

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052324\
 Data File : BG061274.D
 Acq On : 23 May 2024 16:47
 Operator : MA/JU
 Sample : P2547-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BH5P1

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: BG061274.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.081	11	16	31	rVB	163557	234835	31.51%	2.710%
2	3.340	57	60	64	rBV4	3588	4515	0.61%	0.052%
3	3.610	102	106	116	rVB	24408	39054	5.24%	0.451%
4	3.745	125	129	138	rBV	49599	80852	10.85%	0.933%
5	5.966	502	507	511	rVB	1843	2848	0.38%	0.033%
6	7.065	687	694	705	rVB2	66016	110548	14.84%	1.276%
7	7.218	713	720	726	rBV	42647	69479	9.32%	0.802%
8	7.423	749	755	760	rBV	60901	96361	12.93%	1.112%
9	7.476	760	764	771	rVB	6349	9249	1.24%	0.107%
10	7.887	827	834	840	rVB	71754	116795	15.67%	1.348%
11	8.598	948	955	967	rVB	79131	133735	17.95%	1.543%
12	9.039	1024	1030	1038	rBV	67376	112784	15.14%	1.302%
13	9.597	1120	1125	1131	rBV2	5063	7397	0.99%	0.085%
14	9.762	1147	1153	1160	rBV	60806	102963	13.82%	1.188%
15	10.314	1239	1247	1256	rBV	81373	145036	19.46%	1.674%
16	10.449	1266	1270	1283	rBV2	13600	25503	3.42%	0.294%
17	10.678	1302	1309	1316	rBV	107709	184052	24.70%	2.124%
18	10.813	1325	1332	1342	rVB2	85375	148447	19.92%	1.713%
19	11.419	1429	1435	1442	rBV2	15392	28138	3.78%	0.325%
20	12.265	1576	1579	1584	rVB2	1541	1903	0.26%	0.022%
21	13.921	1856	1861	1868	rVB	191758	272662	36.59%	3.147%
22	14.209	1904	1910	1920	rVB2	196367	299255	40.16%	3.454%
23	14.515	1956	1962	1970	rVV	198095	305971	41.06%	3.531%
24	14.580	1970	1973	1978	rVB2	4051	6121	0.82%	0.071%
25	14.721	1991	1997	1998	rBV	131621	159793	21.44%	1.844%
26	14.744	1998	2001	2012	rVB3	184135	283698	38.07%	3.274%
27	14.926	2028	2032	2038	rVB2	3179	4852	0.65%	0.056%
28	15.514	2124	2132	2138	rBV	274221	405704	54.45%	4.682%
29	15.567	2138	2141	2146	rVV2	4865	6908	0.93%	0.080%
30	15.643	2149	2154	2161	rVV2	94440	141469	18.98%	1.633%
31	15.866	2188	2192	2197	rVB2	2071	2666	0.36%	0.031%
32	16.013	2212	2217	2223	rVB2	1732	3078	0.41%	0.036%
33	16.066	2223	2226	2230	rBV2	1014	1831	0.25%	0.021%
34	16.242	2252	2256	2261	rVB3	1645	3004	0.40%	0.035%
35	16.571	2310	2312	2316	rVB2	1507	1810	0.24%	0.021%
36	16.806	2348	2352	2355	rBV2	1616	2667	0.36%	0.031%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
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37	16.924	2367	2372	2377	rBV	5177	6140	0.82%	0.071%
38	17.018	2386	2388	2391	rVB5	1478	1775	0.24%	0.020%
39	17.082	2396	2399	2404	rVB2	4323	5642	0.76%	0.065%
40	17.182	2413	2416	2420	rVB3	3856	4470	0.60%	0.052%
41	17.265	2424	2430	2434	rVV	303644	443140	59.47%	5.114%
42	17.306	2434	2437	2442	rVV	101198	147960	19.86%	1.708%
43	17.364	2442	2447	2459	rVV2	344335	517343	69.43%	5.970%
44	17.458	2459	2463	2467	rVV3	2042	3117	0.42%	0.036%
45	17.541	2473	2477	2481	rBV3	1261	2063	0.28%	0.024%
46	17.670	2493	2499	2505	rBV2	10695	16611	2.23%	0.192%
47	17.770	2511	2516	2517	rBV3	1782	2177	0.29%	0.025%
48	17.846	2527	2529	2535	rVB3	2106	2759	0.37%	0.032%
49	17.934	2540	2544	2550	rBV2	9957	13039	1.75%	0.150%
50	18.140	2571	2579	2583	rBV	12582	19795	2.66%	0.228%
51	18.193	2583	2588	2592	rVV	18371	28017	3.76%	0.323%
52	18.228	2592	2594	2599	rVV2	4721	6856	0.92%	0.079%
53	18.275	2599	2602	2605	rVV	5468	7028	0.94%	0.081%
54	18.340	2606	2613	2616	rVV3	18146	34038	4.57%	0.393%
55	18.369	2616	2618	2624	rVB3	8646	12489	1.68%	0.144%
56	18.457	2630	2633	2636	rVB3	2622	2794	0.37%	0.032%
57	18.498	2637	2640	2643	rVB2	1855	1847	0.25%	0.021%
58	18.645	2660	2665	2668	rVV2	9785	12231	1.64%	0.141%
59	18.686	2668	2672	2676	rVB	11042	14107	1.89%	0.163%
60	18.886	2698	2706	2710	rVB4	3526	6712	0.90%	0.077%
61	18.921	2710	2712	2713	rBV	2788	2221	0.30%	0.026%
62	18.939	2713	2715	2719	rVV2	3431	3895	0.52%	0.045%
63	18.974	2719	2721	2725	rVB2	2995	3475	0.47%	0.040%
64	19.062	2730	2736	2739	rBV3	9203	17614	2.36%	0.203%
65	19.109	2740	2744	2750	rVB4	22512	28600	3.84%	0.330%
66	19.204	2755	2760	2770	rVB6	10183	24522	3.29%	0.283%
67	19.321	2772	2780	2787	rVB	198265	273913	36.76%	3.161%
68	19.386	2788	2791	2793	rBV	2233	2278	0.31%	0.026%
69	19.421	2794	2797	2802	rVB4	4755	4548	0.61%	0.052%
70	19.468	2804	2805	2808	rBV2	1967	1802	0.24%	0.021%
71	19.562	2818	2821	2824	rBV2	5233	6919	0.93%	0.080%
72	19.656	2831	2837	2841	rBV	489206	745163	100.00%	8.600%
73	19.756	2850	2854	2859	rBV3	7925	9958	1.34%	0.115%
74	19.862	2868	2872	2879	rVB	9328	15476	2.08%	0.179%
75	19.920	2879	2882	2887	rVB3	8646	13952	1.87%	0.161%
76	20.008	2891	2897	2898	rBV3	12268	20674	2.77%	0.239%

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 Title : SVOA CALIBRATION

77	20.055	2904	2905	2909	rVB3	3950	4032	0.54%	0.047%
78	20.097	2909	2912	2915	rBV5	2031	2893	0.39%	0.033%
79	20.138	2915	2919	2921	rVV4	7921	11991	1.61%	0.138%
80	20.179	2923	2926	2933	rVB2	19675	31770	4.26%	0.367%
81	20.279	2939	2943	2946	rBV4	11606	13850	1.86%	0.160%
82	20.332	2948	2952	2957	rVB2	14351	23958	3.22%	0.276%
83	20.478	2973	2977	2981	rBV2	7937	13210	1.77%	0.152%
84	20.520	2981	2984	2989	rVV4	8565	12452	1.67%	0.144%
85	20.696	3010	3014	3017	rVB6	5662	7452	1.00%	0.086%
86	20.843	3037	3039	3044	rBV6	4055	4862	0.65%	0.056%
87	20.931	3050	3054	3060	rVB2	19133	29910	4.01%	0.345%
88	21.107	3082	3084	3088	rVB	14092	13751	1.85%	0.159%
89	21.148	3088	3091	3094	rVV2	10441	15467	2.08%	0.178%
90	21.189	3094	3098	3105	rVV4	27756	72189	9.69%	0.833%
91	21.254	3107	3109	3116	rVB	23355	32890	4.41%	0.380%
92	21.518	3146	3154	3158	rBV2	351897	687690	92.29%	7.936%
93	21.560	3158	3161	3171	rVB	85053	134274	18.02%	1.550%
94	21.712	3182	3187	3193	rBV4	19822	28908	3.88%	0.334%
95	22.929	3390	3394	3405	rVB5	31396	80365	10.78%	0.927%
96	23.134	3426	3429	3435	rVB2	24900	43249	5.80%	0.499%
97	23.622	3505	3512	3520	rBV2	56564	180874	24.27%	2.087%
98	24.321	3624	3631	3634	rBV	27172	60701	8.15%	0.701%
99	24.392	3635	3643	3650	rVV2	224855	547304	73.45%	6.316%
100	24.603	3669	3679	3689	rBV2	195754	567880	76.21%	6.554%

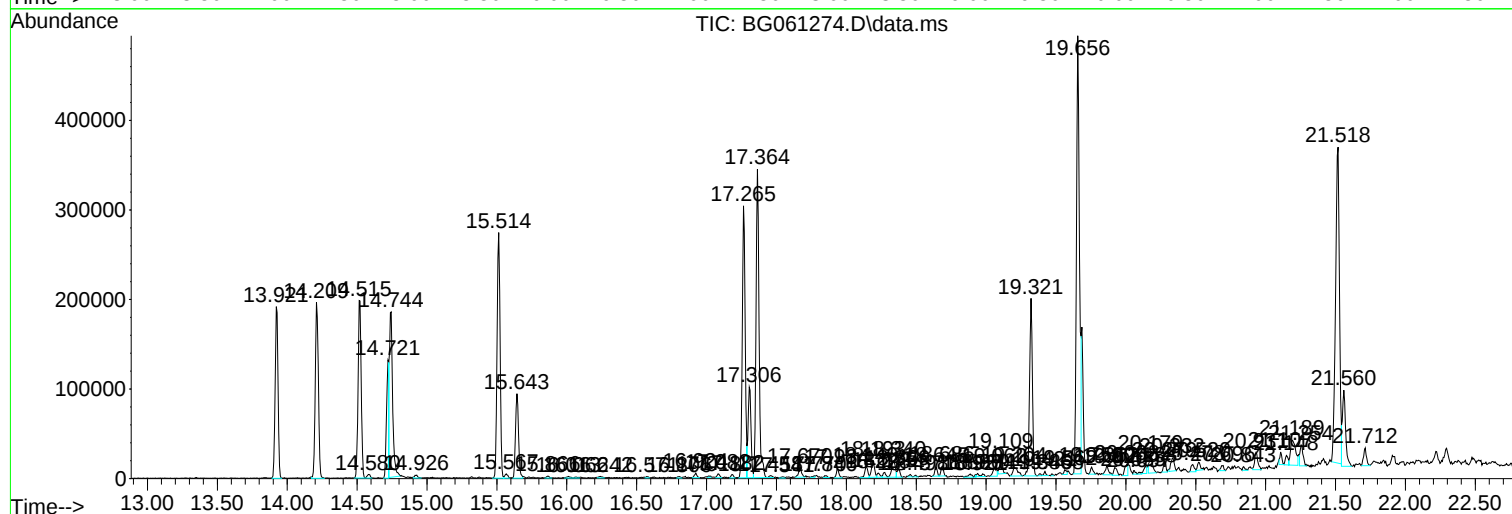
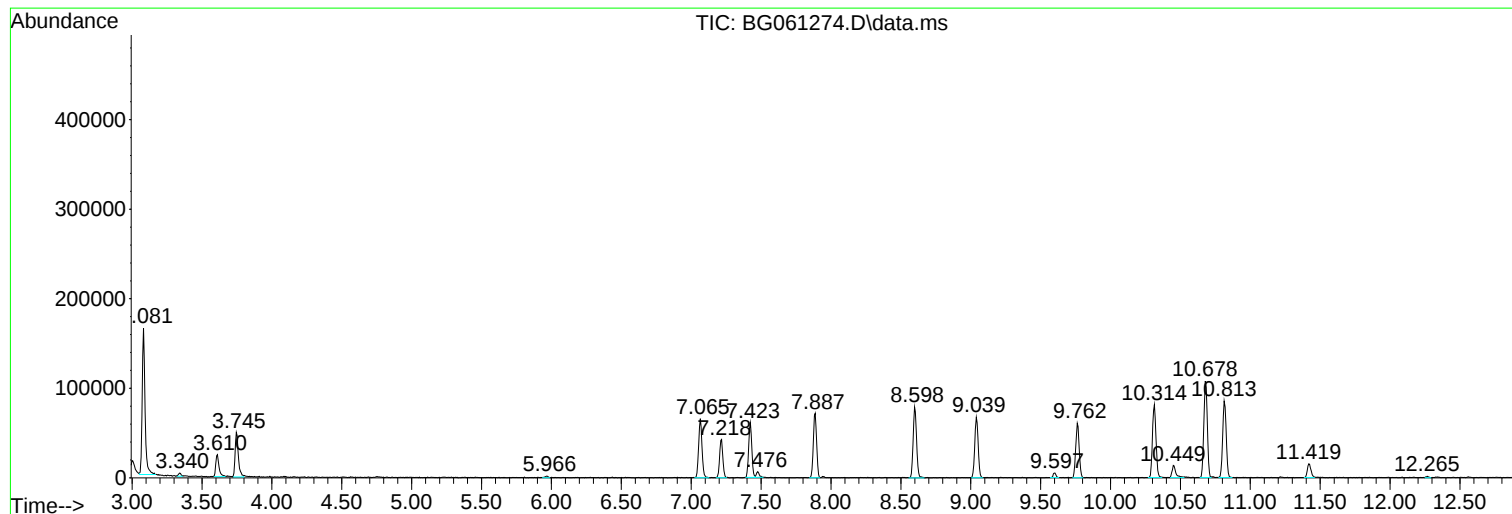
Sum of corrected areas: 8665065

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ALS Vial : 10 Sample Multiplier: 1

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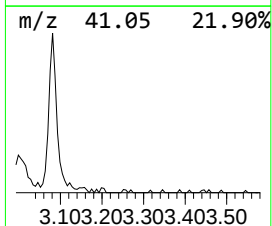
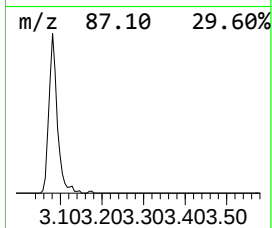
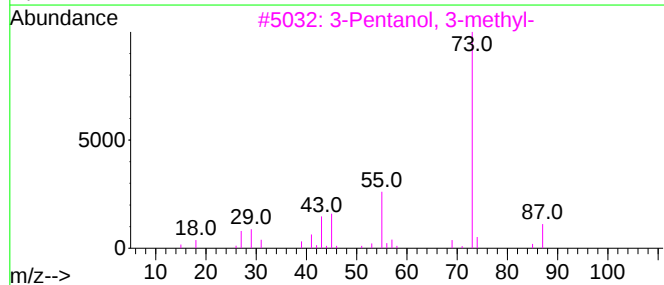
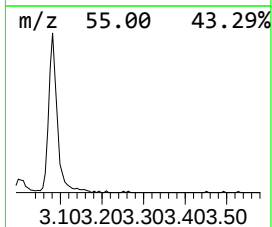
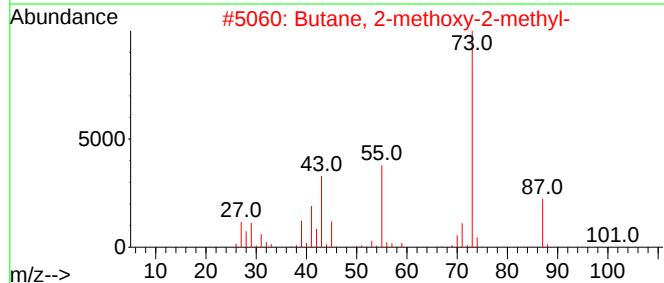
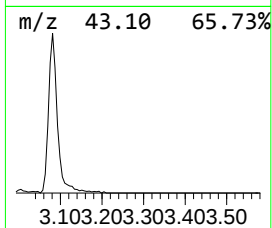
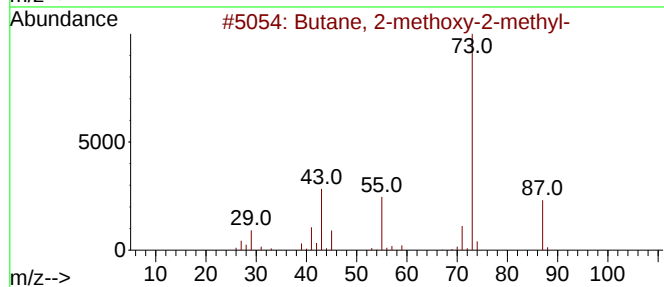
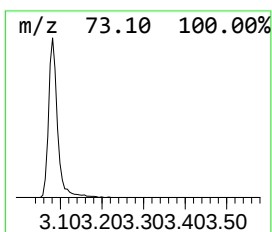
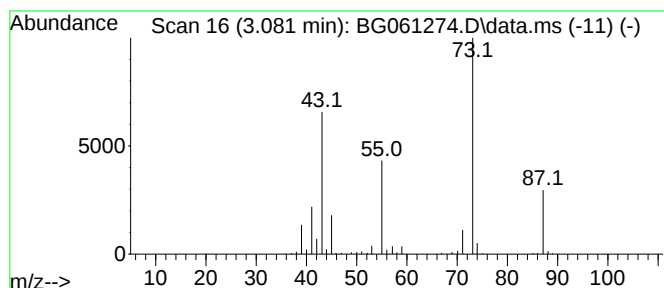
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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.
3.081	40.21 ng/ul	234835	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
3			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
4			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
5			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28



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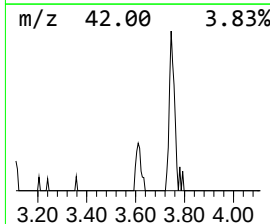
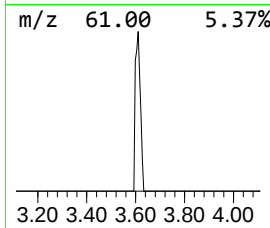
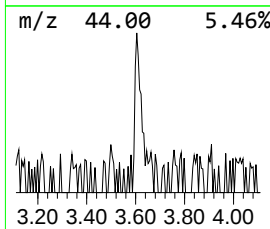
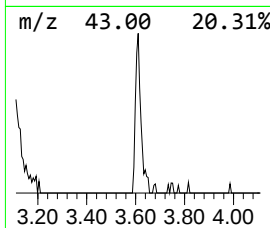
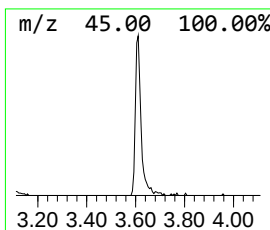
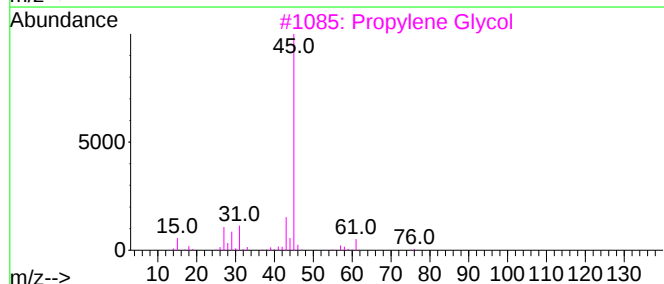
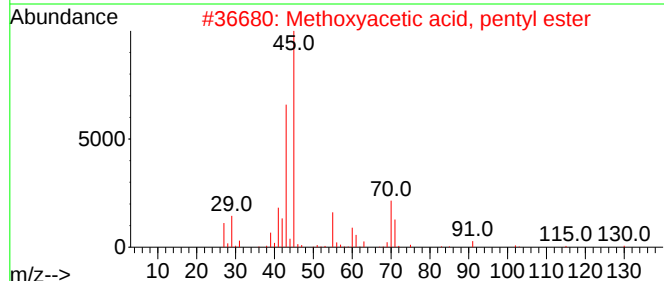
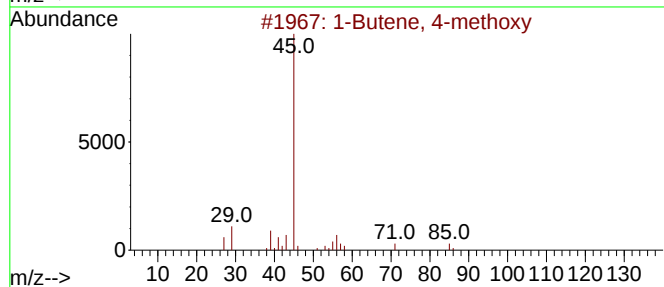
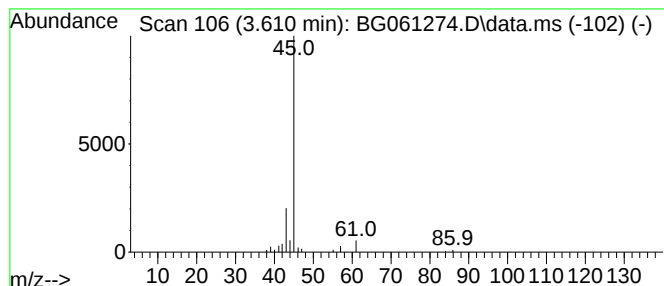
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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.
3.610	6.69 ng/ul	39054	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butene, 4-methoxy	86	C5H10O	004696-30-4	9
2			Methoxyacetic acid, pentyl ester	160	C8H16O3	168920-35-2	9
3			Propylene Glycol	76	C3H8O2	000057-55-6	9
4			Acetic acid, cyano-, 2-methoxyet...	143	C6H9NO3	010258-54-5	9
5			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



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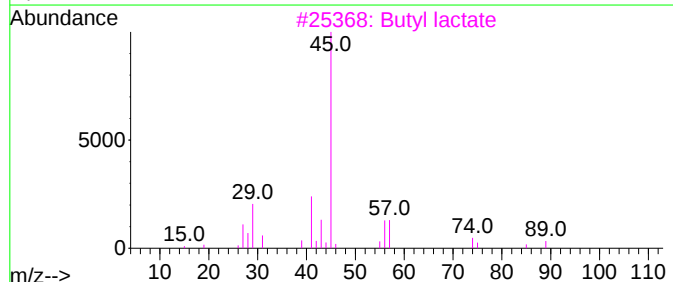
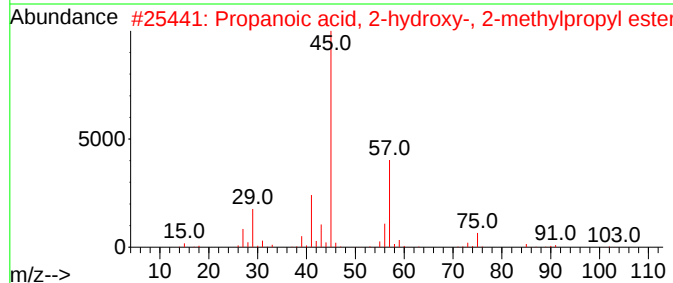
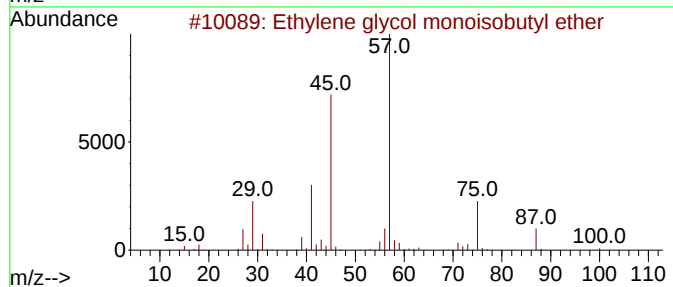
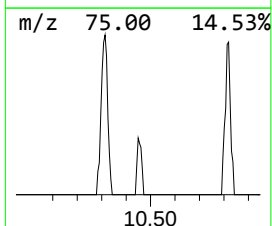
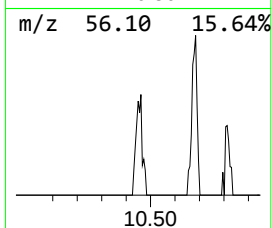
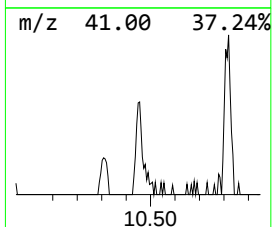
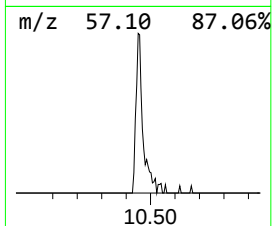
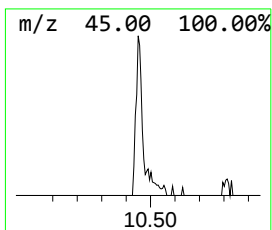
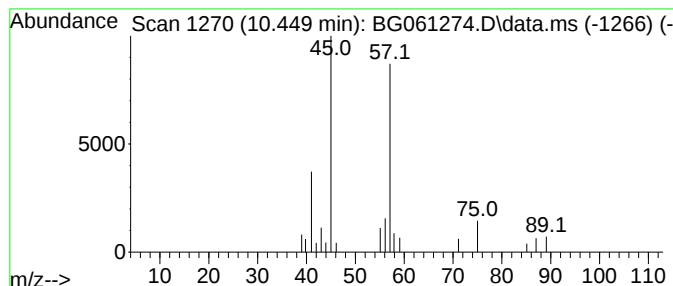
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.449	2.77 ng/ul	25503	Naphthalene-d8	10.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	40
2			Propanoic acid, 2-hydroxy-, 2-me...	146	C7H14O3	000585-24-0	39
3			Butyl lactate	146	C7H14O3	000138-22-7	38
4			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	36
5			2-Pentanol, 3-methyl-	102	C6H14O	000565-60-6	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052324\
Data File : BG061274.D
Acq On : 23 May 2024 16:47
Operator : MA/JU
Sample : P2547-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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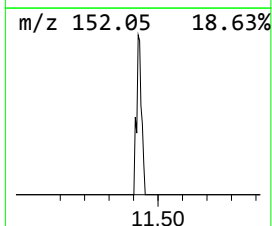
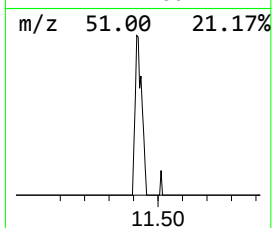
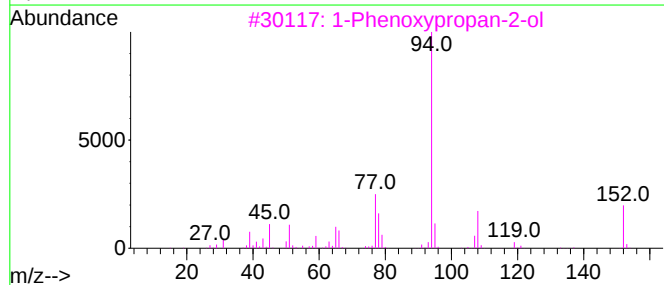
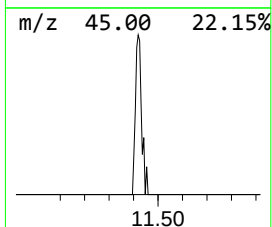
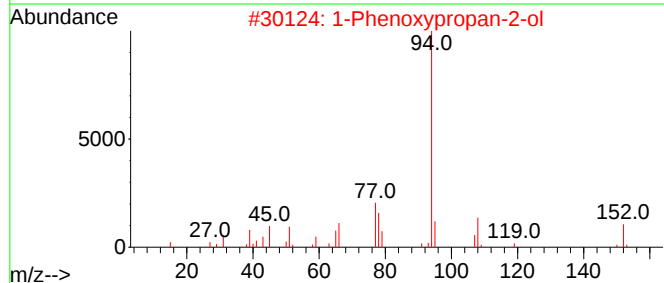
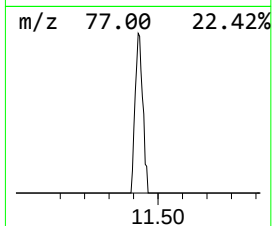
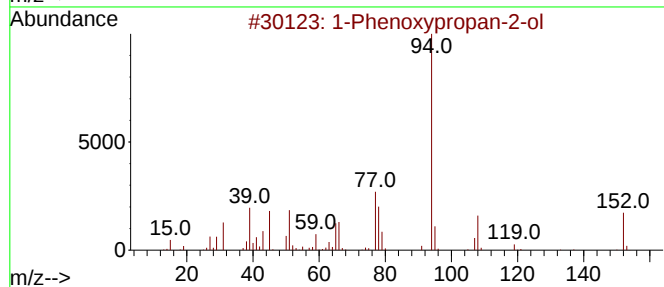
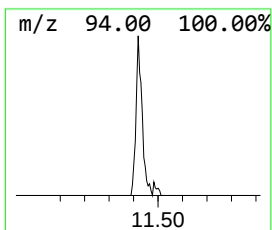
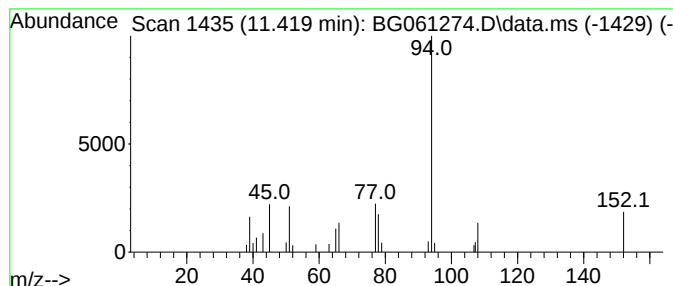
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Quant Title : SVOA CALIBRATION

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.
11.419	3.06 ng/ul	28138	Naphthalene-d8	10.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	80
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052324\
Data File : BG061274.D
Acq On : 23 May 2024 16:47
Operator : MA/JU
Sample : P2547-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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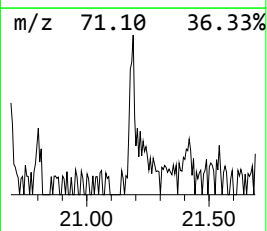
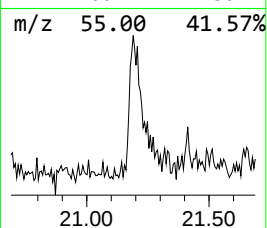
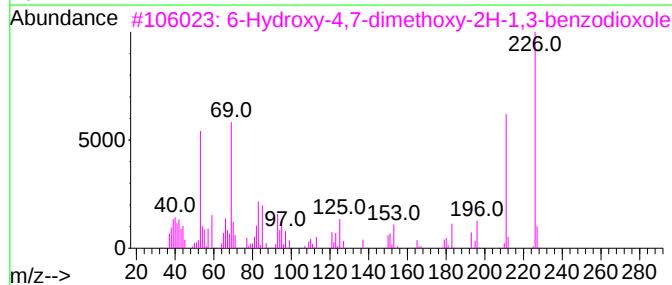
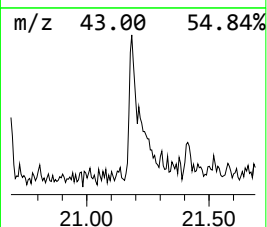
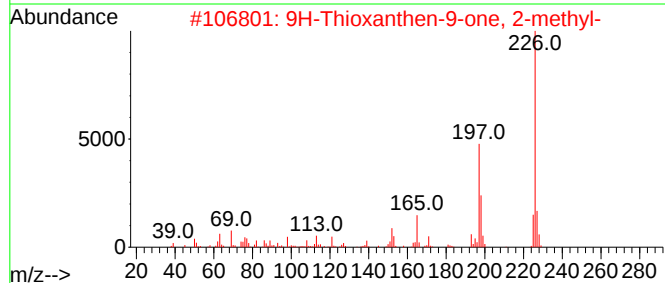
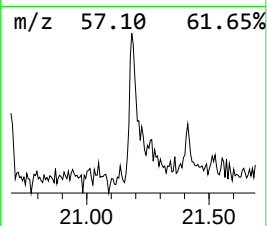
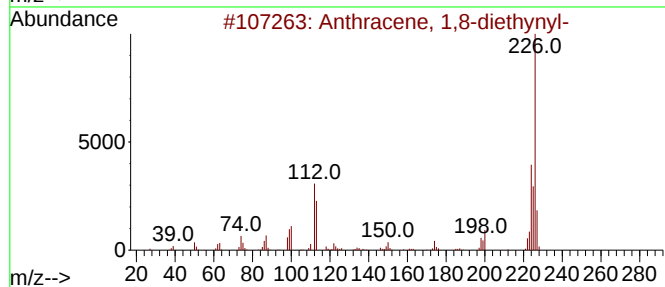
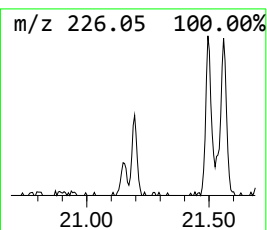
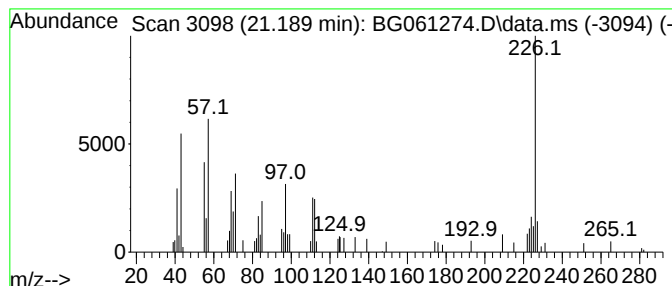
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.189	2.10 ng/ul	72189	Chrysene-d12	21.518

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1,8-diethynyl-	226	C18H10	078053-58-4	46
2			9H-Thioxanthen-9-one, 2-methyl-	226	C14H10O5	015774-82-0	43
3			6-Hydroxy-4,7-dimethoxy-2H-1,3-b...	226	C10H10O6	849798-24-9	43
4			9H-Xanthen-9-one, 3-methoxy-	226	C14H10O3	003722-52-9	38
5			2-Fluoroanthra-9,10-quinone	226	C14H7FO2	000572-84-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052324\
Data File : BG061274.D
Acq On : 23 May 2024 16:47
Operator : MA/JU
Sample : P2547-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH5P1

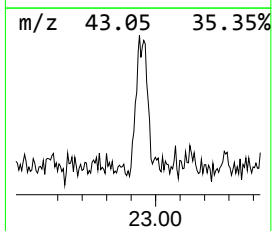
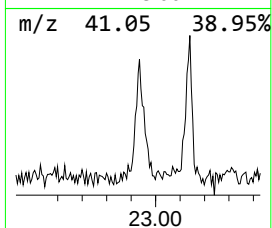
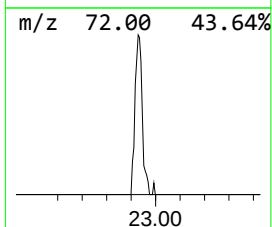
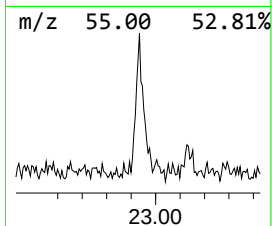
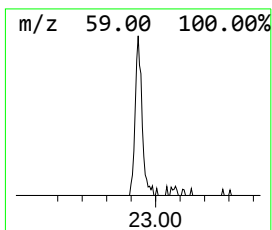
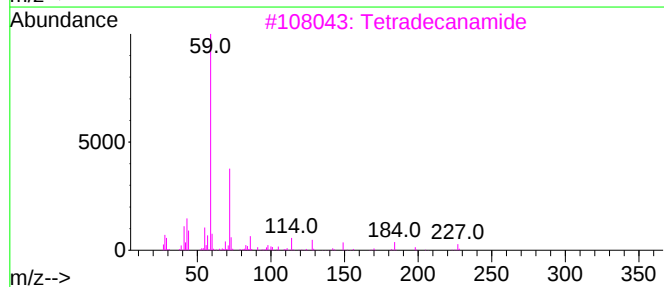
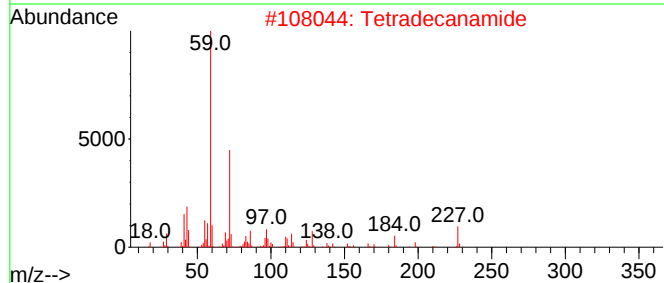
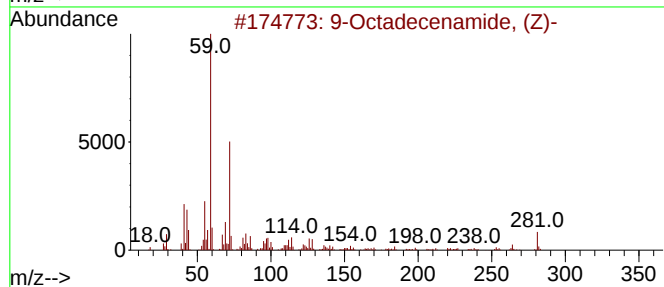
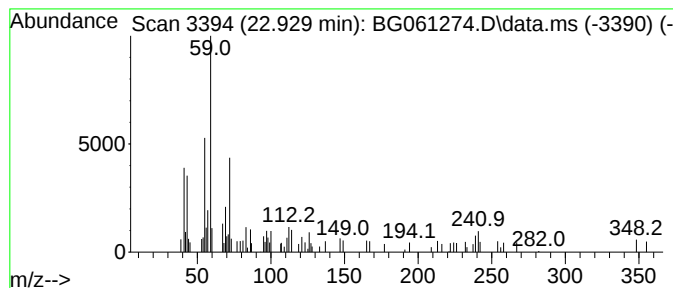
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.929	2.34 ng/ul	80365	Chrysene-d12	21.518

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	59
2	Tetradecanamide			227	C14H29NO	000638-58-4	59
3	Tetradecanamide			227	C14H29NO	000638-58-4	59
4	Hexadecanamide			255	C16H33NO	000629-54-9	59
5	Butanal, oxime			87	C4H9NO	000110-69-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052324\
Data File : BG061274.D
Acq On : 23 May 2024 16:47
Operator : MA/JU
Sample : P2547-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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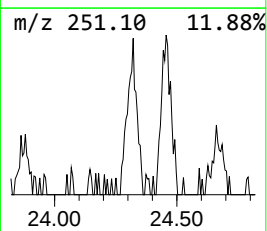
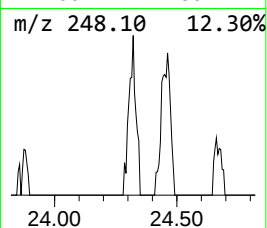
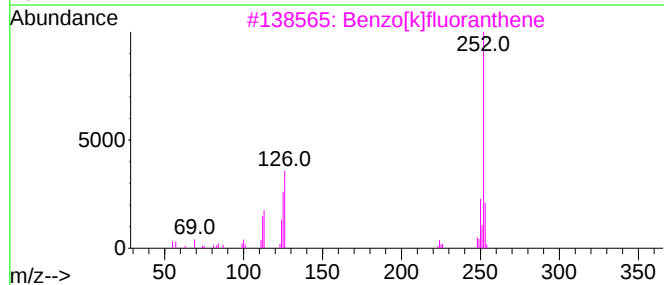
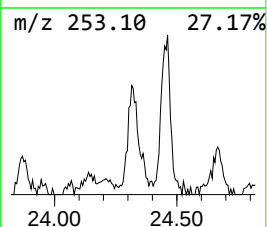
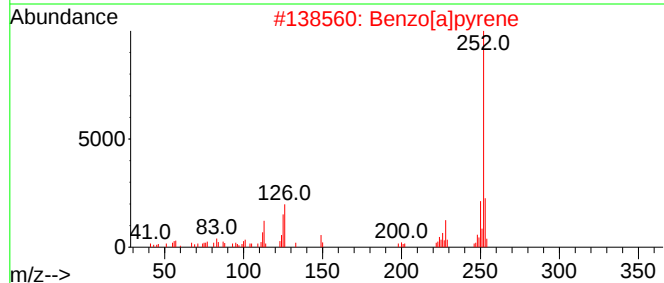
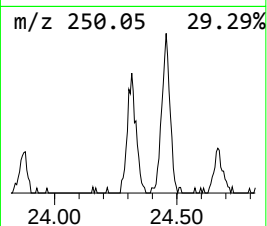
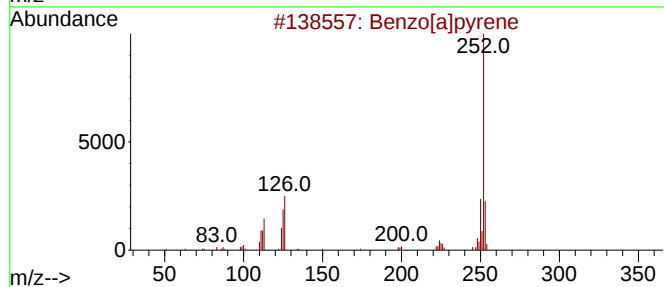
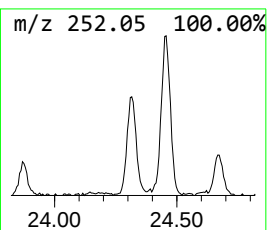
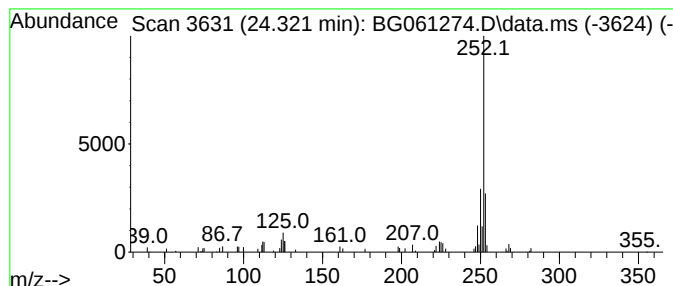
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Perylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.321	2.14 ng/ul	60701	Perylene-d12	24.603

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	90
2			Benzo[a]pyrene	252	C20H12	000050-32-8	76
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	76
4			Perylene	252	C20H12	000198-55-0	76
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052324\
Data File : BG061274.D
Acq On : 23 May 2024 16:47
Operator : MA/JU
Sample : P2547-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BH5P1

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
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unknown-01	3.610	6.7	ng/ul	39054	1	7.881	116795	20.0
unknown-02	10.449	2.8	ng/ul	25503	2	10.678	184052	20.0
1-Phenoxypropan...	11.419	3.1	ng/ul	28138	2	10.678	184052	20.0
unknown-03	21.189	2.1	ng/ul	72189	5	21.518	687690	20.0
9-Octadecenamid...	22.929	2.3	ng/ul	80365	5	21.518	687690	20.0
Perylene	24.321	2.1	ng/ul	60701	6	24.603	567880	20.0