

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062124\
 Data File : BG061629.D
 Acq On : 21 Jun 2024 20:58
 Operator : MA/JU
 Sample : P2754-11
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4BQ0

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-BG062024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061629.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.472	27	31	34	rVB4	15282	19935	1.03%	0.088%
2	3.613	53	55	60	rBV4	5300	7621	0.39%	0.034%
3	3.736	72	76	83	rBV2	25893	42510	2.19%	0.189%
4	3.883	97	101	117	rVB	222015	380296	19.61%	1.687%
5	4.236	157	161	166	rVB2	10643	13170	0.68%	0.058%
6	4.682	235	237	240	rBV3	4091	4585	0.24%	0.020%
7	4.800	253	257	259	rBV4	3639	3273	0.17%	0.015%
8	4.841	262	264	267	rBV3	2361	3825	0.20%	0.017%
9	5.146	313	316	322	rBV3	8183	14082	0.73%	0.062%
10	5.540	382	383	387	rVB4	3393	3818	0.20%	0.017%
11	6.063	470	472	475	rBV3	4579	4957	0.26%	0.022%
12	6.139	481	485	489	rBV3	9872	15604	0.80%	0.069%
13	6.322	513	516	520	rBV4	3757	4729	0.24%	0.021%
14	7.003	629	632	633	rBV	4477	3927	0.20%	0.017%
15	7.215	660	668	675	rBV2	293878	496027	25.58%	2.200%
16	7.403	692	700	707	rBV	194030	321982	16.60%	1.428%
17	7.538	721	723	726	rBV3	3976	5415	0.28%	0.024%
18	7.602	726	734	741	rVV	257020	454722	23.45%	2.017%
19	7.661	741	744	751	rVB3	9743	17318	0.89%	0.077%
20	7.790	761	766	769	rBV6	1851	3204	0.17%	0.014%
21	7.873	776	780	781	rBV2	2847	3551	0.18%	0.016%
22	8.078	809	815	821	rBV	203857	359081	18.51%	1.593%
23	8.137	821	825	828	rVB5	5705	7974	0.41%	0.035%
24	8.313	853	855	857	rBV3	4367	3889	0.20%	0.017%
25	8.396	867	869	872	rVB3	3496	3355	0.17%	0.015%
26	8.548	893	895	899	rVB4	2810	3695	0.19%	0.016%
27	8.689	918	919	922	rBV	4196	3367	0.17%	0.015%
28	8.766	922	932	939	rVV	361822	607885	31.34%	2.697%
29	9.236	1005	1012	1020	rVV	261966	467024	24.08%	2.072%
30	9.400	1038	1040	1042	rVB3	5664	4850	0.25%	0.022%
31	9.641	1079	1081	1084	rBV3	5387	4336	0.22%	0.019%
32	9.800	1102	1108	1116	rBV2	30066	53838	2.78%	0.239%
33	9.958	1127	1135	1143	rBV2	220691	392837	20.25%	1.743%
34	10.393	1206	1209	1213	rVB3	3092	5059	0.26%	0.022%
35	10.493	1216	1226	1233	rBV	348685	627007	32.33%	2.781%
36	10.611	1243	1246	1250	rBV5	4074	6157	0.32%	0.027%

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Integrator: RTE

Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
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37	10.758	1268	1271	1274	rBV2	3095	4128	0.21%	0.018%
38	10.887	1286	1293	1300	rVV	327277	582597	30.04%	2.584%
39	10.940	1300	1302	1305	rVV3	9335	11870	0.61%	0.053%
40	11.016	1307	1315	1322	rVV	355916	633157	32.65%	2.809%

41	11.181	1341	1343	1347	rBV4	3766	4311	0.22%	0.019%
42	11.410	1379	1382	1389	rVB7	12930	21205	1.09%	0.094%
43	11.621	1412	1418	1426	rBV2	49121	93224	4.81%	0.414%
44	11.709	1430	1433	1434	rBV2	4999	3985	0.21%	0.018%
45	11.874	1457	1461	1463	rBV3	5023	6160	0.32%	0.027%

46	11.903	1464	1466	1469	rVB3	7395	9050	0.47%	0.040%
47	12.291	1527	1532	1534	rBV3	4571	6711	0.35%	0.030%
48	12.320	1534	1537	1539	rVV4	3470	4043	0.21%	0.018%
49	12.350	1540	1542	1546	rVB4	6482	7194	0.37%	0.032%
50	12.456	1557	1560	1562	rBV3	7263	8274	0.43%	0.037%

51	12.538	1571	1574	1578	rVB3	7804	11821	0.61%	0.052%
52	12.749	1608	1610	1617	rVB8	7855	13560	0.70%	0.060%
53	12.849	1625	1627	1629	rVB2	4295	3266	0.17%	0.014%
54	13.243	1691	1694	1700	rVB5	12663	17624	0.91%	0.078%
55	13.490	1734	1736	1739	rBV4	5426	6036	0.31%	0.027%

56	13.519	1739	1741	1742	rBV2	4784	3572	0.18%	0.016%
57	13.619	1756	1758	1761	rVB4	7018	7622	0.39%	0.034%
58	13.684	1768	1769	1773	rVB4	5035	6097	0.31%	0.027%
59	13.760	1780	1782	1785	rBV4	5413	5435	0.28%	0.024%
60	14.018	1824	1826	1827	rBV2	9454	6804	0.35%	0.030%

61	14.036	1827	1829	1833	rVB5	6250	4852	0.25%	0.022%
62	14.107	1833	1841	1848	rBV	668014	951591	49.06%	4.221%
63	14.177	1851	1853	1858	rVB5	11605	16217	0.84%	0.072%
64	14.336	1876	1880	1883	rVV3	6590	10731	0.55%	0.048%
65	14.394	1884	1890	1897	rVB	744017	1159449	59.78%	5.143%

66	14.535	1912	1914	1917	rBV4	6277	7941	0.41%	0.035%
67	14.700	1936	1942	1948	rVB2	553821	835143	43.06%	3.705%
68	14.765	1949	1953	1957	rVB2	37340	51912	2.68%	0.230%
69	14.829	1960	1964	1965	rBV4	8143	12218	0.63%	0.054%
70	14.870	1965	1971	1978	rBV	339493	539079	27.80%	2.391%

71	15.100	2006	2010	2019	rVB2	36463	62924	3.24%	0.279%
72	15.223	2027	2031	2035	rBV7	7358	12554	0.65%	0.056%
73	15.370	2054	2056	2058	rVB3	5928	5282	0.27%	0.023%
74	15.493	2071	2077	2081	rBV9	9514	21061	1.09%	0.093%
75	15.693	2105	2111	2117	rBV	951426	1435162	74.00%	6.366%

76	15.746	2117	2120	2123	rVV2	45766	62818	3.24%	0.279%
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77	15.793	2123	2128	2134	rVB	195082	261748	13.50%	1.161%
78	15.951	2151	2155	2160	rVB4	59881	83201	4.29%	0.369%
79	16.040	2168	2170	2173	rVB4	11076	11510	0.59%	0.051%
80	16.557	2255	2258	2264	rVB4	30591	44021	2.27%	0.195%
81	16.674	2274	2278	2282	rVB2	39832	57726	2.98%	0.256%
82	16.980	2326	2330	2334	rBV7	18796	31117	1.60%	0.138%
83	17.256	2374	2377	2382	rVB2	40081	61573	3.17%	0.273%
84	17.320	2385	2388	2391	rBV4	28493	32325	1.67%	0.143%
85	17.356	2391	2394	2400	rVB5	48234	63564	3.28%	0.282%
86	17.444	2402	2409	2413	rBV	723320	1140774	58.82%	5.061%
87	17.485	2413	2416	2421	rVB	427295	561614	28.96%	2.491%
88	17.544	2421	2426	2430	rBV2	1026816	1551722	80.01%	6.884%
89	17.620	2436	2439	2444	rVB6	36203	43964	2.27%	0.195%
90	18.031	2505	2509	2517	rVB4	66018	125325	6.46%	0.556%
91	18.331	2555	2560	2564	rBV3	272950	370780	19.12%	1.645%
92	18.513	2586	2591	2596	rBV4	96571	176561	9.10%	0.783%
93	19.107	2689	2692	2697	rVB5	52143	56320	2.90%	0.250%
94	19.488	2753	2757	2762	rVB	505667	682365	35.18%	3.027%
95	19.600	2773	2776	2785	rBV4	125790	186614	9.62%	0.828%
96	19.823	2808	2814	2817	rBV	1233936	1939464	100.00%	8.604%
97	21.363	3074	3076	3083	rVB3	184786	281576	14.52%	1.249%
98	21.704	3128	3134	3140	rBV3	644383	1244176	64.15%	5.519%
99	24.712	3639	3646	3654	rVB3	568275	1446678	74.59%	6.418%
100	24.935	3678	3684	3693	rVB2	418186	1046141	53.94%	4.641%

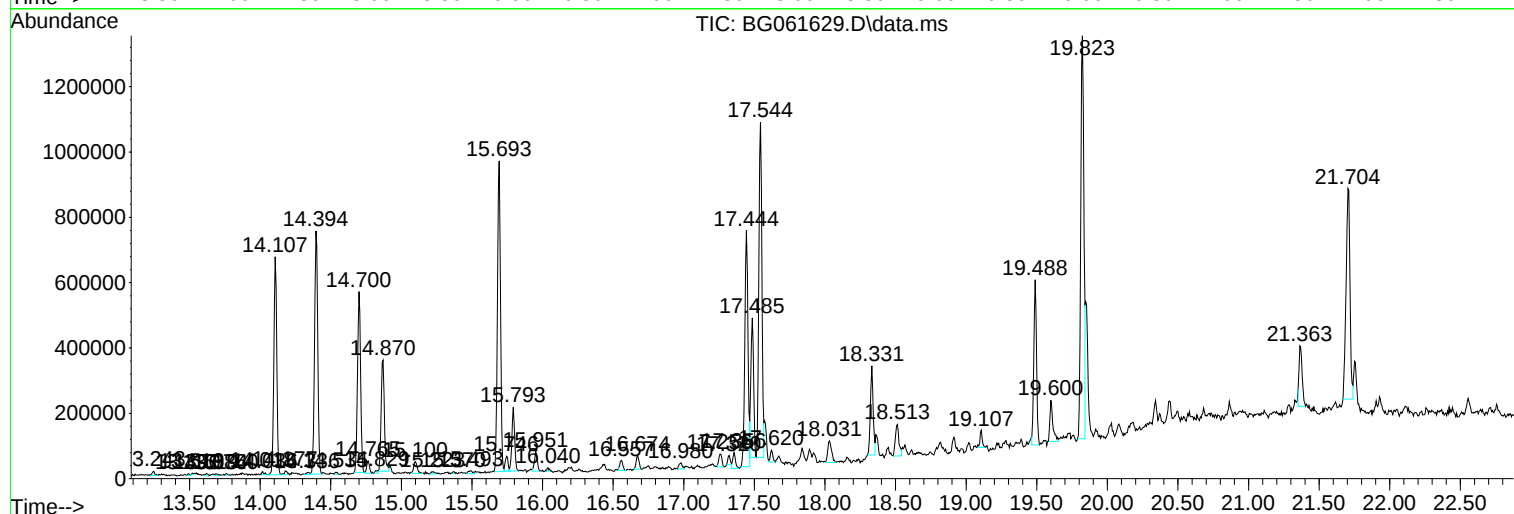
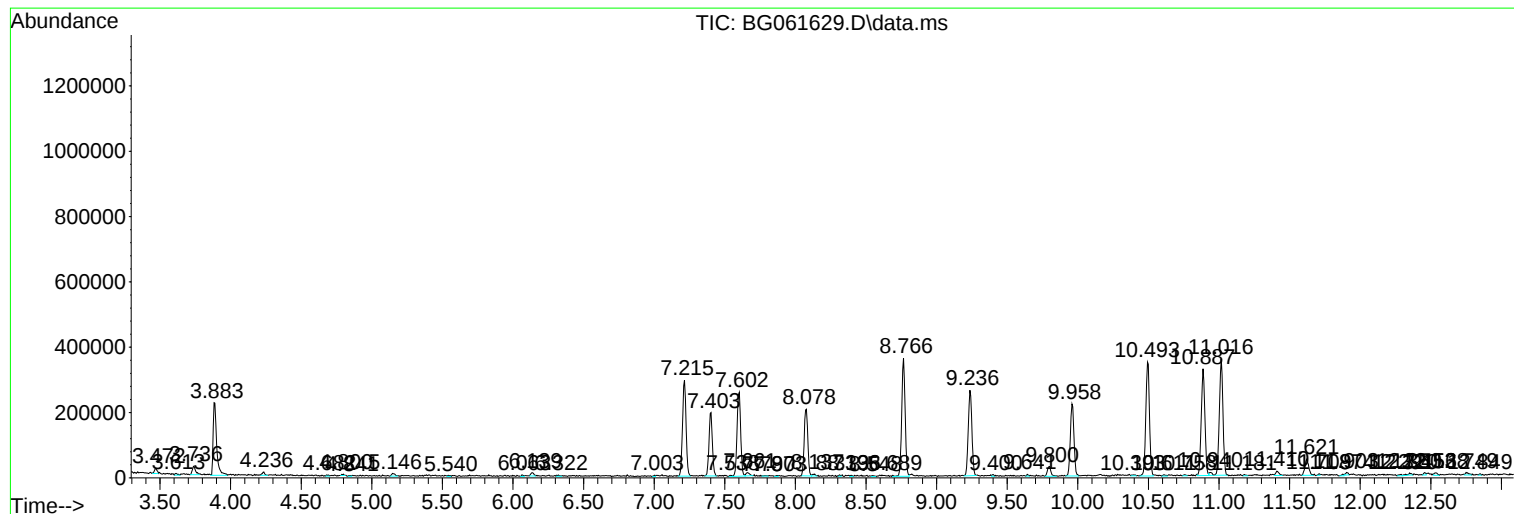
Sum of corrected areas: 22542444

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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



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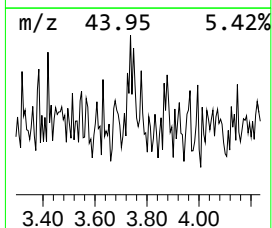
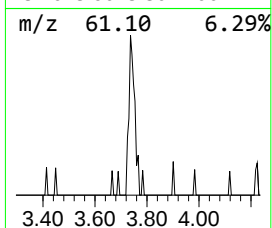
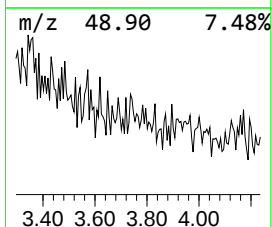
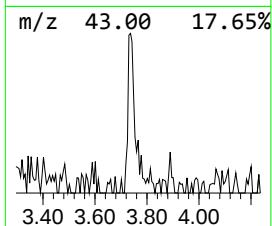
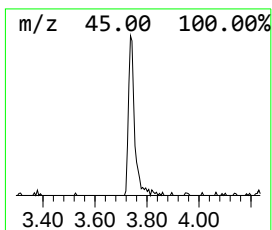
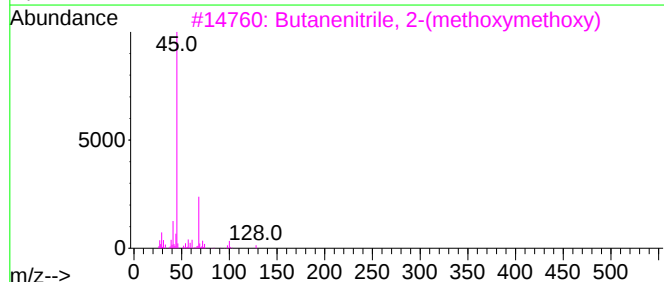
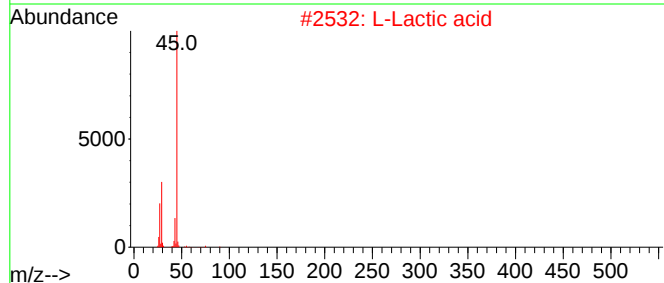
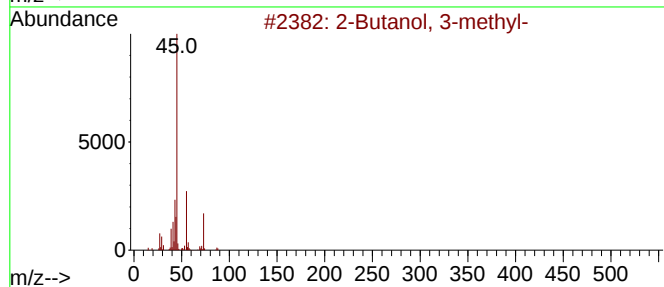
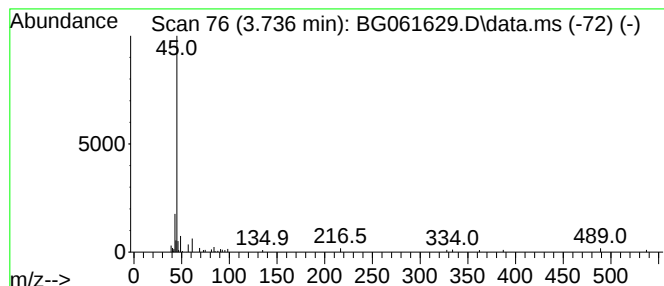
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.736	2.37 ng/ul	42510	1,4-Dichlorobenzene-d4	8.072

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 3-methyl-			88	C5H12O	000598-75-4	39
2	L-Lactic acid			90	C3H6O3	000079-33-4	9
3	Butanenitrile, 2-(methoxymethoxy)			129	C6H11NO2	1000196-86-5	9
4	Methoxyacetic acid, pentyl ester			160	C8H16O3	168920-35-2	9
5	2-Heptanol, (S)-			116	C7H16O	006033-23-4	9



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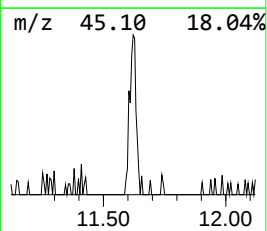
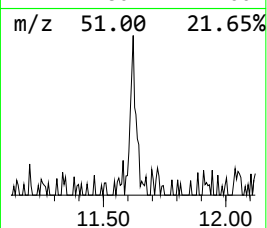
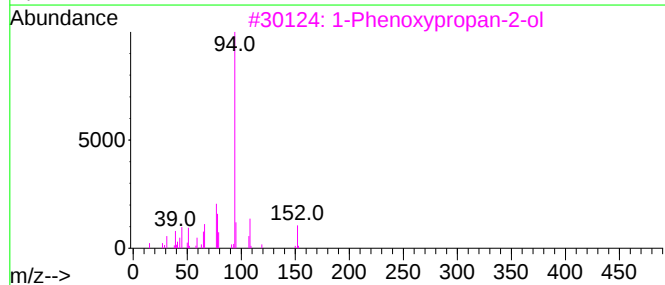
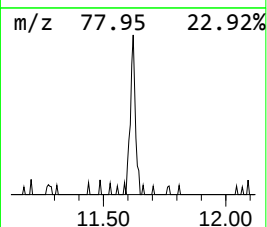
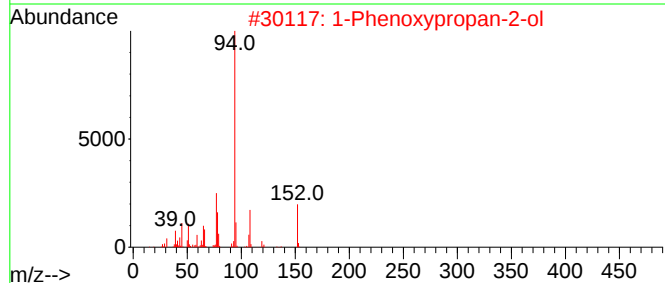
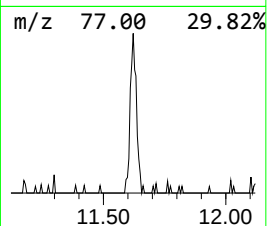
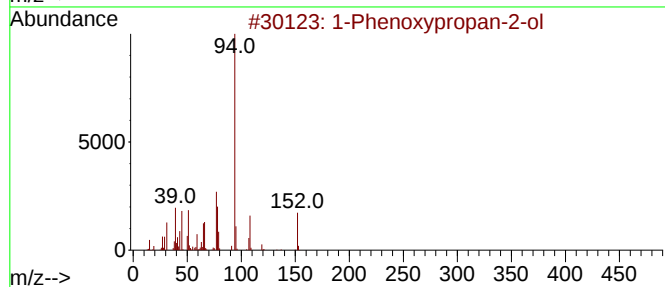
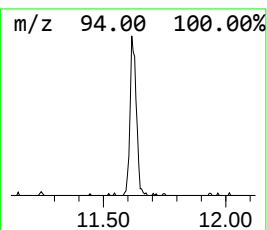
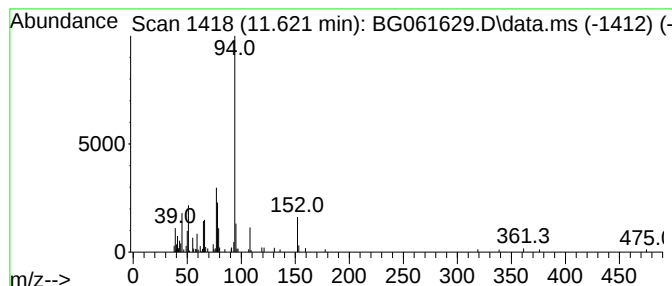
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.621	3.20 ng/ul	93224	Naphthalene-d8	10.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			2-Phenoxybutyric acid	180	C10H12O3	013794-14-4	59
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	58



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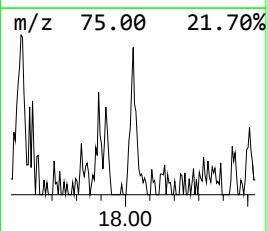
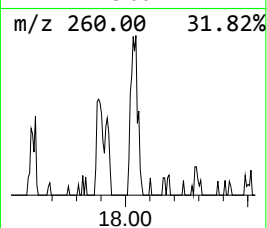
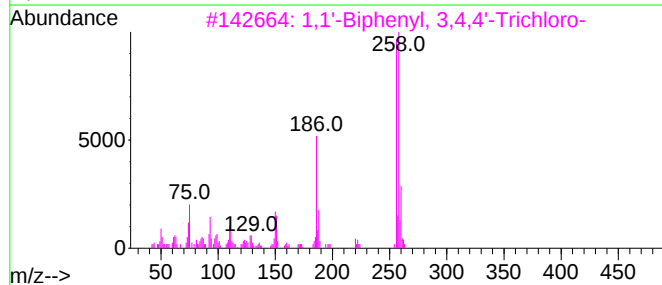
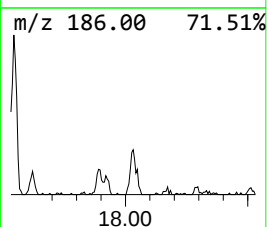
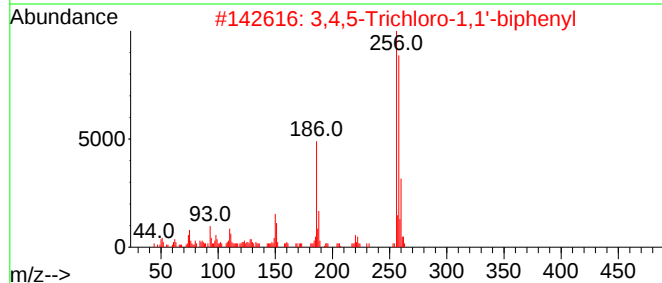
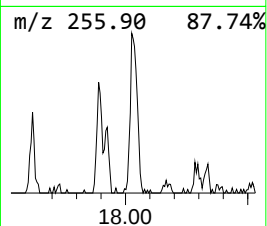
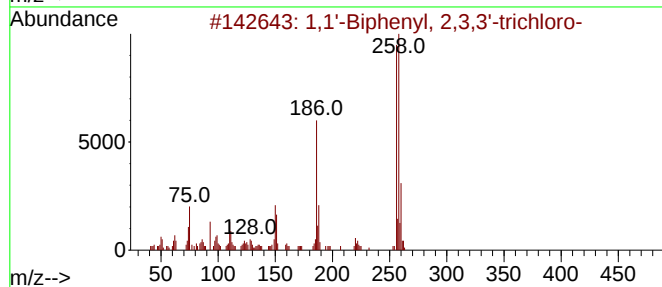
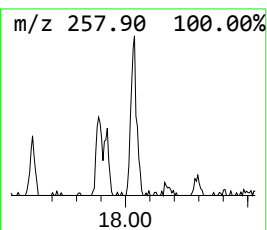
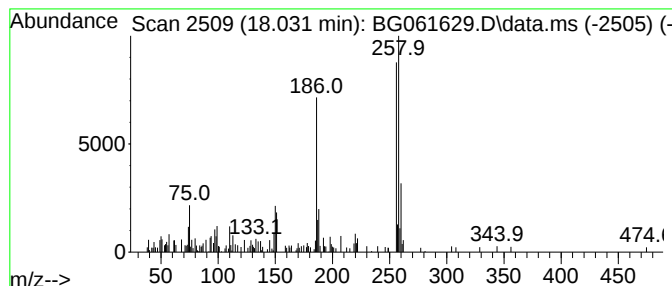
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
18.031	2.20 ng/ul	125325	Phenanthrene-d10			17.444
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-		256	C12H7Cl3	038444-84-7	99
2	3,4,5-Trichloro-1,1'-biphenyl		256	C12H7Cl3	053555-66-1	98
3	1,1'-Biphenyl, 3,4,4'-Trichloro-		256	C12H7Cl3	038444-90-5	98
4	1,1'-Biphenyl, 3,3',5-trichloro-		256	C12H7Cl3	038444-87-0	97
5	2,4',6-Trichloro-1,1'-biphenyl		256	C12H7Cl3	038444-77-8	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062124\
Data File : BG061629.D
Acq On : 21 Jun 2024 20:58
Operator : MA/JU
Sample : P2754-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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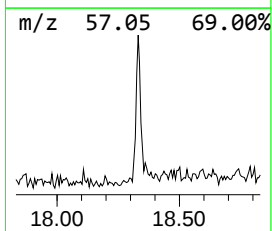
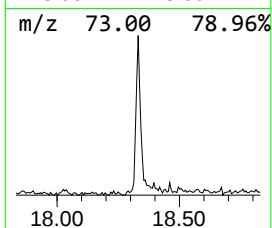
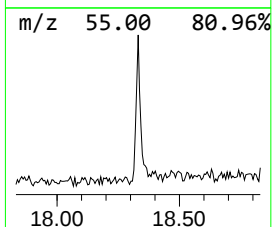
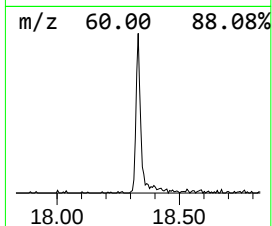
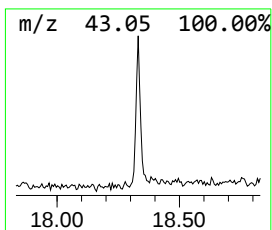
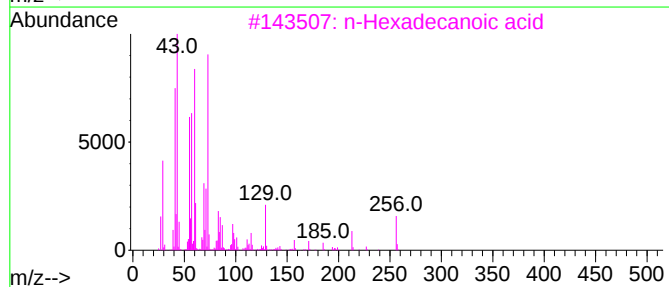
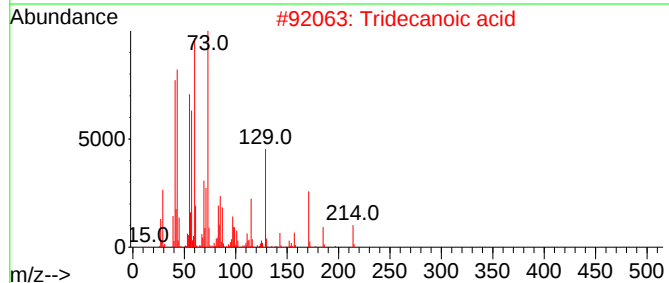
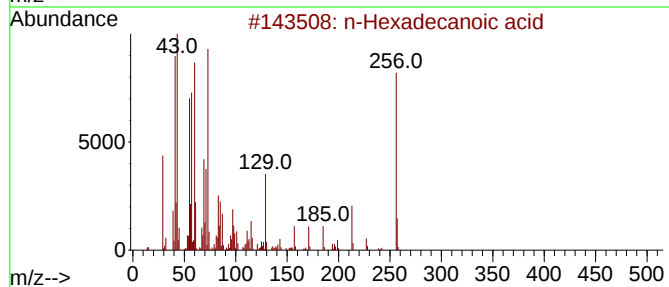
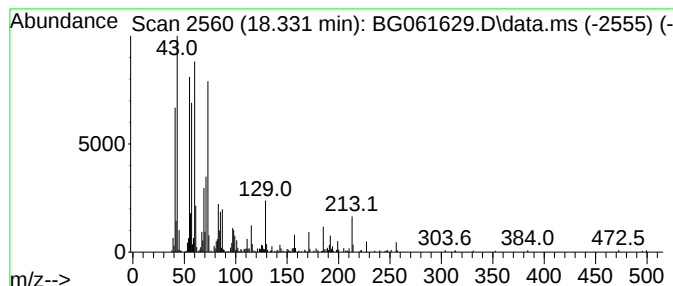
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.331	6.50 ng/ul	370780	Phenanthrene-d10	17.444

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Tridecanoic acid	214	C13H26O2	000638-53-9	89
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	74
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062124\
Data File : BG061629.D
Acq On : 21 Jun 2024 20:58
Operator : MA/JU
Sample : P2754-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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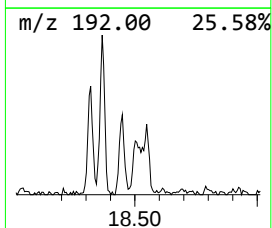
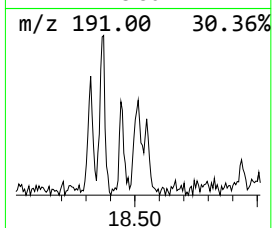
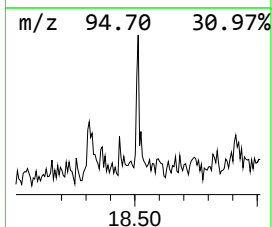
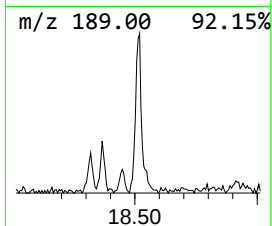
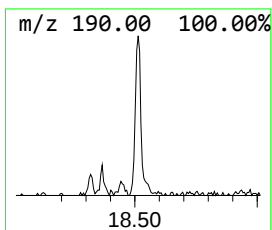
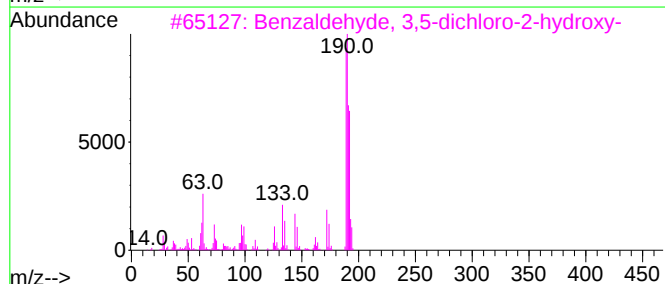
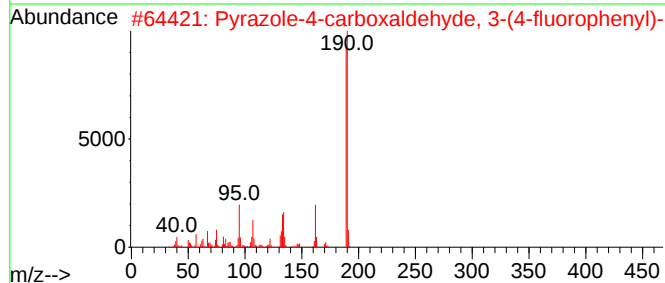
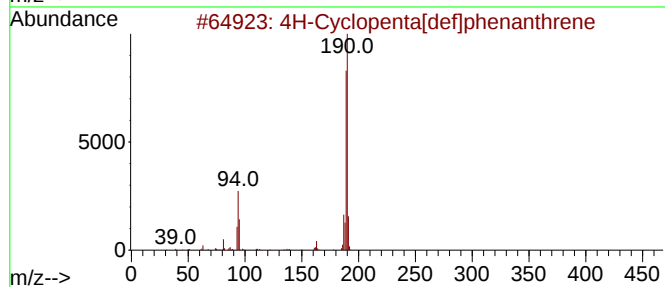
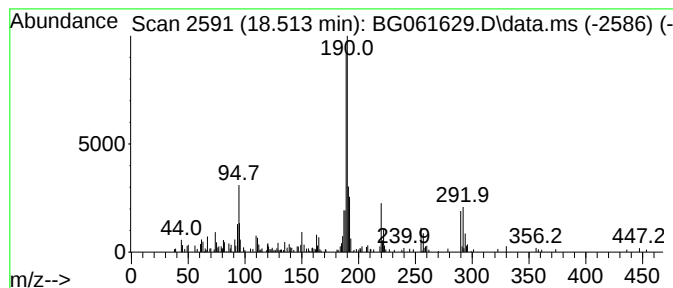
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.513	3.10 ng/ul	176561	Phenanthrene-d10	17.444

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
2			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	43
3			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	43
4			Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	43
5			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062124\
Data File : BG061629.D
Acq On : 21 Jun 2024 20:58
Operator : MA/JU
Sample : P2754-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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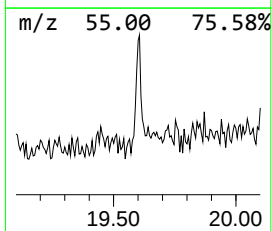
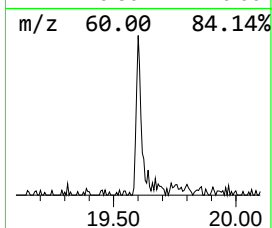
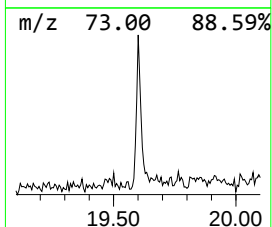
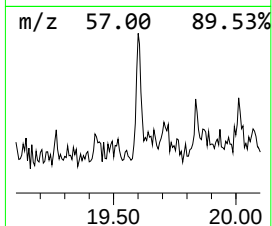
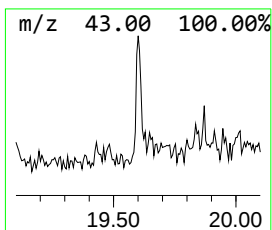
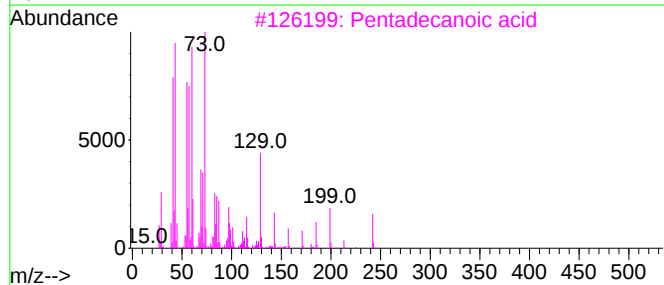
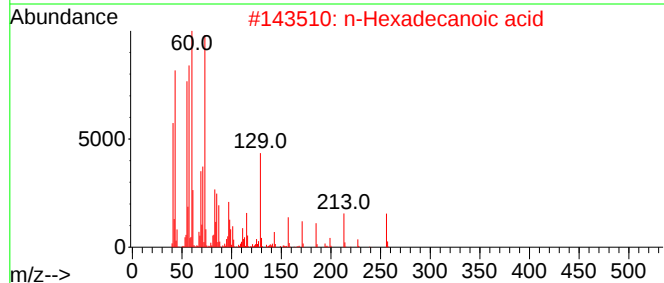
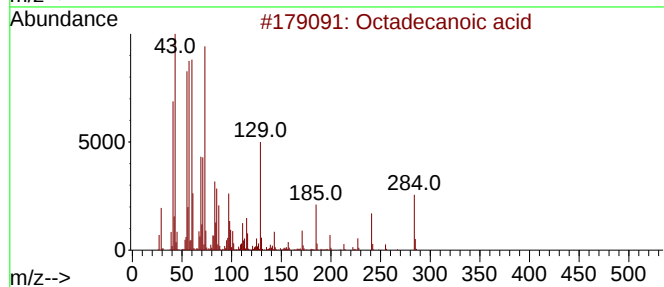
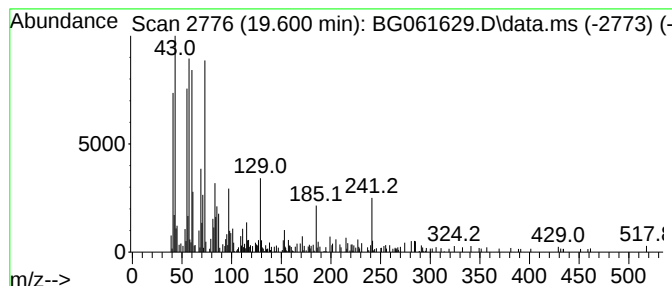
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.600	3.00 ng/ul	186614	Chrysene-d12	21.709

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4			Octadecanoic acid	284	C18H36O2	000057-11-4	83
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062124\
Data File : BG061629.D
Acq On : 21 Jun 2024 20:58
Operator : MA/JU
Sample : P2754-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

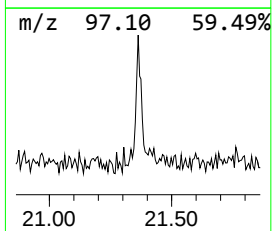
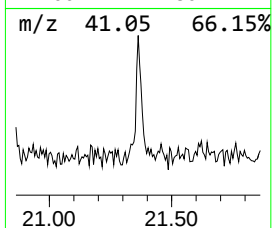
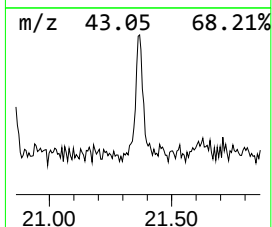
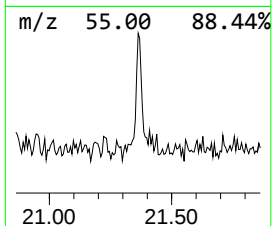
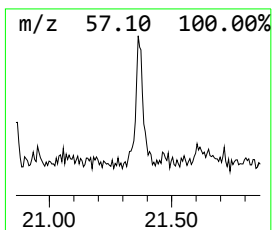
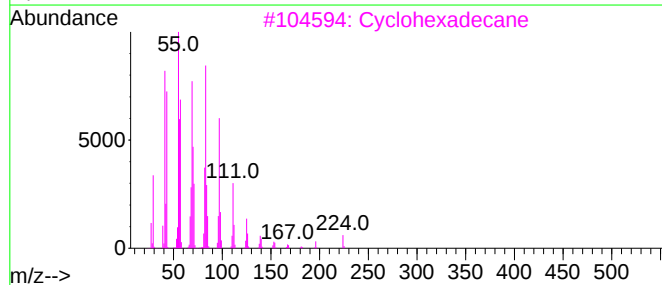
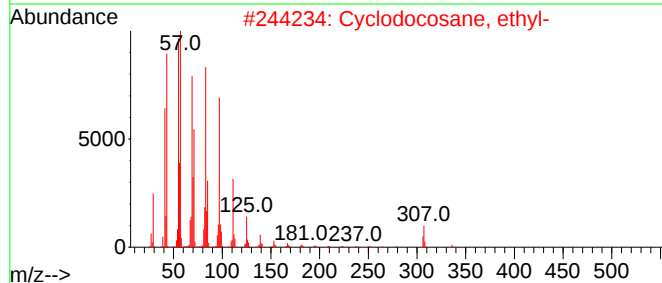
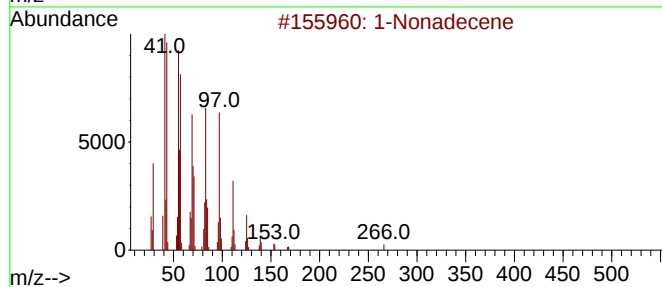
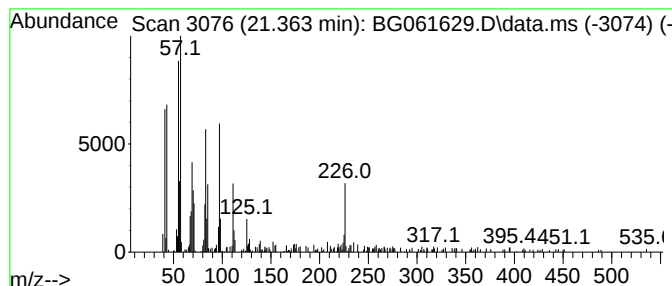
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.363	4.53 ng/ul	281576	Chrysene-d12		21.709	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Nonadecene		266	C19H38	018435-45-5	93
2	Cyclodocosane, ethyl-		336	C24H48	1000151-22-6	89
3	Cyclohexadecane		224	C16H32	000295-65-8	89
4	Z-5-Nonadecene		266	C19H38	1000131-11-8	86
5	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062124\
Data File : BG061629.D
Acq On : 21 Jun 2024 20:58
Operator : MA/JU
Sample : P2754-11
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4BQ0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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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1-Phenoxypropan...	11.621	3.2	ng/ul	93224	2	10.887	582597	20.0
1,1'-Biphenyl, ...	18.031	2.2	ng/ul	125325	4	17.444	1140770	20.0
n-Hexadecanoic ...	18.331	6.5	ng/ul	370780	4	17.444	1140770	20.0
4H-Cyclopenta[d...	18.513	3.1	ng/ul	176561	4	17.444	1140770	20.0
Octadecanoic acid	19.600	3.0	ng/ul	186614	5	21.709	1244180	20.0
(DEL) Alkane: S...	21.363	4.5	ng/ul	281576	5	21.709	1244180	20.0