%
70	20.638	3001	3004	3007	rVB5	5877	5520	0.32%	0.050%
71	20.685	3007	3012	3015	rVB6	7616	9938	0.57%	0.090%
72	20.726	3016	3019	3021	rBV4	5068	6381	0.37%	0.058%
73	20.843	3035	3039	3040	rBV4	5662	7576	0.44%	0.068%
74	20.926	3048	3053	3061	rVB2	24081	39765	2.29%	0.360%
75	21.096	3076	3082	3087	rBV5	17374	41005	2.36%	0.371%
76	21.143	3087	3090	3092	rVV3	14756	18428	1.06%	0.167%

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77	21.178	3092	3096	3104	rVV6	32550	88617	5.10%	0.801%
78	21.249	3106	3108	3113	rVB	18774	22980	1.32%	0.208%
79	21.507	3144	3152	3157	rBV3	409735	765855	44.08%	6.924%
80	21.554	3157	3160	3168	rVB	81269	128449	7.39%	1.161%
81	21.701	3181	3185	3190	rVB6	20207	29040	1.67%	0.263%
82	21.901	3216	3219	3225	rBV6	11130	18226	1.05%	0.165%
83	21.971	3227	3231	3234	rBV6	7435	14346	0.83%	0.130%
84	22.212	3270	3272	3276	rVB3	17409	19653	1.13%	0.178%
85	22.283	3280	3284	3291	rVB5	19246	34265	1.97%	0.310%
86	22.477	3312	3317	3321	rBV8	13203	23116	1.33%	0.209%
87	22.917	3387	3392	3393	rBV5	12659	17666	1.02%	0.160%
88	23.252	3444	3449	3450	rBV5	9757	14081	0.81%	0.127%
89	23.617	3503	3511	3518	rBV2	60531	189855	10.93%	1.717%
90	23.846	3548	3550	3552	rBV3	8585	9347	0.54%	0.085%
91	24.310	3620	3629	3633	rBV2	31832	89017	5.12%	0.805%
92	24.380	3633	3641	3648	rVV2	253553	668413	38.47%	6.043%
93	24.451	3649	3653	3661	rVB	51589	109648	6.31%	0.991%
94	24.592	3666	3677	3688	rBV	247475	698720	40.21%	6.317%
95	25.667	3858	3860	3870	rVB8	17235	35951	2.07%	0.325%
96	26.954	4073	4079	4081	rBV6	8866	17443	1.00%	0.158%
97	27.823	4225	4227	4232	rBV6	4057	5824	0.34%	0.053%
98	28.082	4262	4271	4285	rVV3	23583	106952	6.16%	0.967%
99	28.511	4336	4344	4345	rBV7	8913	22936	1.32%	0.207%
100	32.406	5005	5007	5010	rBV3	4057	4813	0.28%	0.044%

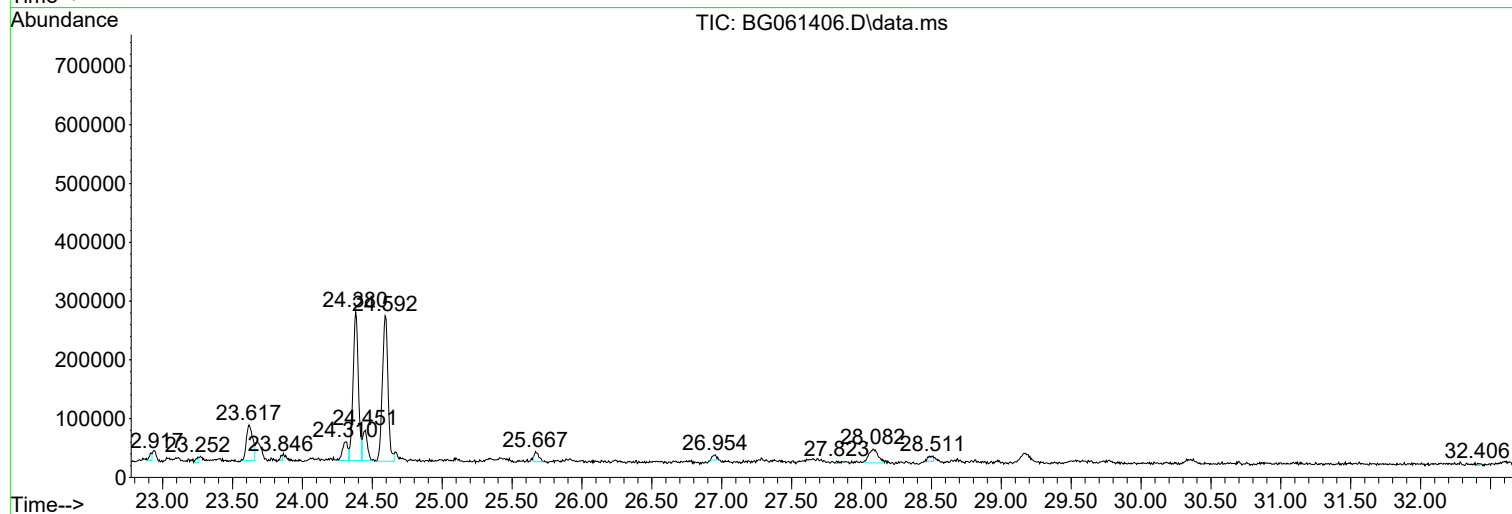
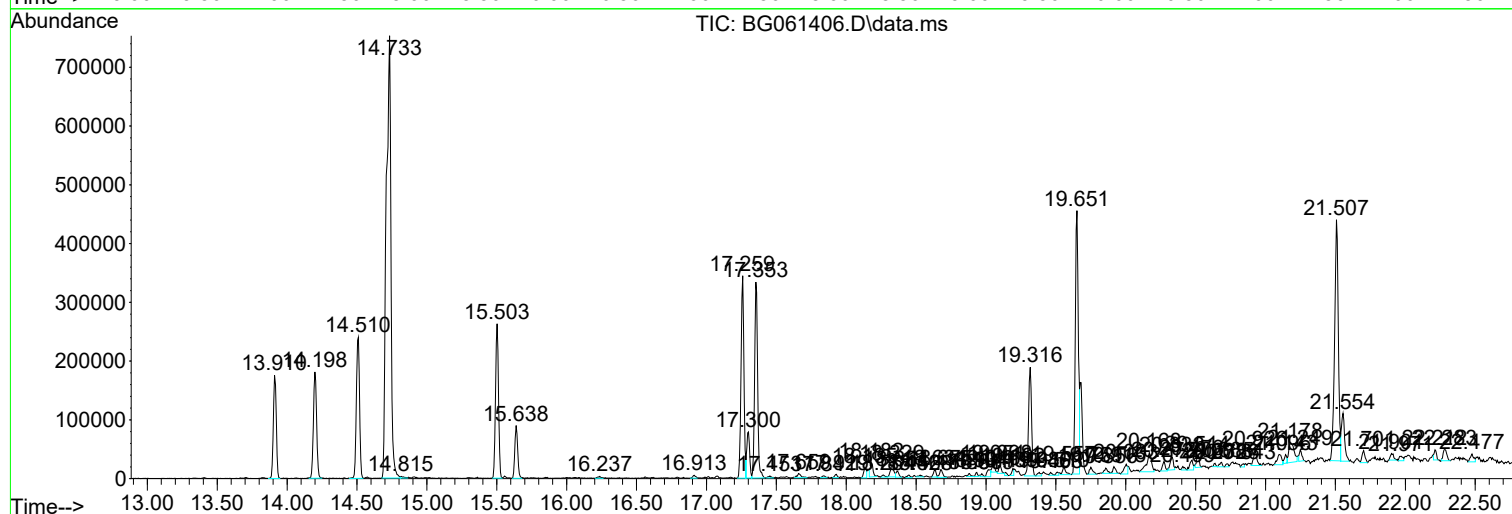
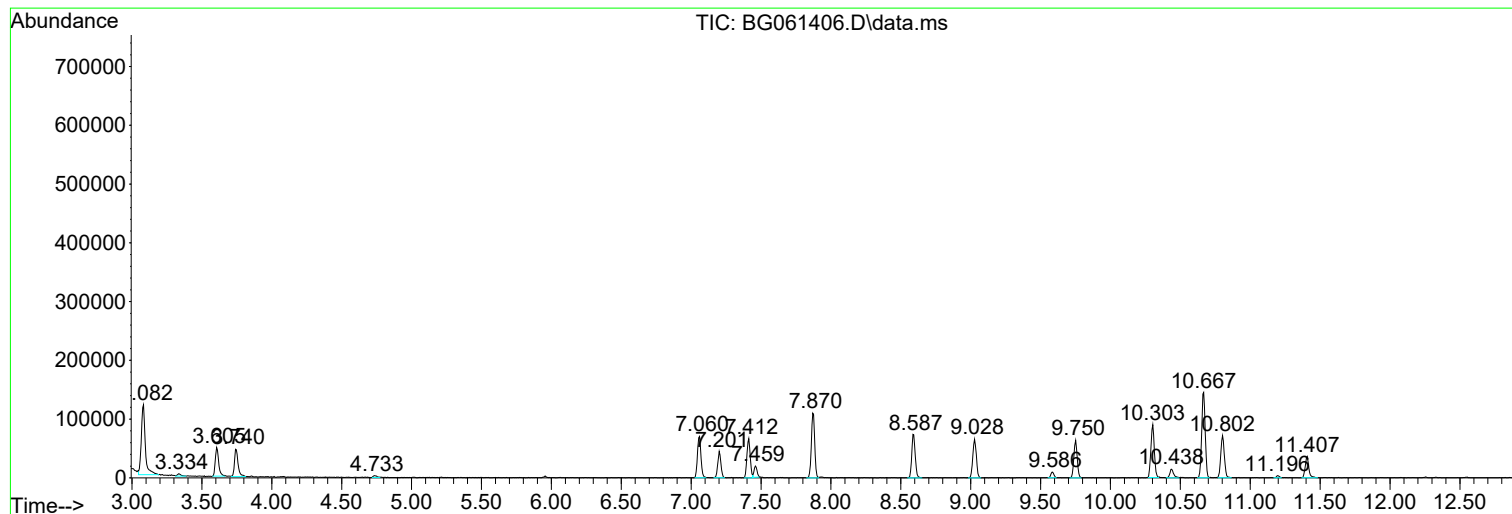
Sum of corrected areas: 11060194

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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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



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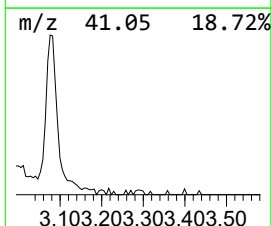
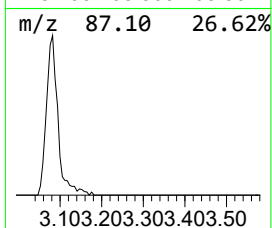
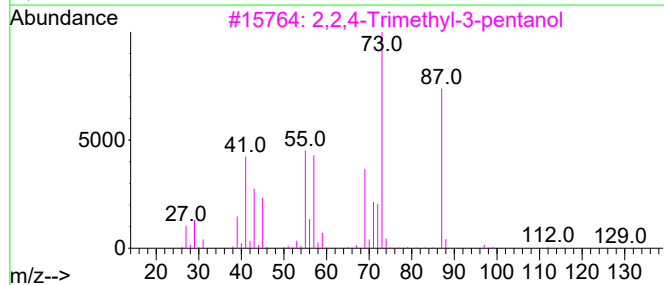
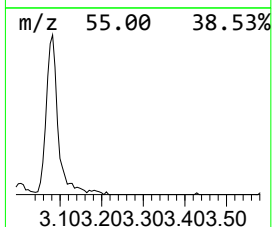
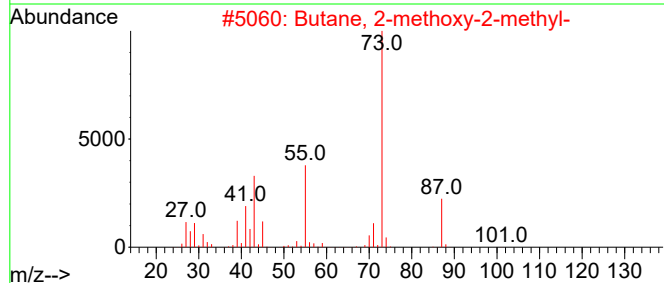
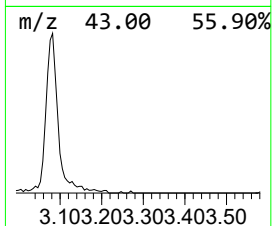
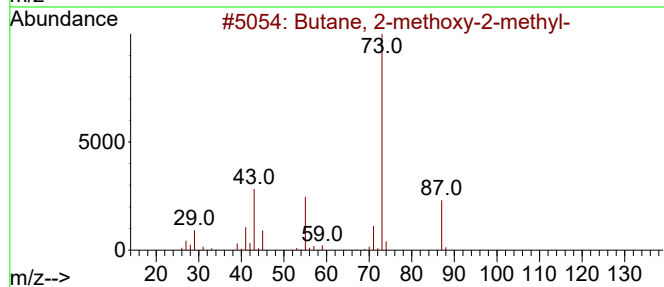
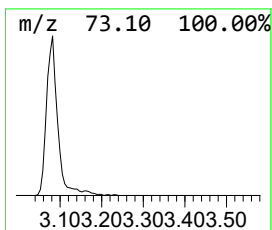
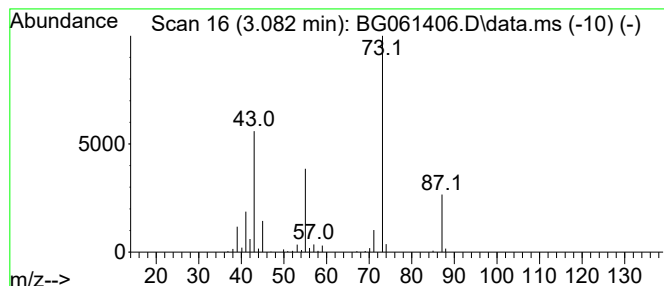
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.082	25.45 ng/ul	237088	1,4-Dichlorobenzene-d4	7.876

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	45
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
4			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	38
5			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38



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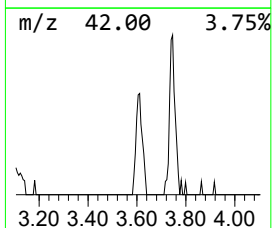
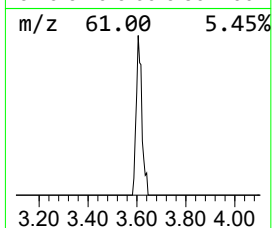
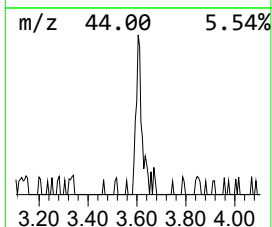
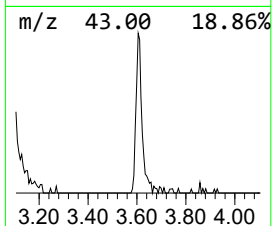
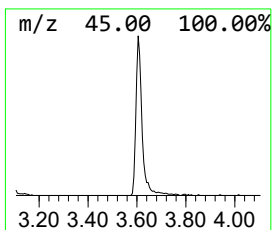
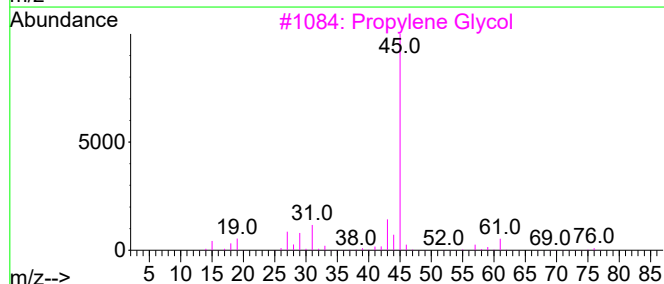
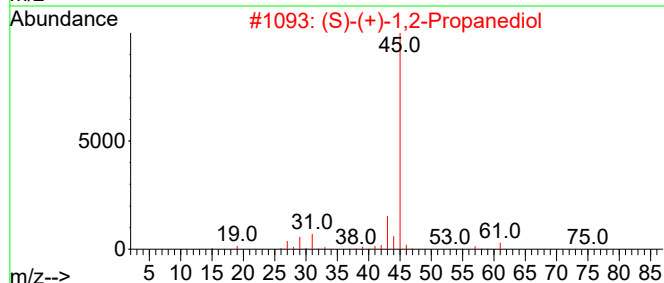
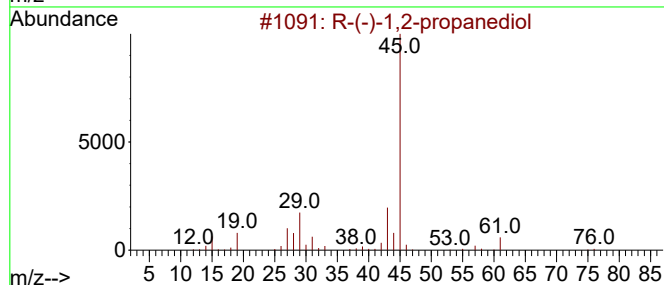
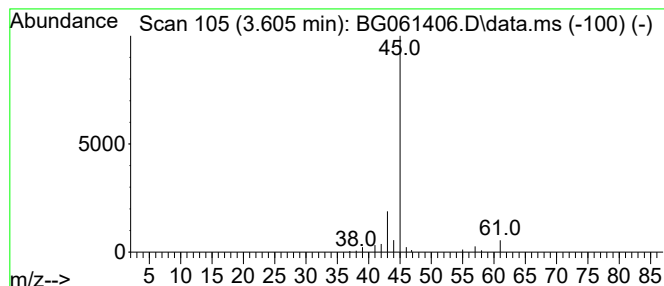
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.605	8.67 ng/ul	80753	1,4-Dichlorobenzene-d4	7.876

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	43
2	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	43
3	Propylene Glycol			76	C3H8O2	000057-55-6	9
4	Propylene Glycol			76	C3H8O2	000057-55-6	9
5	Isopropyl Alcohol			60	C3H8O	000067-63-0	9



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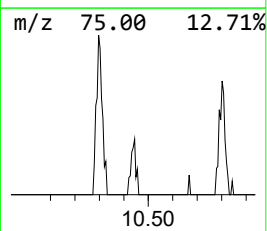
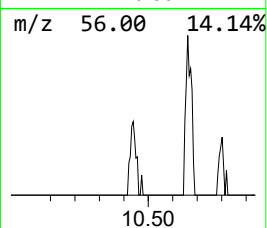
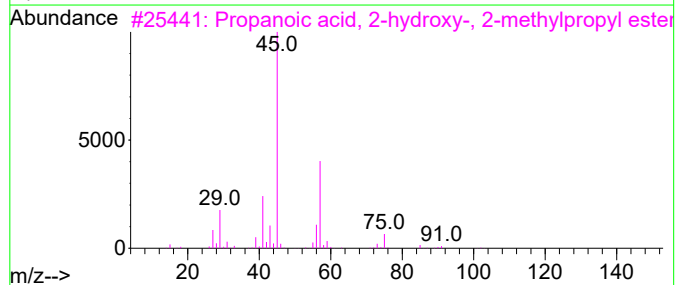
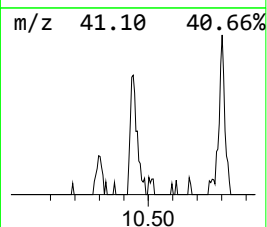
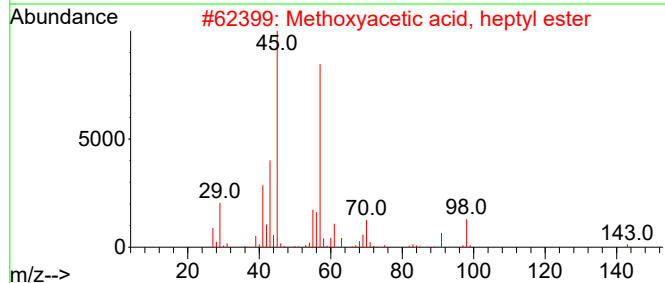
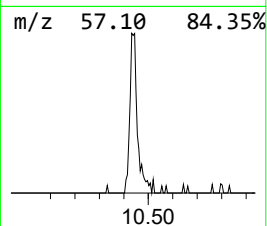
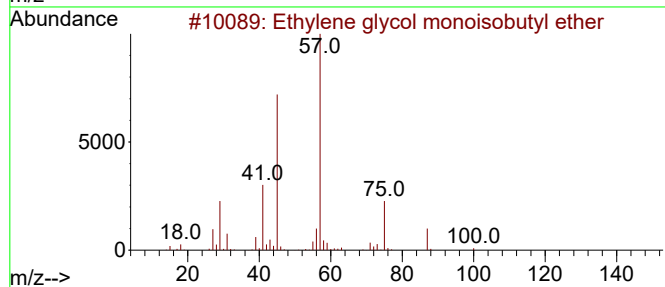
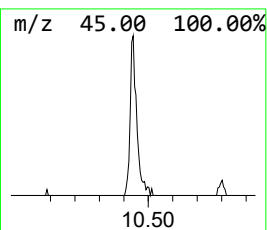
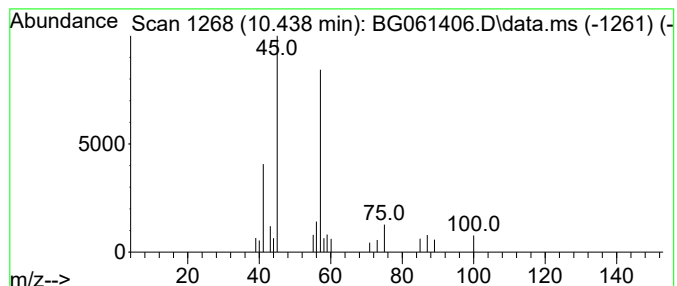
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethylene glycol monoisobuty... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.438	2.18 ng/ul	27918	Naphthalene-d8	10.667

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	50
2			Methoxyacetic acid, heptyl ester	188	C10H20O3	168920-36-3	42
3			Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	39
4			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	23
5			1-Methoxy-3-methyl-3-butene	100	C6H12O	034752-58-4	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053124\
Data File : BG061406.D
Acq On : 1 Jun 2024 00:21
Operator : MA/JU
Sample : P2588-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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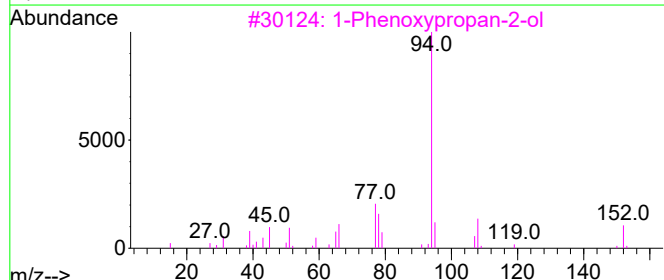
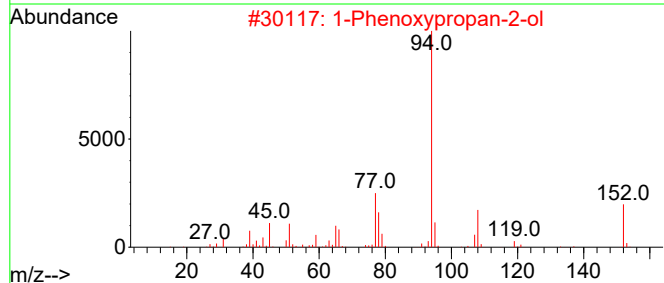
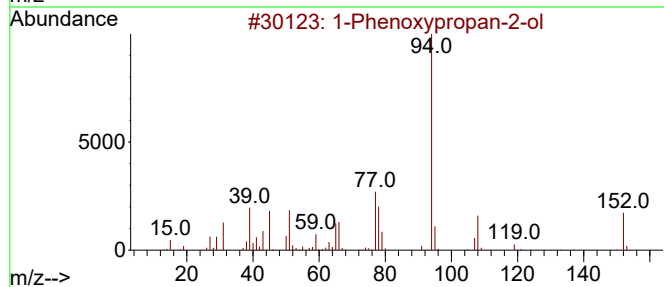
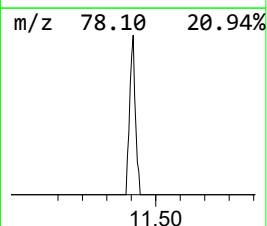
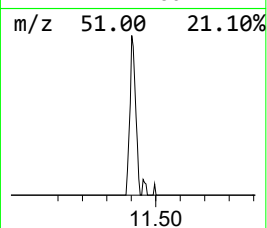
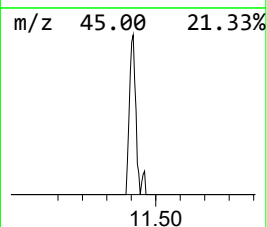
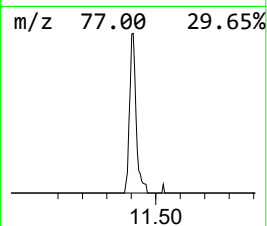
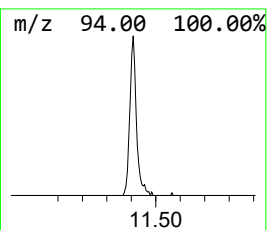
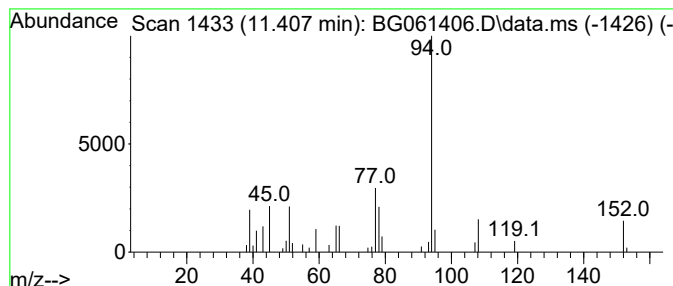
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.407	5.06 ng/ul	64707	Naphthalene-d8	10.667

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053124\
Data File : BG061406.D
Acq On : 1 Jun 2024 00:21
Operator : MA/JU
Sample : P2588-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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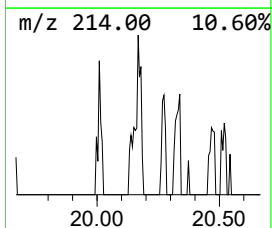
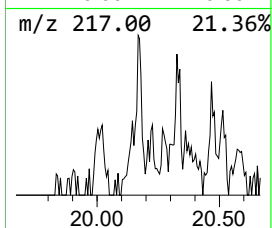
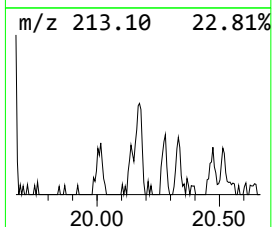
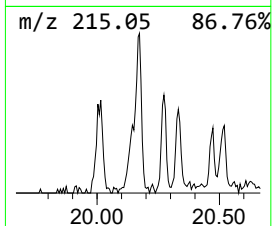
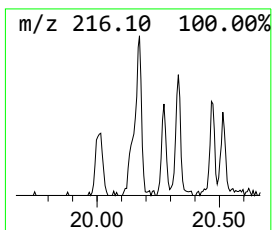
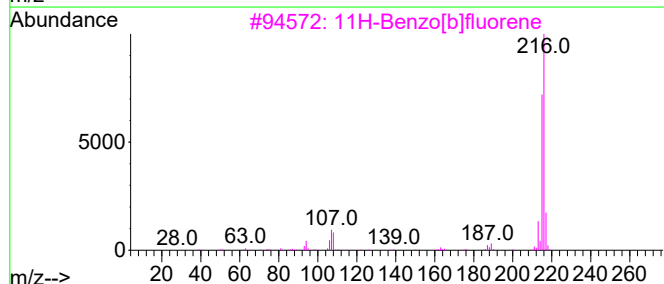
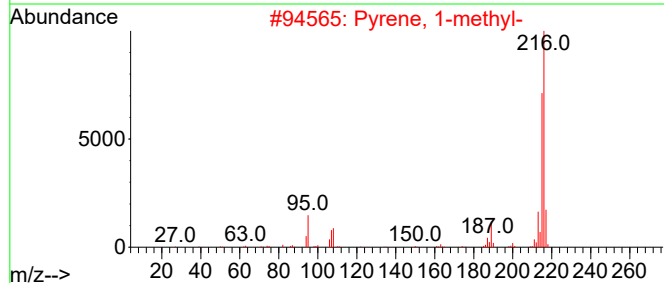
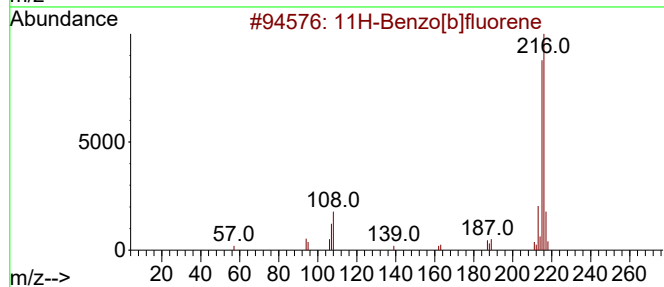
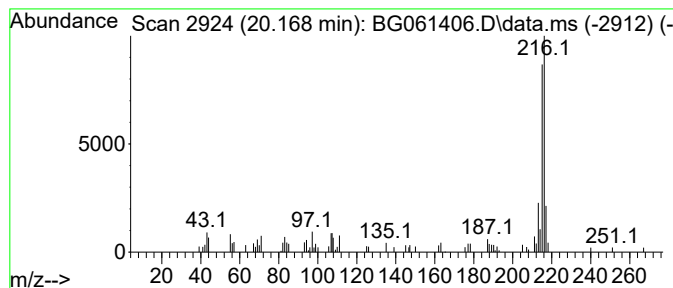
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 11H-Benzo[b]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.168	2.04 ng/ul	78241	Chrysene-d12	21.507

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053124\
Data File : BG061406.D
Acq On : 1 Jun 2024 00:21
Operator : MA/JU
Sample : P2588-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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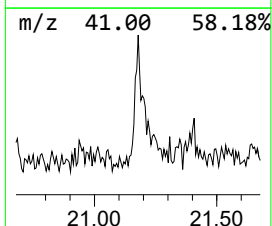
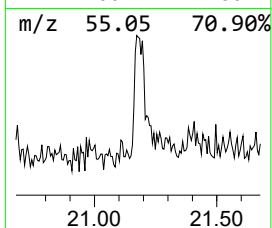
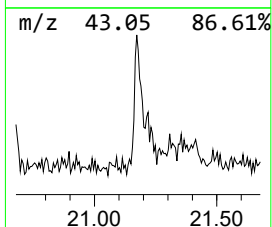
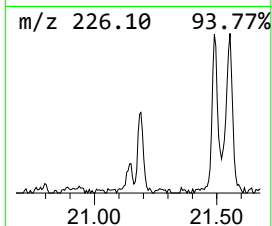
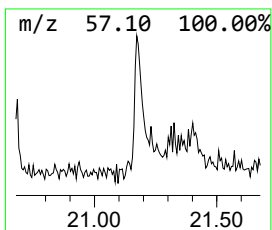
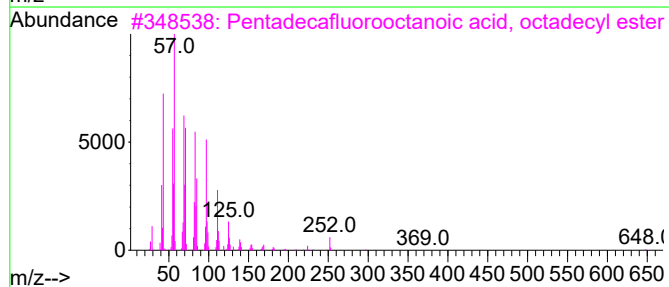
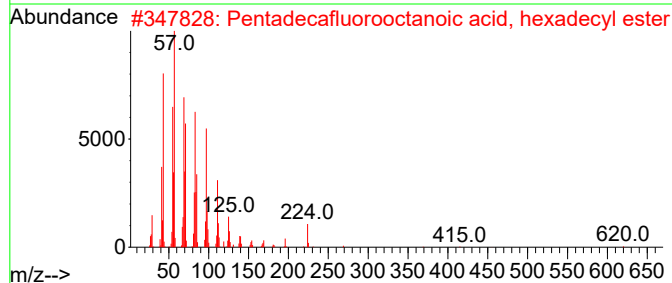
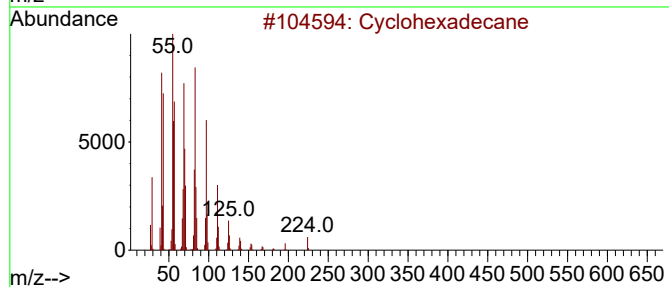
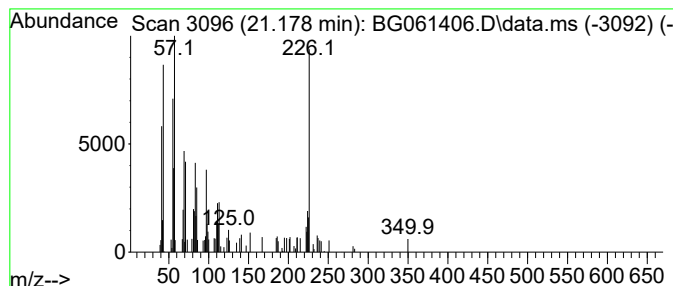
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic21.178 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.178	2.31 ng/ul	88617	Chrysene-d12	21.507

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	87
2			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	55
3			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	55
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	55
5			Carbonic acid, isobutyl pentadec...	328	C20H40O3	1000314-61-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053124\
Data File : BG061406.D
Acq On : 1 Jun 2024 00:21
Operator : MA/JU
Sample : P2588-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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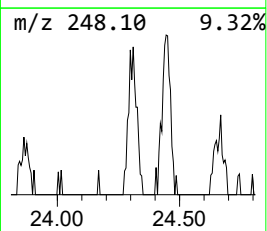
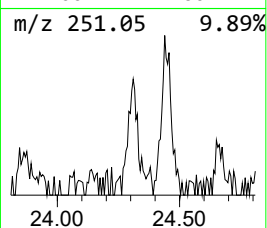
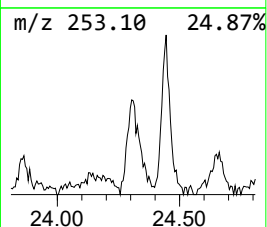
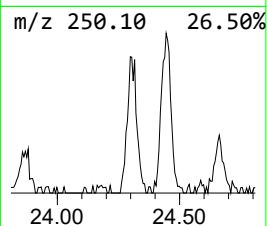
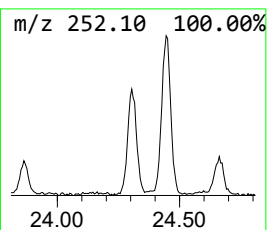
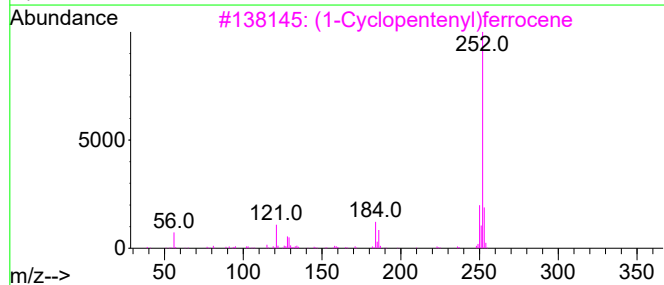
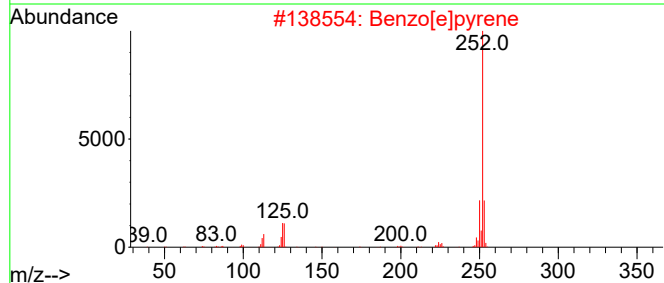
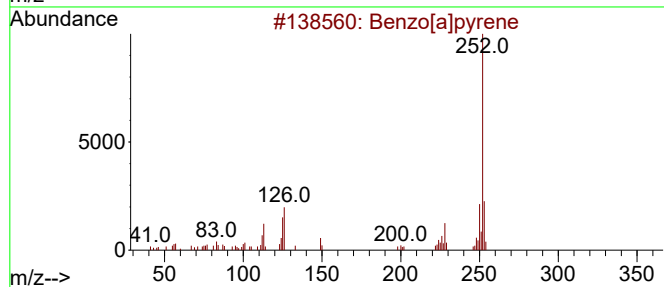
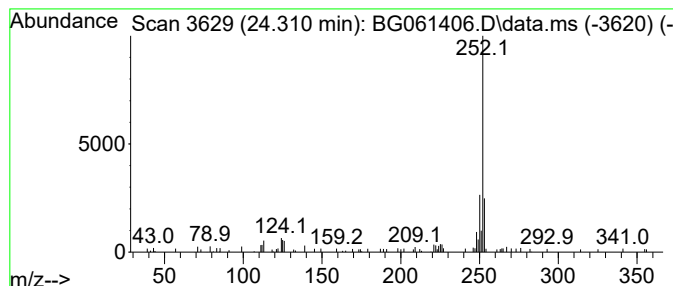
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.310	2.55 ng/ul	89017	Perylene-d12	24.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	76
2			Benzo[e]pyrene	252	C20H12	000192-97-2	72
3			(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	72
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	70
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG053124\
Data File : BG061406.D
Acq On : 1 Jun 2024 00:21
Operator : MA/JU
Sample : P2588-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH5S5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.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
Butane, 2-metho...	3.082	25.4	ng/ul	237088	1	7.876	186318	20.0
unknown-01	3.605	8.7	ng/ul	80753	1	7.876	186318	20.0
Ethylene glycol...	10.438	2.2	ng/ul	27918	2	10.667	255615	20.0
1-Phenoxypropan...	11.407	5.1	ng/ul	64707	2	10.667	255615	20.0
11H-Benzo[b]flu...	20.168	2.0	ng/ul	78241	5	21.507	765855	20.0
(DEL) Alkane: C...	21.178	2.3	ng/ul	88617	5	21.507	765855	20.0
Benzo[e]pyrene	24.310	2.5	ng/ul	89017	6	24.592	698720	20.0