

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
 Data File : BG061681.D
 Acq On : 24 Jun 2024 18:55
 Operator : MA/JU
 Sample : P2776-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4CY6

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: BG061681.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.467	27	30	33	rBV4	12708	15289	0.75%	0.069%
2	3.737	72	76	83	rVV	29664	47308	2.33%	0.212%
3	3.896	101	103	108	rVV3	23654	36705	1.81%	0.165%
4	4.001	118	121	125	rVB6	7953	9990	0.49%	0.045%
5	4.054	128	130	134	rBV4	3153	4706	0.23%	0.021%
6	4.113	137	140	142	rBV3	3424	3882	0.19%	0.017%
7	4.231	158	160	165	rVB3	9022	11344	0.56%	0.051%
8	4.601	219	223	228	rBV6	4625	8199	0.40%	0.037%
9	4.742	241	247	251	rBV5	9566	16912	0.83%	0.076%
10	4.795	251	256	261	rVV2	13074	22020	1.08%	0.099%
11	5.147	311	316	321	rBV	40141	60190	2.97%	0.270%
12	5.388	355	357	360	rVB3	4898	4211	0.21%	0.019%
13	5.459	368	369	374	rBV2	2839	4035	0.20%	0.018%
14	5.817	427	430	433	rBV4	3263	4691	0.23%	0.021%
15	6.134	481	484	492	rVB2	14176	19875	0.98%	0.089%
16	6.510	544	548	549	rVB	3616	4011	0.20%	0.018%
17	6.675	574	576	580	rVB	5353	7466	0.37%	0.033%
18	7.215	662	668	682	rVB	295656	502451	24.75%	2.254%
19	7.351	686	691	693	rBV2	7037	10511	0.52%	0.047%
20	7.398	693	699	708	rVB	189933	303809	14.97%	1.363%
21	7.597	727	733	740	rVV	245841	414908	20.44%	1.862%
22	7.656	740	743	747	rVV4	5917	7599	0.37%	0.034%
23	7.862	774	778	783	rVV5	2107	4199	0.21%	0.019%
24	8.073	808	814	821	rBV	236142	378951	18.67%	1.700%
25	8.126	821	823	827	rVB4	6642	9141	0.45%	0.041%
26	8.520	885	890	894	rBV5	3353	7241	0.36%	0.032%
27	8.643	909	911	913	rBV2	3877	4076	0.20%	0.018%
28	8.766	925	932	941	rBV	324223	543474	26.77%	2.438%
29	8.896	950	954	960	rVV4	9559	18564	0.91%	0.083%
30	8.966	963	966	969	rVB2	3166	4342	0.21%	0.019%
31	9.119	989	992	993	rBV	5188	5069	0.25%	0.023%
32	9.237	1005	1012	1021	rVB	255557	443088	21.83%	1.988%
33	9.307	1021	1024	1027	rBV3	4735	5509	0.27%	0.025%
34	9.395	1035	1039	1044	rBV5	3034	6075	0.30%	0.027%
35	9.442	1045	1047	1052	rVB3	2737	3800	0.19%	0.017%
36	9.642	1079	1081	1083	rBV2	5490	4424	0.22%	0.020%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

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Max Peaks: 100

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
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37	9.724	1092	1095	1098	rVB4	3561	5228	0.26%	0.023%
38	9.795	1102	1107	1113	rVB3	37434	58588	2.89%	0.263%
39	9.959	1127	1135	1144	rBV	223131	395644	19.49%	1.775%
40	10.083	1154	1156	1159	rVB2	5120	4667	0.23%	0.021%

41	10.500	1219	1227	1235	rVB	341071	615897	30.34%	2.763%
42	10.647	1247	1252	1257	rBV4	16956	30885	1.52%	0.139%
43	10.788	1273	1276	1279	rBV2	3842	4232	0.21%	0.019%
44	10.888	1285	1293	1300	rVV	362654	639865	31.52%	2.871%
45	11.017	1307	1315	1326	rBV	198686	378149	18.63%	1.697%

46	11.346	1368	1371	1373	rBV2	3987	4262	0.21%	0.019%
47	11.416	1376	1383	1389	rVB6	19800	32173	1.58%	0.144%
48	11.622	1412	1418	1427	rVV2	59460	121205	5.97%	0.544%
49	11.775	1441	1444	1447	rBV2	2933	4546	0.22%	0.020%
50	12.398	1546	1550	1551	rBV4	8506	8303	0.41%	0.037%

51	12.462	1557	1561	1565	rBV3	6898	8620	0.42%	0.039%
52	12.680	1596	1598	1602	rVB3	5100	6094	0.30%	0.027%
53	12.797	1615	1618	1622	rVB3	4383	6321	0.31%	0.028%
54	12.844	1622	1626	1628	rBV4	5499	6827	0.34%	0.031%
55	13.009	1652	1654	1658	rBV4	3818	4033	0.20%	0.018%

56	13.255	1694	1696	1700	rVB3	4837	4174	0.21%	0.019%
57	13.526	1739	1742	1748	rVB3	7506	12200	0.60%	0.055%
58	13.602	1754	1755	1759	rBV4	5504	5916	0.29%	0.027%
59	14.107	1834	1841	1847	rVB	631369	936901	46.16%	4.204%
60	14.178	1850	1853	1855	rBV2	5479	5733	0.28%	0.026%

61	14.248	1864	1865	1870	rVB3	4310	4992	0.25%	0.022%
62	14.295	1870	1873	1874	rBV2	4933	3899	0.19%	0.017%
63	14.348	1877	1882	1884	rBV5	6036	10278	0.51%	0.046%
64	14.395	1884	1890	1897	rVV	711300	1113921	54.88%	4.998%
65	14.601	1920	1925	1928	rBV5	7197	9921	0.49%	0.045%

66	14.701	1936	1942	1949	rVB	617344	946516	46.63%	4.247%
67	14.765	1949	1953	1960	rVB6	13122	25541	1.26%	0.115%
68	14.824	1960	1963	1965	rBV3	4268	6540	0.32%	0.029%
69	14.877	1965	1972	1977	rBV	337621	605842	29.85%	2.718%
70	15.036	1995	1999	2005	rVB6	22365	29700	1.46%	0.133%

71	15.112	2007	2012	2017	rVB7	26914	51189	2.52%	0.230%
72	15.159	2017	2020	2021	rBV3	6707	5874	0.29%	0.026%
73	15.582	2090	2092	2093	rBV	5410	4483	0.22%	0.020%
74	15.694	2104	2111	2117	rBV	941638	1462409	72.04%	6.562%
75	15.747	2118	2120	2123	rVV3	14773	13447	0.66%	0.060%

76	15.794	2123	2128	2135	rVB	287762	433151	21.34%	1.943%
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 Title : SVOA CALIBRATION

77	15.905	2145	2147	2148	rBV2	5905	4279	0.21%	0.019%
78	15.993	2161	2162	2163	rBV	7066	3913	0.19%	0.018%
79	16.146	2186	2188	2191	rVB3	4504	4026	0.20%	0.018%
80	16.410	2227	2233	2240	rBV8	11278	22018	1.08%	0.099%
81	16.504	2247	2249	2252	rBV4	4148	4470	0.22%	0.020%
82	17.445	2403	2409	2414	rBV	866811	1308209	64.45%	5.870%
83	17.486	2414	2416	2420	rVV	118580	146712	7.23%	0.658%
84	17.544	2420	2426	2435	rVB2	1060678	1583414	78.01%	7.104%
85	18.044	2505	2511	2515	rBV3	72998	123374	6.08%	0.554%
86	18.197	2535	2537	2538	rBV2	9675	7387	0.36%	0.033%
87	18.332	2555	2560	2565	rBV	297450	438373	21.60%	1.967%
88	18.508	2586	2590	2594	rBV6	62233	88203	4.35%	0.396%
89	18.614	2605	2608	2614	rBV7	25327	45910	2.26%	0.206%
90	18.960	2665	2667	2671	rBV5	18942	21835	1.08%	0.098%
91	19.360	2732	2735	2740	rVB4	54142	75501	3.72%	0.339%
92	19.489	2753	2757	2763	rVB	266231	391757	19.30%	1.758%
93	19.595	2773	2775	2779	rBV3	59300	66931	3.30%	0.300%
94	19.824	2808	2814	2825	rBV2	1172514	2029860	100.00%	9.108%
95	21.363	3072	3076	3083	rVB	447597	661041	32.57%	2.966%
96	21.710	3129	3135	3140	rBV2	722922	1253595	61.76%	5.625%
97	22.550	3274	3278	3292	rVB3	229915	520614	25.65%	2.336%
98	24.078	3533	3538	3542	rBV	220400	380800	18.76%	1.709%
99	24.718	3640	3647	3654	rBV3	381253	937241	46.17%	4.205%
100	24.942	3678	3685	3695	rVB2	425751	1181905	58.23%	5.303%

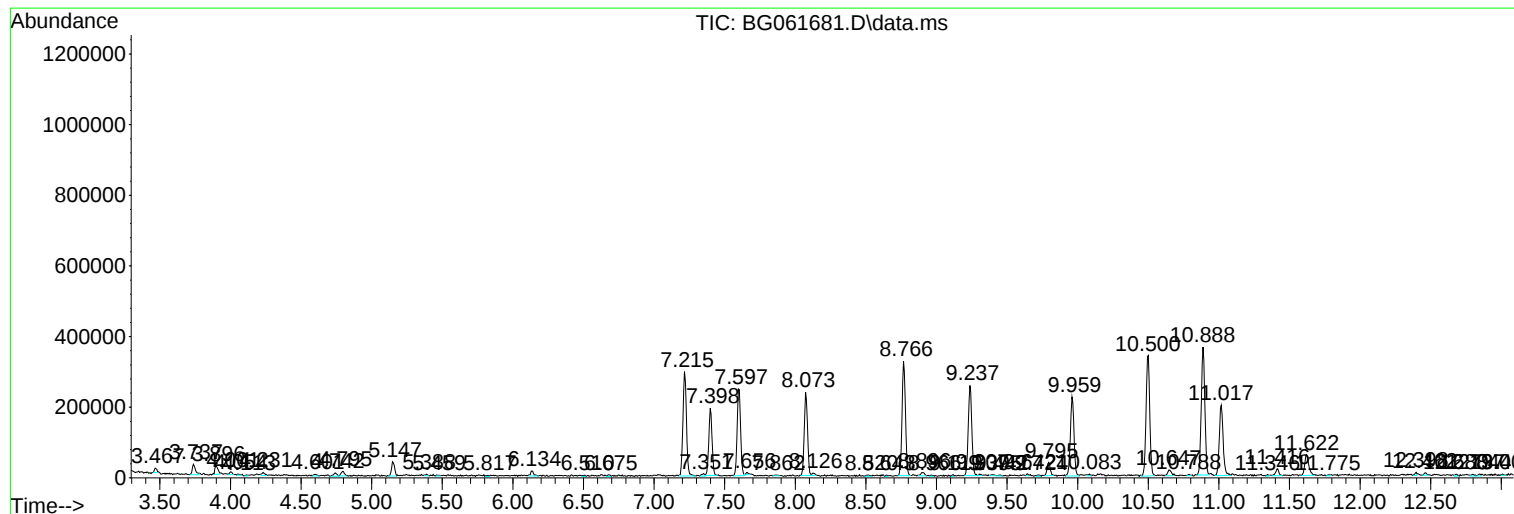
Sum of corrected areas: 22287599

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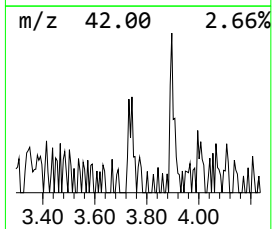
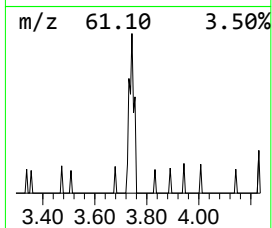
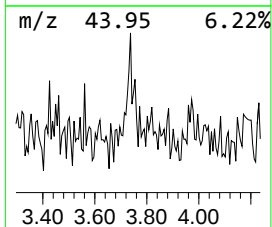
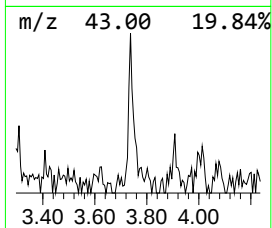
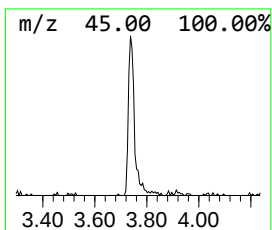
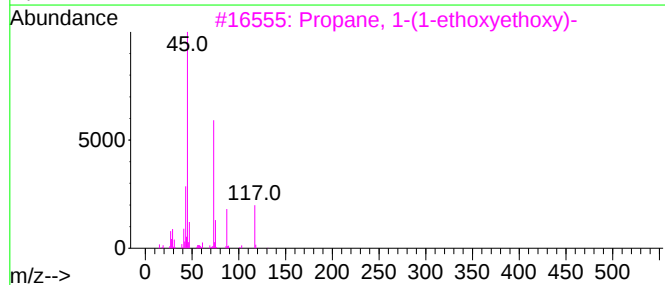
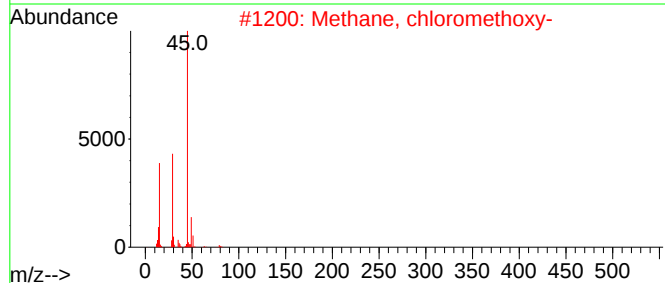
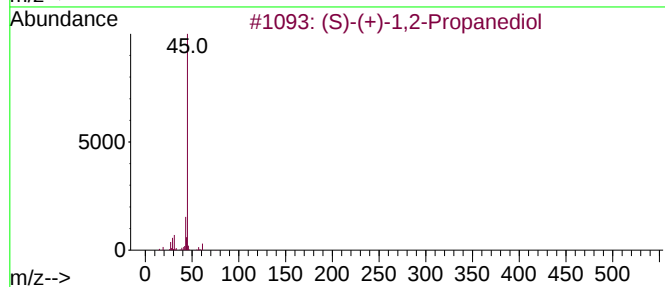
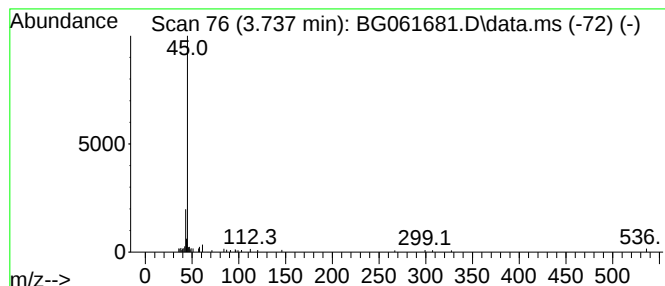
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (S)-(+)-1,2-Propanediol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.737	2.50 ng/ul	47308	1,4-Dichlorobenzene-d4	8.073

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	50
2		Methane, chloromethoxy-	80	C2H5ClO	000107-30-2	39
3		Propane, 1-(1-ethoxyethoxy)-	132	C7H16O2	020680-10-8	9
4		2-Propanol, 1,1,1-trichloro-	162	C3H5Cl3O	000076-00-6	9
5		Propylene Glycol	76	C3H8O2	000057-55-6	9



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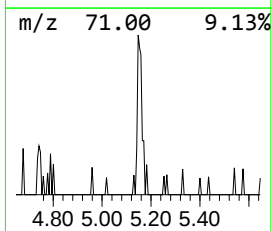
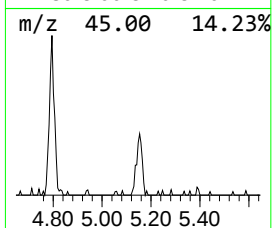
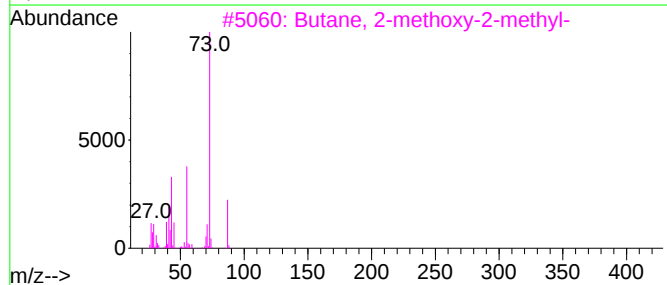
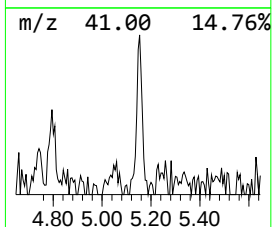
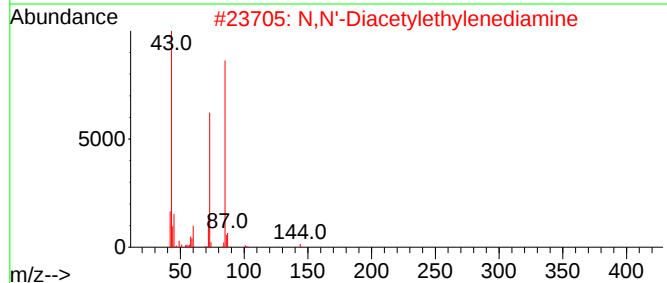
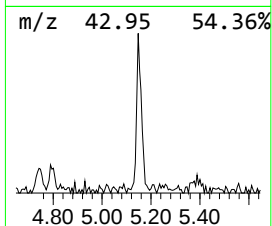
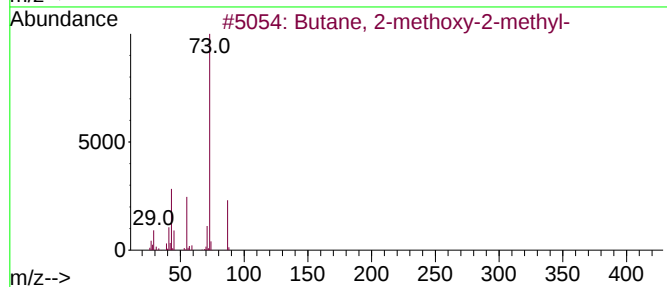
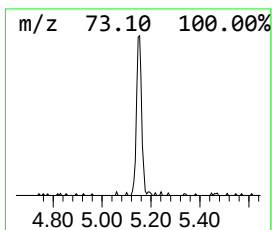
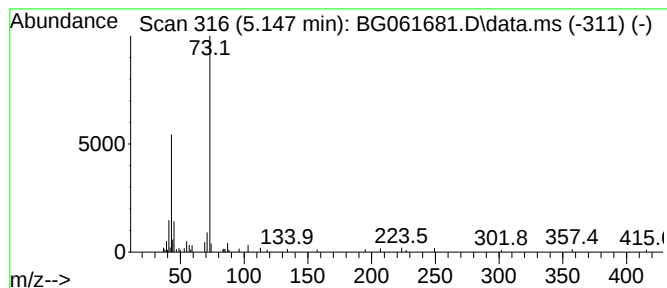
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.147	3.18 ng/ul	60190	1,4-Dichlorobenzene-d4	8.073

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	40
2			N,N'-Diacetylenediamine	144	C6H12N2O2	000871-78-3	38
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	33
4			Ether, 6-bromo-1-ethyloctyl methyl	250	C11H23BrO	020599-95-5	33
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	32



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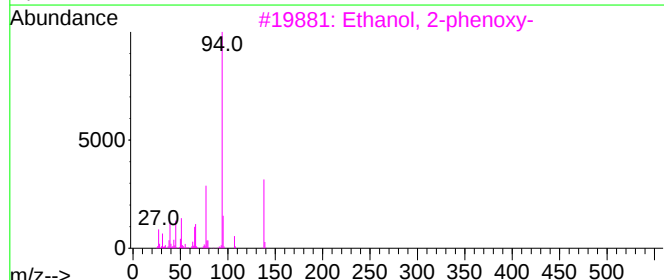
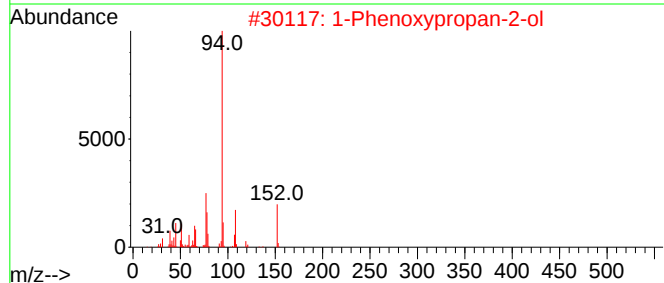
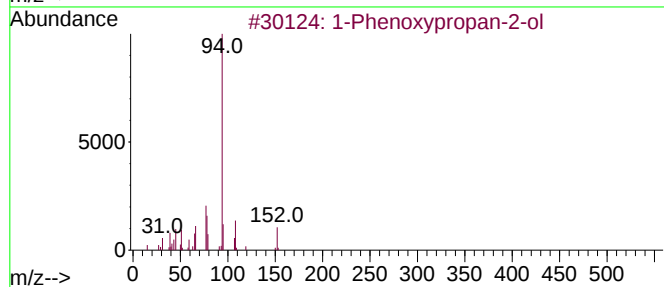
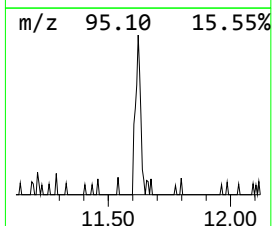
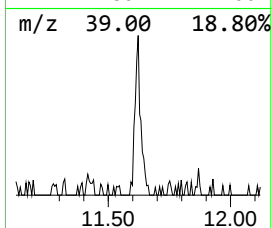
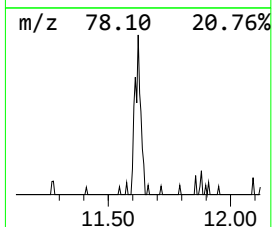
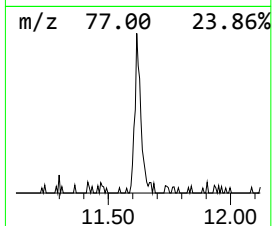
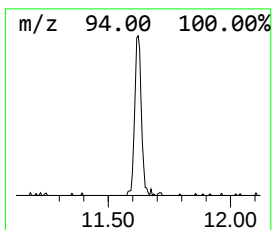
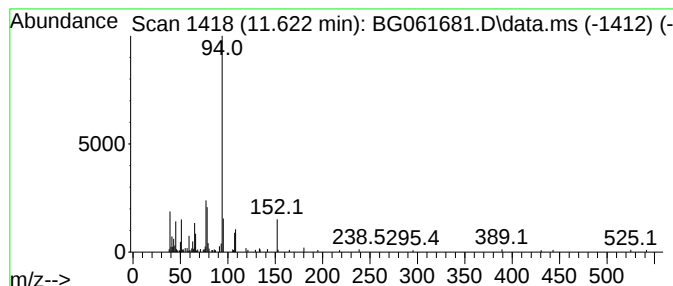
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.622	3.79 ng/ul	121205	Naphthalene-d8	10.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	60
3			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	53
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	53
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061681.D
Acq On : 24 Jun 2024 18:55
Operator : MA/JU
Sample : P2776-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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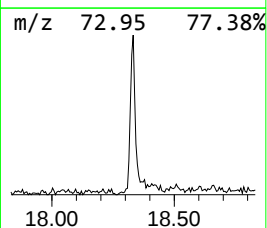
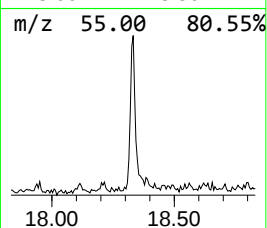
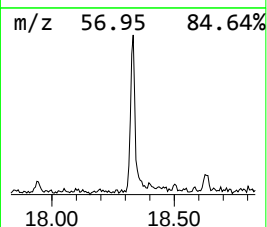
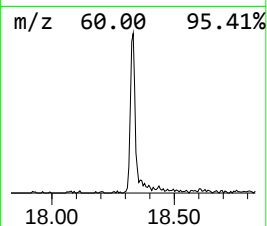
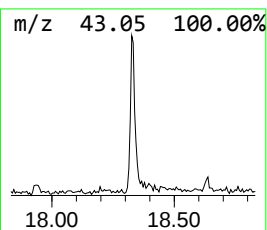
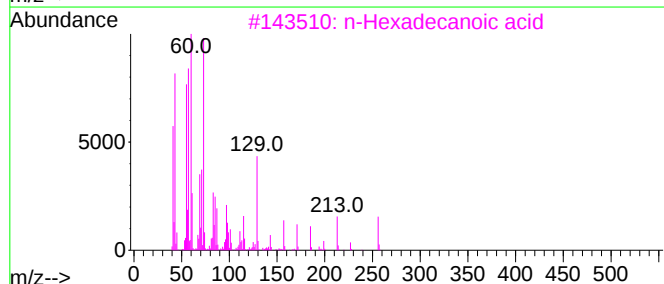
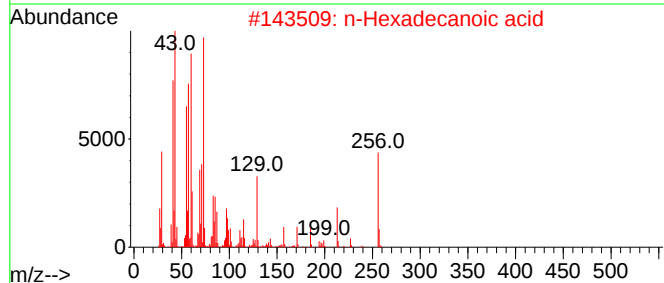
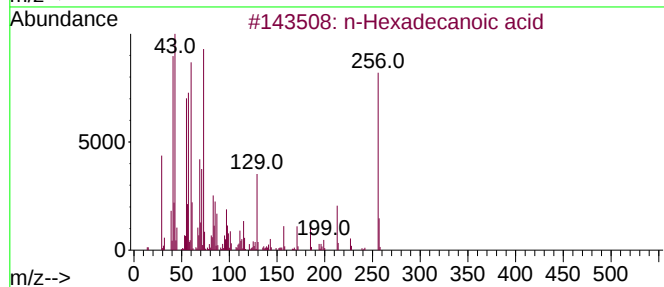
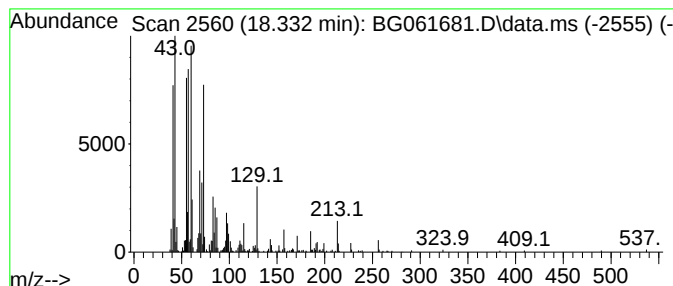
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.332	6.70 ng/ul	438373	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	78
5			Tridecanoic acid	214	C13H26O2	000638-53-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061681.D
Acq On : 24 Jun 2024 18:55
Operator : MA/JU
Sample : P2776-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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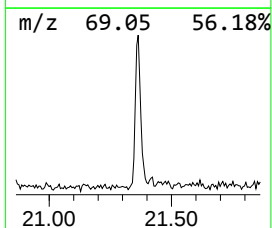
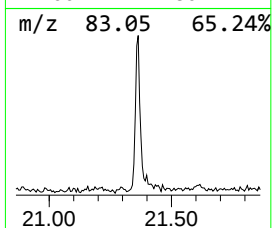
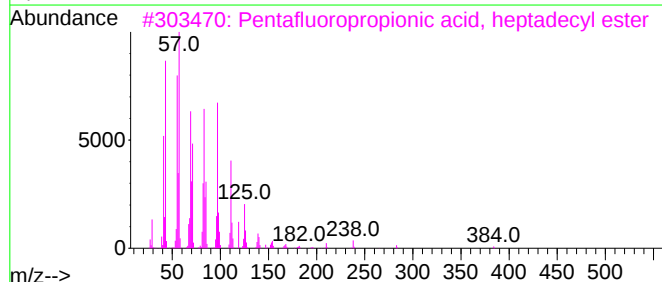
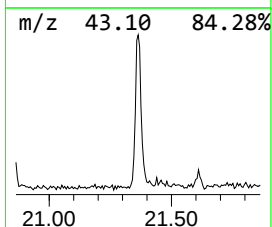
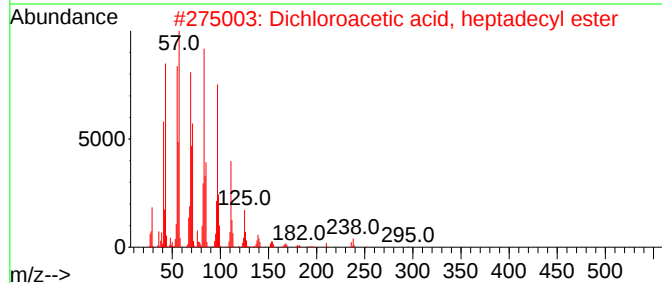
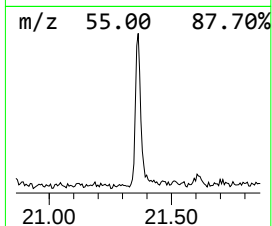
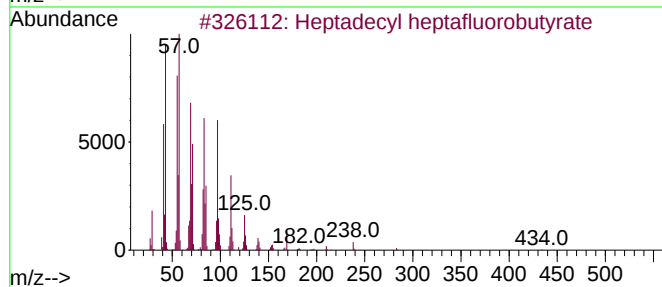
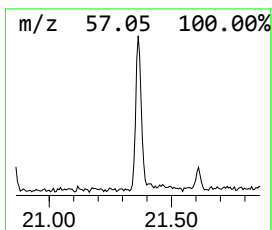
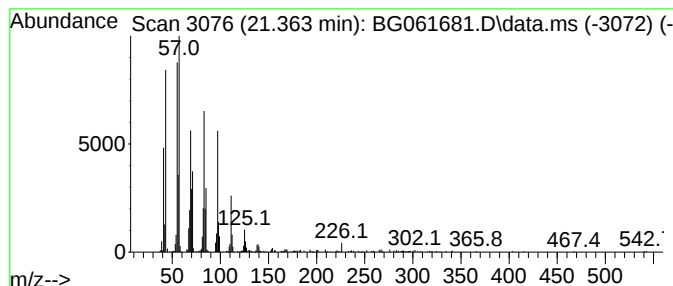
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Heptadecyl heptafluorobutyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.363	10.55 ng/ul	661041	Chrysene-d12	21.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
2			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
3			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
4			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061681.D
Acq On : 24 Jun 2024 18:55
Operator : MA/JU
Sample : P2776-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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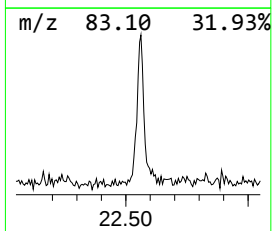
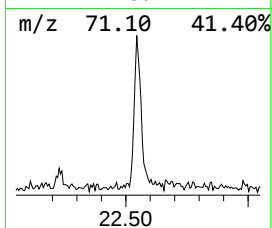
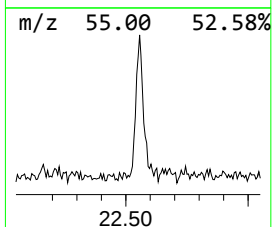
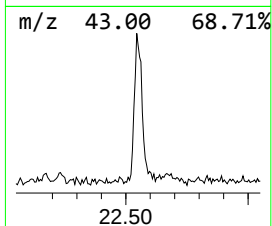
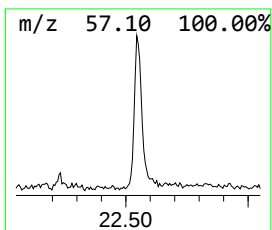
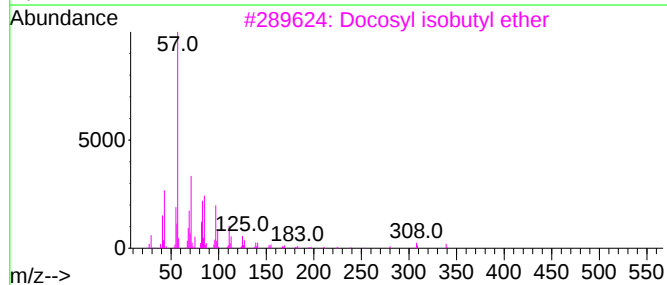
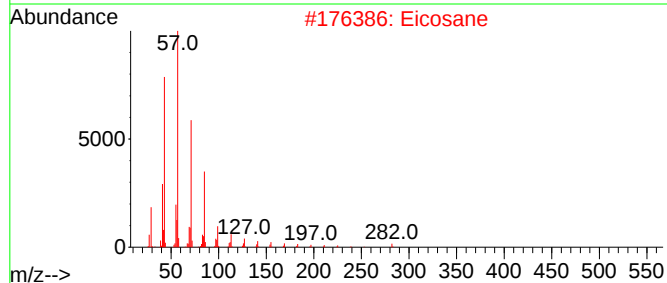
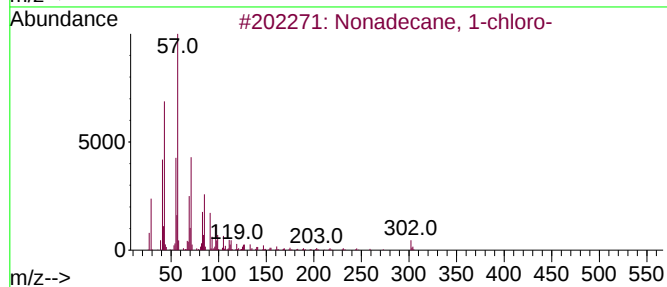
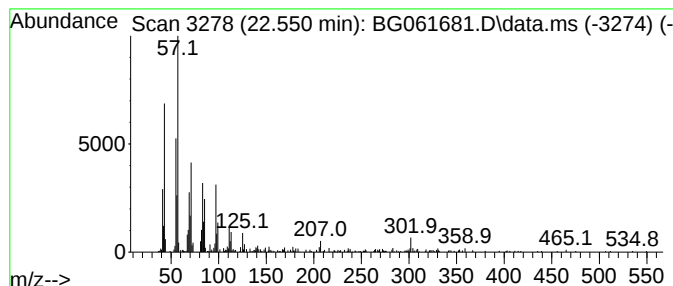
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Nonadecane, 1-chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.550	8.31 ng/ul	520614	Chrysene-d12	21.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	90
2			Eicosane	282	C20H42	000112-95-8	90
3			Docosyl isobutyl ether	382	C26H54O	1000406-33-1	72
4			Isobutyl tetracosyl ether	410	C28H58O	1000406-33-2	68
5			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061681.D
Acq On : 24 Jun 2024 18:55
Operator : MA/JU
Sample : P2776-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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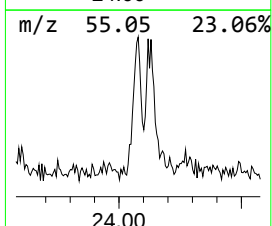
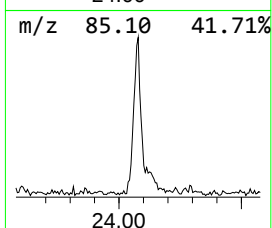
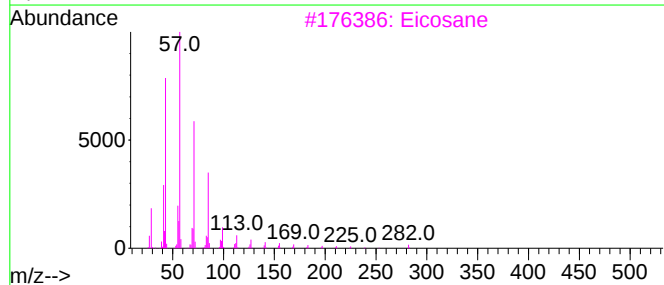
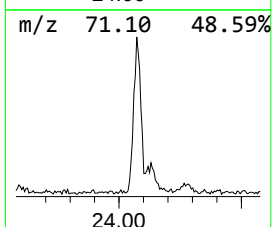
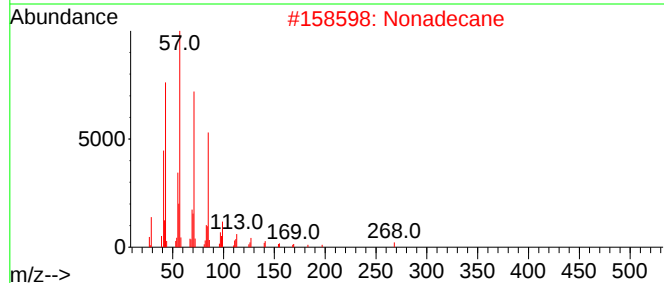
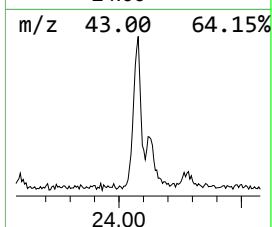
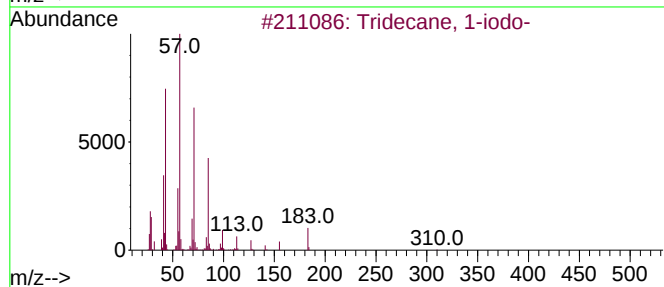
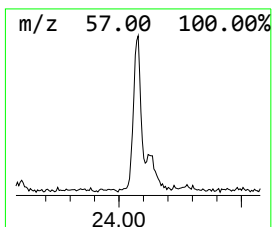
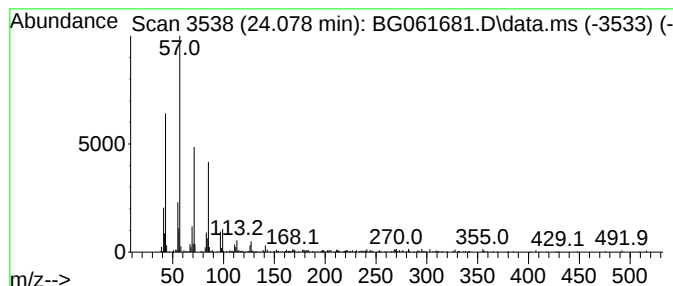
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tridecane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.078	6.44 ng/ul	380800	Perylene-d12	24.942

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
2			Nonadecane	268	C19H40	000629-92-5	93
3			Eicosane	282	C20H42	000112-95-8	91
4			Tricosane, 2-methyl-	338	C24H50	001928-30-9	90
5			Octacosane	394	C28H58	000630-02-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061681.D
Acq On : 24 Jun 2024 18:55
Operator : MA/JU
Sample : P2776-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
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unknown-01	5.147	3.2	ng/ul	60190	1	8.073	378951	20.0
1-Phenoxypropan...	11.622	3.8	ng/ul	121205	2	10.887	639865	20.0
n-Hexadecanoic ...	18.332	6.7	ng/ul	438373	4	17.445	1308210	20.0
Heptadecyl hept...	21.363	10.6	ng/ul	661041	5	21.710	1253600	20.0
Nonadecane, 1-C...	22.550	8.3	ng/ul	520614	5	21.710	1253600	20.0
Tridecane, 1-iodo-	24.078	6.4	ng/ul	380800	6	24.942	1181910	20.0