LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : B0AA8

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_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

Signal : TIC: BP013220.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	3	7	18	rVV	1313521	1607865	14.29%	1.453%
2	3.334	22	25	34	rVB	128547	163541	1.45%	0.148%
3	3.434	39	42	46	rVB2	34430	35697	0.32%	0.032%
4	3.611	69	72	91	rBV	794150	1159977	10.31%	1.048%
5	3.770	96	99	109	rBV	526908	760320	6.76%	0.687%
6	3.993	134	137	142	rVB	28622	36939	0.33%	0.033%
7	4.093	151	154	160	rVB	36785	50651	0.45%	0.046%
8	4.569	231	235	236	rBV	27844	31210	0.28%	0.028%
9	4.605	236	241	248	rVB	3763794	5327406	47.34%	4.813%
10	4.769	267	269	275	rVB4	7937	13608	0.12%	0.012%
11	4.858	281	284	287	rBV4	16705	21023	0.19%	0.019%
12	4.905	287	292	300	rVB	375664	533075	4.74%	0.482%
13	5.040	309	315	321	rBV	91046	138651	1.23%	0.125%
14	5.116	321	328	333	rVB	181255	265851	2.36%	0.240%
15	6.193	507	511	516	rVB6	19584	30002	0.27%	0.027%
16	6.258	517	522	528	rBV	213207	313060	2.78%	0.283%
17	6.334	533	535	541	rVB7	6864	11528	0.10%	0.010%
18	7.105	661	666	678	rVB	241041	391936	3.48%	0.354%
19	7.275	688	695	704	rBV	1327068	2023409	17.98%	1.828%
20	7.358	704	709	716	rVB2	32966	52691	0.47%	0.048%
21	7.475	722	729	739	rBV	1836375	2883166	25.62%	2.605%
22	7.558	742	743	751	rVB8	11849	16528	0.15%	0.015%
23	7.658	756	760	767	rVB3	22115	36896	0.33%	0.033%
24	7.858	789	794	802	rBV4	21462	43203	0.38%	0.039%
25	7.952	802	810	816	rVV	1606437	2571937	22.86%	2.324%
26	8.010	816	820	827	rVB	43934	70881	0.63%	0.064%
27	8.569	908	915	921	rVB	7837	20677	0.18%	0.019%
28	8.657	923	930	940	rBV	541063	866856	7.70%	0.783%
29	8.822	954	958	964	rVB4	20971	35609	0.32%	0.032%
30	9.116	997	1008	1017	rBV	1488171	2450573	21.78%	2.214%
31	9.228	1020	1027	1039	rVB	1135774	1876794	16.68%	1.696%
32	9.516	1072	1076	1083	rVB6	21241	38858	0.35%	0.035%
33	9.840	1121	1131	1142	rBV	1189211	1912279	16.99%	1.728%
34	10.381	1215	1223	1237	rBV	1837159	2989356	26.57%	2.701%
35	10.540	1242	1250	1258	rVB6	46493	113546	1.01%	0.103%
36	10.634	1258	1266	1280	rBV	347799	651620	5.79%	0.589%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

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_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

37	10.763	1280	1288	1297	rVV	2104744	3517100	31.25%	3.178%
38	10.916	1306	1314	1322	rVB8	20840	52912	0.47%	0.048%
39	11.504	1406	1414	1421	rBV	57643	101778	0.90%	0.092%
40	11.716	1439	1450	1461	rBV2	32426	110426	0.98%	0.100%
41	11.810	1461	1466	1473	rVB6	20695	46969	0.42%	0.042%
42	11.940	1480	1488	1492	rBV	209175	326915	2.91%	0.295%
43	12.004	1492	1499	1503	rVV	310516	786384	6.99%	0.710%
44	12.046	1503	1506	1520	rVB	233466	393242	3.49%	0.355%
45	12.222	1531	1536	1541	rBV	28221	42572	0.38%	0.038%
46	12.275	1541	1545	1552	rVV	27259	54228	0.48%	0.049%
47	12.346	1553	1557	1562	rVB	25294	36420	0.32%	0.033%
48	12.457	1570	1576	1579	rVB4	14976	25834	0.23%	0.023%
49	12.622	1602	1604	1611	rVB8	8705	15723	0.14%	0.014%
50	12.746	1620	1625	1629	rBV8	7944	14335	0.13%	0.013%
51	12.869	1641	1646	1652	rBV2	21379	33547	0.30%	0.030%
52	12.945	1656	1659	1663	rVV6	7597	11832	0.11%	0.011%
53	13.010	1663	1670	1676	rVB3	26301	50497	0.45%	0.046%
54	13.322	1719	1723	1731	rBV	34526	57673	0.51%	0.052%
55	13.404	1731	1737	1741	rVV5	12062	18765	0.17%	0.017%
56	13.951	1820	1830	1832	rBV2	87003	141700	1.26%	0.128%
57	13.998	1832	1838	1849	rVB	3480301	4761169	42.31%	4.302%
58	14.287	1880	1887	1897	rBV	1003558	1497975	13.31%	1.353%
59	14.475	1916	1919	1923	rVB4	13195	16272	0.14%	0.015%
60	14.592	1932	1939	1948	rBV	3056274	4407077	39.16%	3.982%
61	14.722	1957	1961	1965	rVB4	13564	19006	0.17%	0.017%
62	14.792	1965	1973	1989	rBV	337119	642210	5.71%	0.580%
63	14.998	2004	2008	2015	rVB9	16469	29199	0.26%	0.026%
64	15.222	2039	2046	2059	rVB	200948	348022	3.09%	0.314%
65	15.428	2075	2081	2087	rBV	10045	19719	0.18%	0.018%
66	15.587	2100	2108	2118	rBV	5021615	7384726	65.62%	6.672%
67	15.704	2122	2128	2137	rBV	1503678	2058445	18.29%	1.860%
68	15.822	2144	2148	2156	rVB5	21429	40027	0.36%	0.036%
69	15.945	2165	2169	2175	rVB2	33398	44547	0.40%	0.040%
70	16.275	2219	2225	2236	rBV2	433263	728043	6.47%	0.658%
71	16.363	2236	2240	2246	rVB7	8907	19021	0.17%	0.017%
72	16.998	2343	2348	2358	rBV2	296035	442094	3.93%	0.399%
73	17.092	2360	2364	2368	rBV7	14991	21228	0.19%	0.019%
74	17.145	2368	2373	2380	rVB7	12449	25730	0.23%	0.023%
75	17.222	2381	2386	2397	rVB2	66219	94452	0.84%	0.085%
76	17.345	2401	2407	2417	rBV	3845913	5359896	47.63%	4.843%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

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_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

77	17.445	2417	2424	2434	rBV	5032843	6959343	61.84%	6.288%
78	17.533	2434	2439	2443	rVV2	269042	409268	3.64%	0.370%
79	17.586	2443	2448	2465	rVB	1482836	2323301	20.65%	2.099%
80	17.828	2487	2489	2492	rBV2	13595	13490	0.12%	0.012%
81	17.998	2516	2518	2522	rBV4	10092	13827	0.12%	0.012%
82	18.157	2540	2545	2551	rBV	228977	326376	2.90%	0.295%
83	18.216	2551	2555	2560	rBV2	98513	133999	1.19%	0.121%
84	18.281	2562	2566	2571	rVV2	46737	79934	0.71%	0.072%
85	18.498	2600	2603	2607	rBV2	23897	33477	0.30%	0.030%
86	18.551	2607	2612	2618	rBV4	102909	182164	1.62%	0.165%
87	18.698	2633	2637	2642	rBV2	74118	115147	1.02%	0.104%
88	19.257	2728	2732	2737	rVB2	81472	113164	1.01%	0.102%
89	19.322	2739	2743	2747	rBV	117920	166046	1.48%	0.150%
90	19.680	2798	2804	2814	rBV	7734822	10537393	93.64%	9.520%
91	20.863	3001	3005	3010	rVB	525678	650431	5.78%	0.588%
92	21.433	3097	3102	3108	rBV	5183040	6564842	58.34%	5.931%
93	23.751	3489	3496	3511	rVB2	5620060	11252922	100.00%	10.167%
94	23.910	3516	3523	3533	rVB2	3633448	7464383	66.33%	6.744%

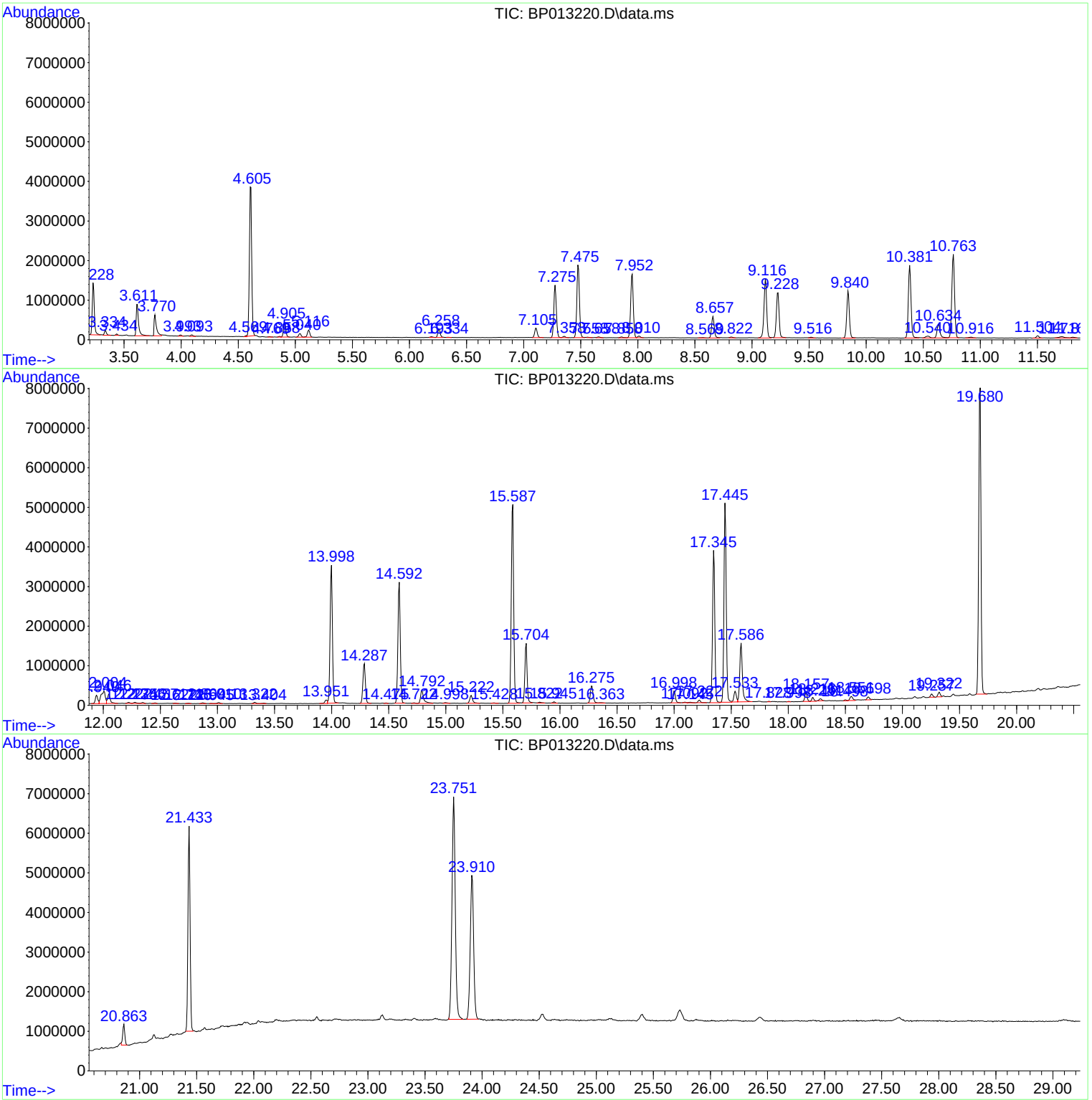
Sum of corrected areas: 110684436

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

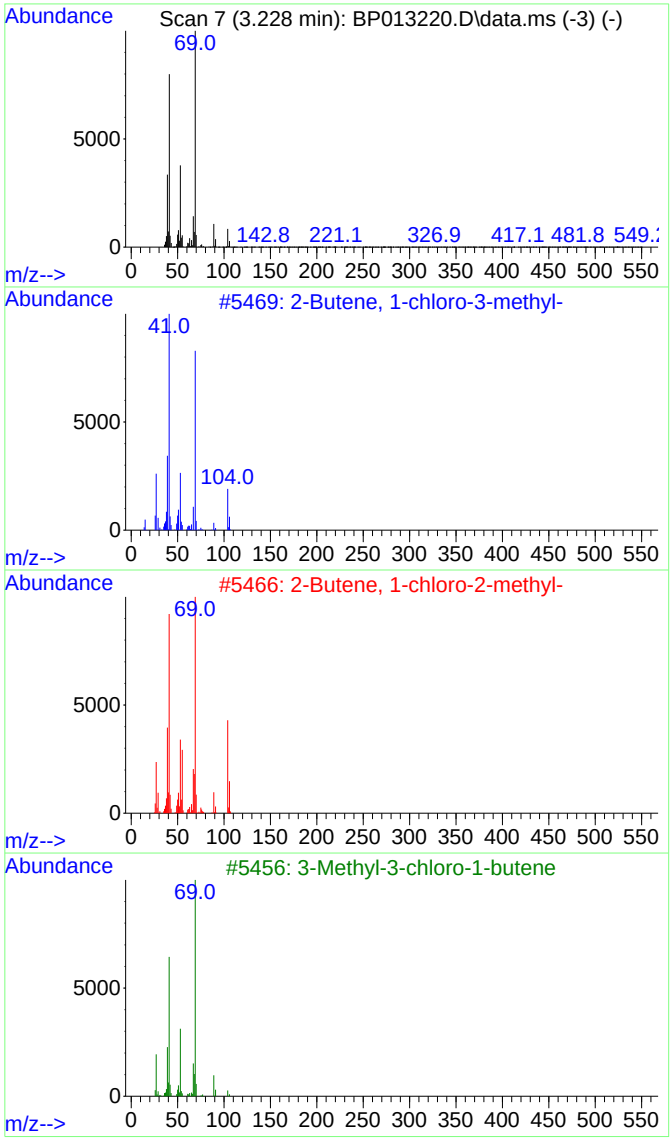
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

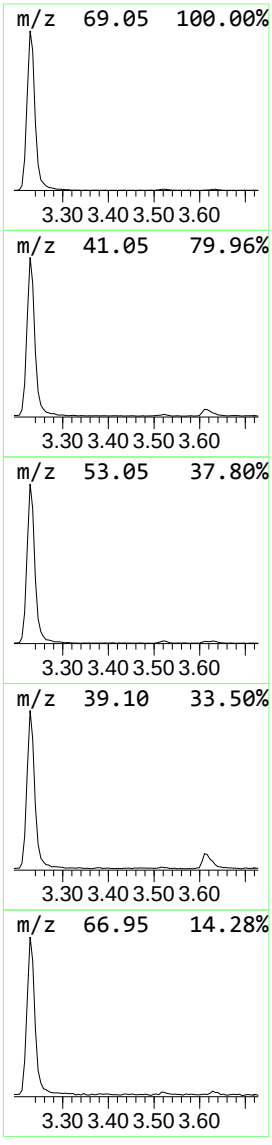
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Butene, 1-chloro-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.228	12.50 ng/ul	1607870	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 1-chloro-3-methyl-	104	C5H9Cl	000503-60-6	80
2	2-Butene, 1-chloro-2-methyl-	104	C5H9Cl	013417-43-1	72
3	3-Methyl-3-chloro-1-butene	104	C5H9Cl	002190-48-9	72
4	2-Butene, 1-chloro-3-methyl-	104	C5H9Cl	000503-60-6	64
5	2-Butene, 1-chloro-2-methyl-	104	C5H9Cl	013417-43-1	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

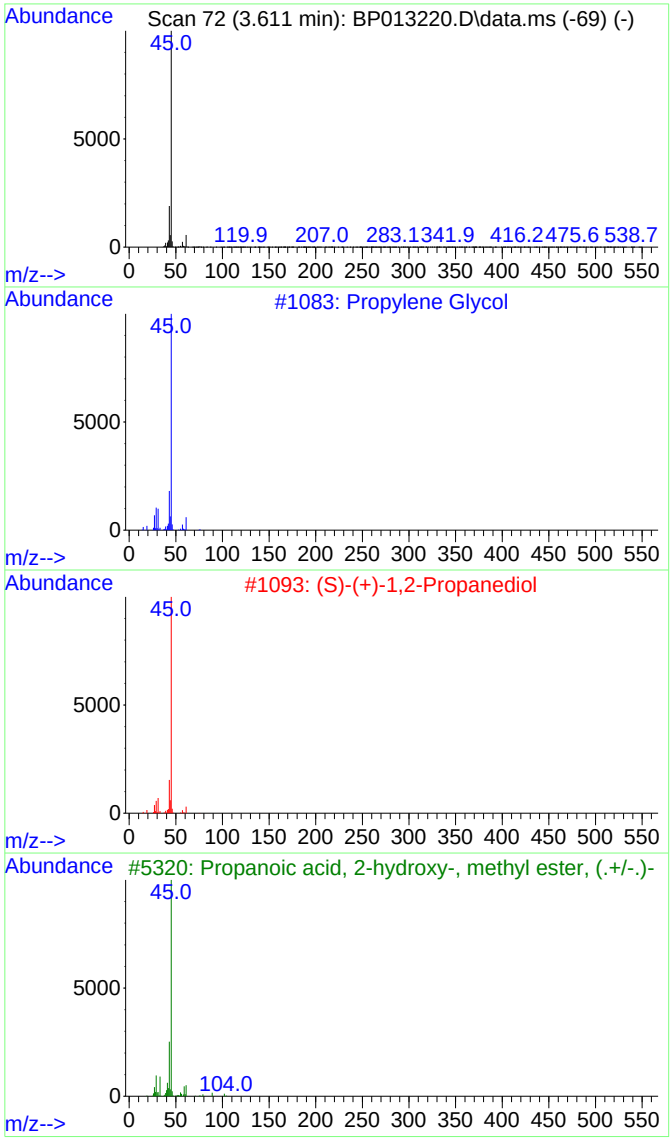
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

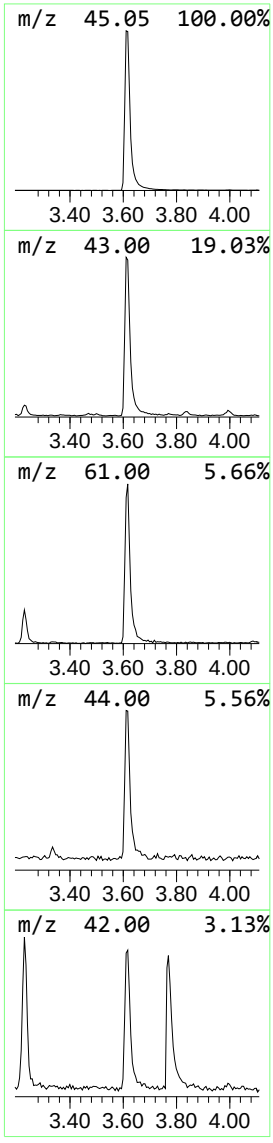
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Propylene Glycol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.611	9.02 ng/ul	1159980	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	86
2	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
3	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
4	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9
5	Propylene Glycol	76	C3H8O2	000057-55-6	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

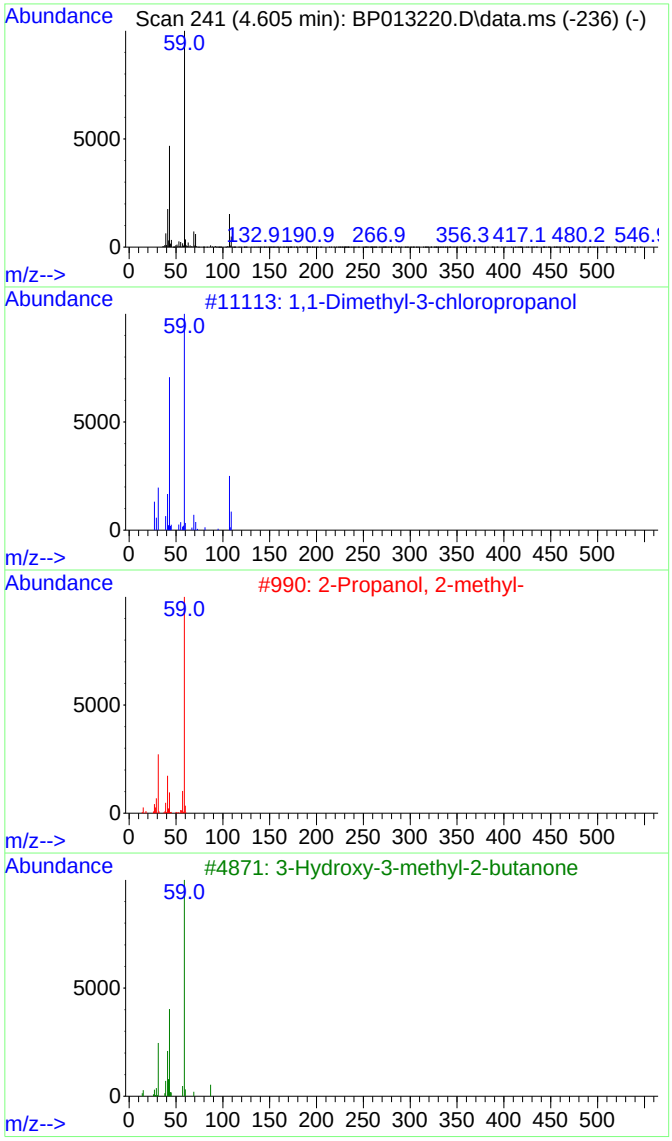
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

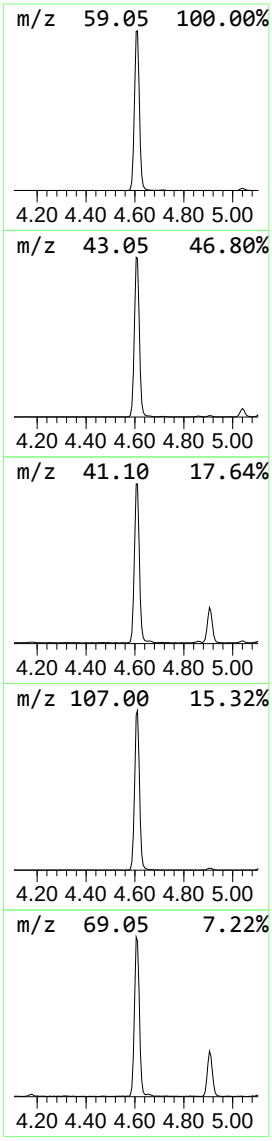
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1,1-Dimethyl-3-chloropropanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.605	41.43 ng/ul	5327410	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1-Dimethyl-3-chloropropanol	122	C5H11ClO	001985-88-2	74
2	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	50
3	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	45
4	3-Pentanol	88	C5H12O	000584-02-1	42
5	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

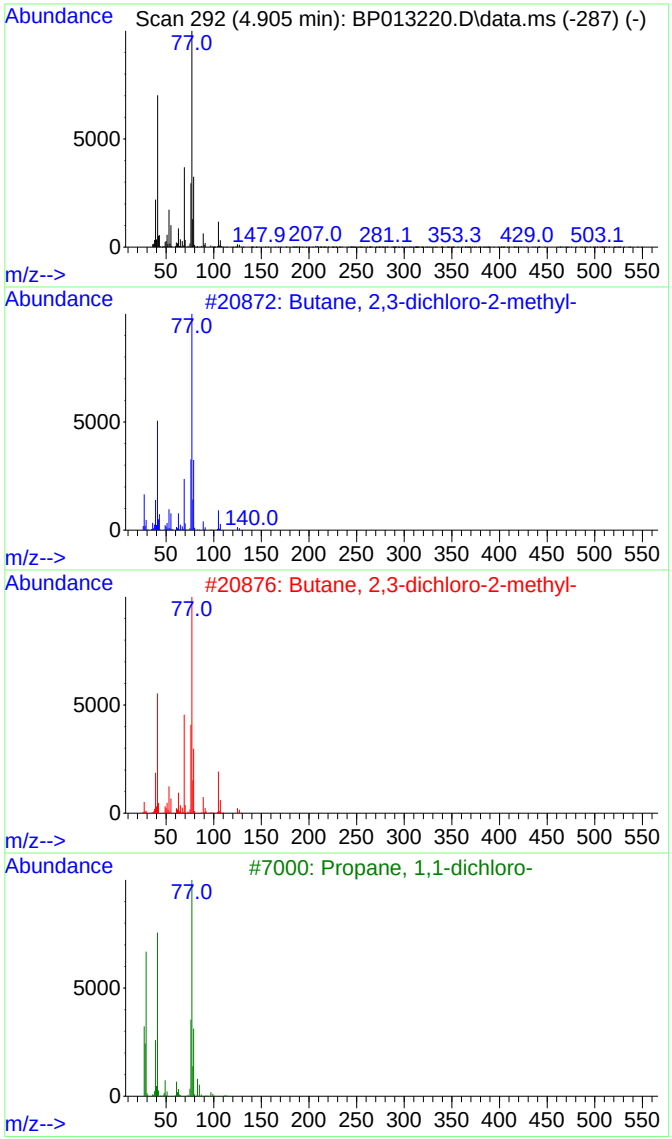
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

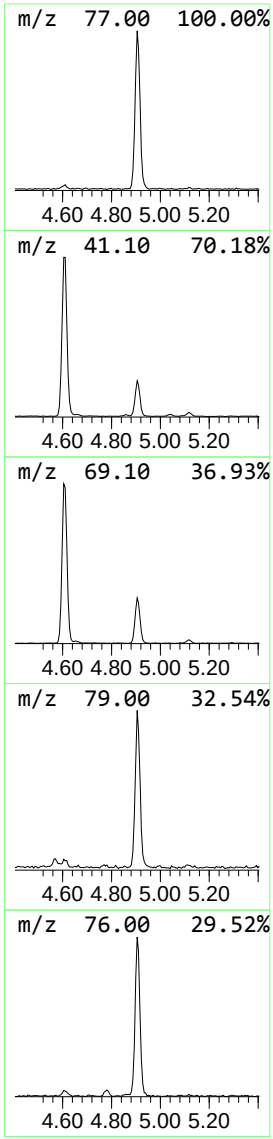
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butane, 2,3-dichloro-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.905	4.15 ng/ul	533075	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	90
2	Butane, 2,3-dichloro-2-methyl-	140	C5H10Cl2	000507-45-9	83
3	Propane, 1,1-dichloro-	112	C3H6Cl2	000078-99-9	39
4	Propane, 2-bromo-1-chloro-	156	C3H6BrCl	003017-95-6	28
5	Allyl chloride	76	C3H5Cl	000107-05-1	11





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

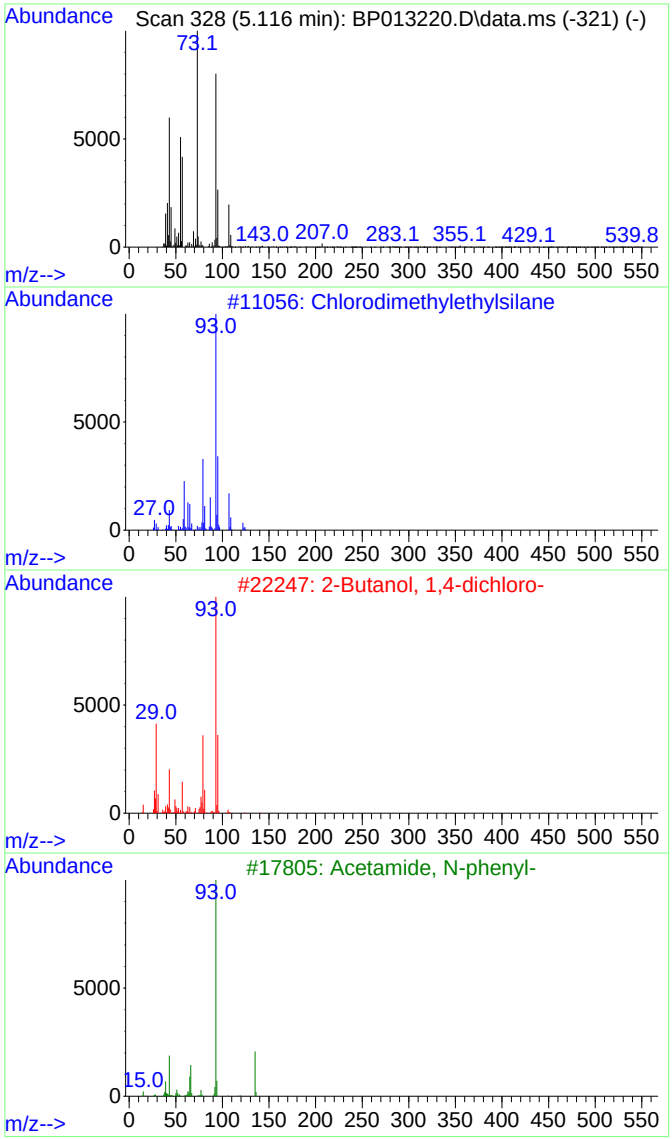
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

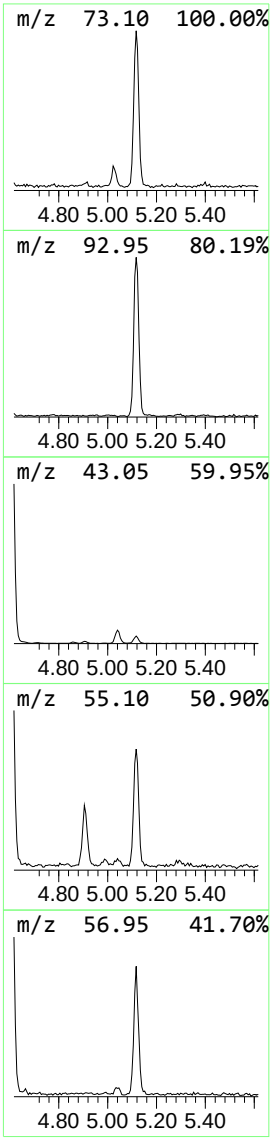
****

Peak Number 5 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.116	2.07 ng/ul	265851	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chlorodimethylethylsilane	122	C4H11ClSi	006917-76-6	36
2	2-Butanol, 1,4-dichloro-	142	C4H8Cl2O	002419-74-1	36
3	Acetamide, N-phenyl-	135	C8H9NO	000103-84-4	32
4	Acetamide, N-phenyl-	135	C8H9NO	000103-84-4	23
5	Acetamide, N-phenyl-	135	C8H9NO	000103-84-4	23





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

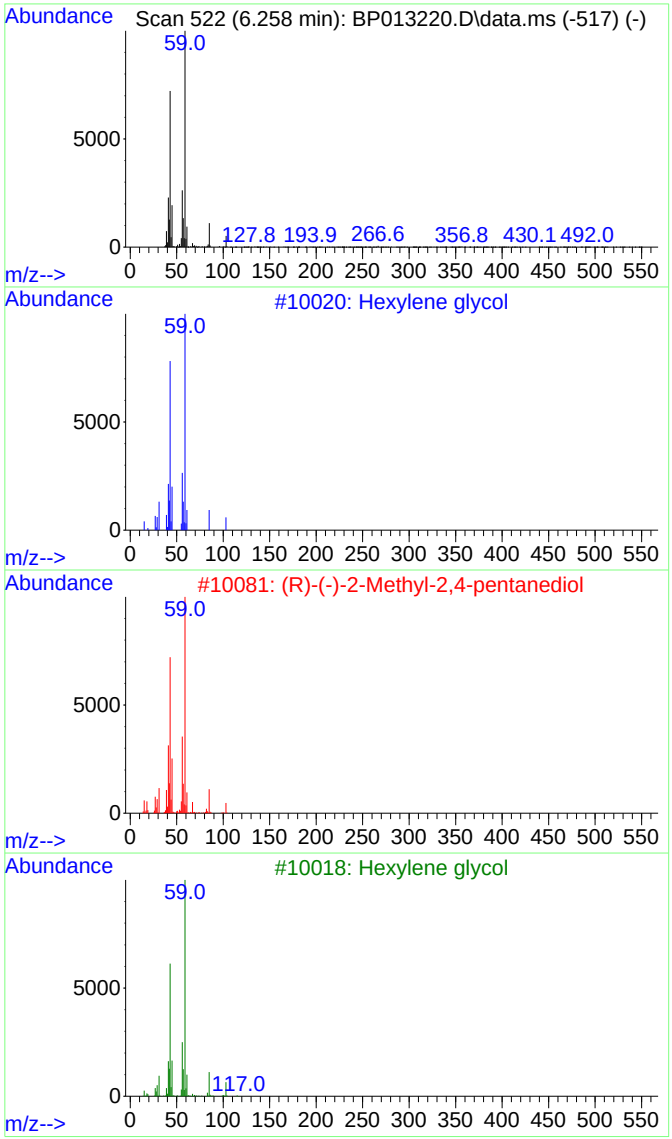
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

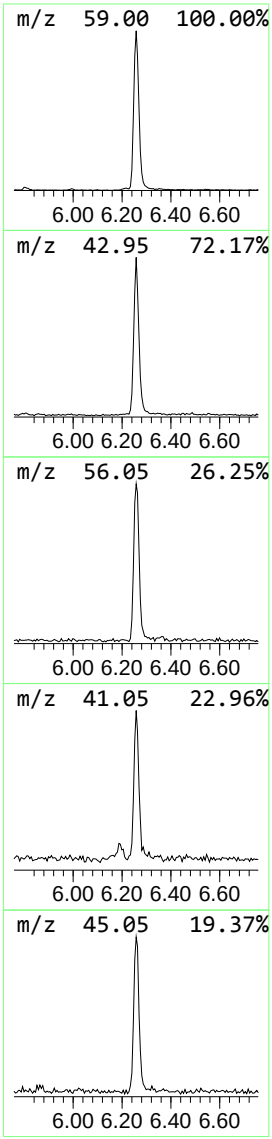
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Hexylene glycol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.258	2.43 ng/ul	313060	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexylene glycol	118	C6H14O2	000107-41-5	90
2	(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	90
3	Hexylene glycol	118	C6H14O2	000107-41-5	90
4	Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	72
5	Hexylene glycol	118	C6H14O2	000107-41-5	53





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Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

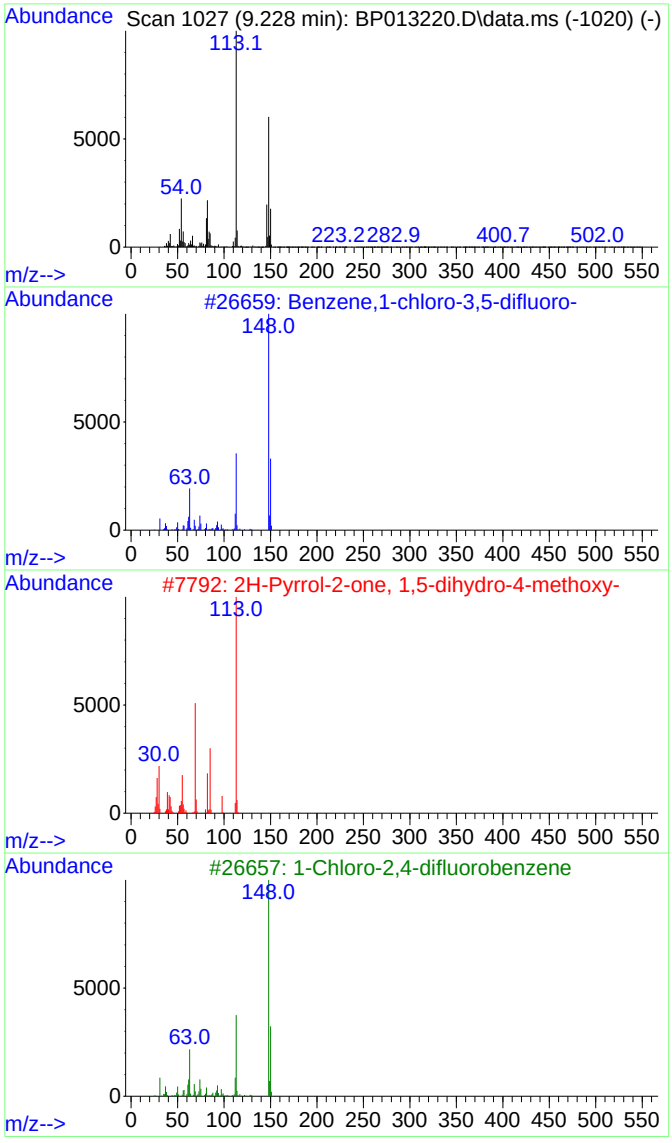
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

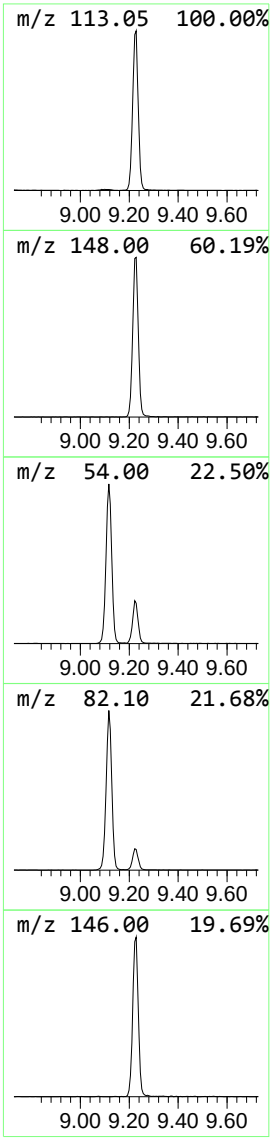
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene,1-chloro-3,5-difluoro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.228	14.59 ng/ul	1876790	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,1-chloro-3,5-difluoro-	148	C6H3ClF2	001435-43-4	52
2	2H-Pyrrol-2-one, 1,5-dihydro-4-m...	113	C5H7NO2	069778-83-2	43
3	1-Chloro-2,4-difluorobenzene	148	C6H3ClF2	001435-44-5	40
4	3-Fluoro-2-hydroxypyridine	113	C5H4FNO	001547-29-1	30
5	3-Fluoro-5-hydroxypyridine	113	C5H4FNO	209328-55-2	27





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Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

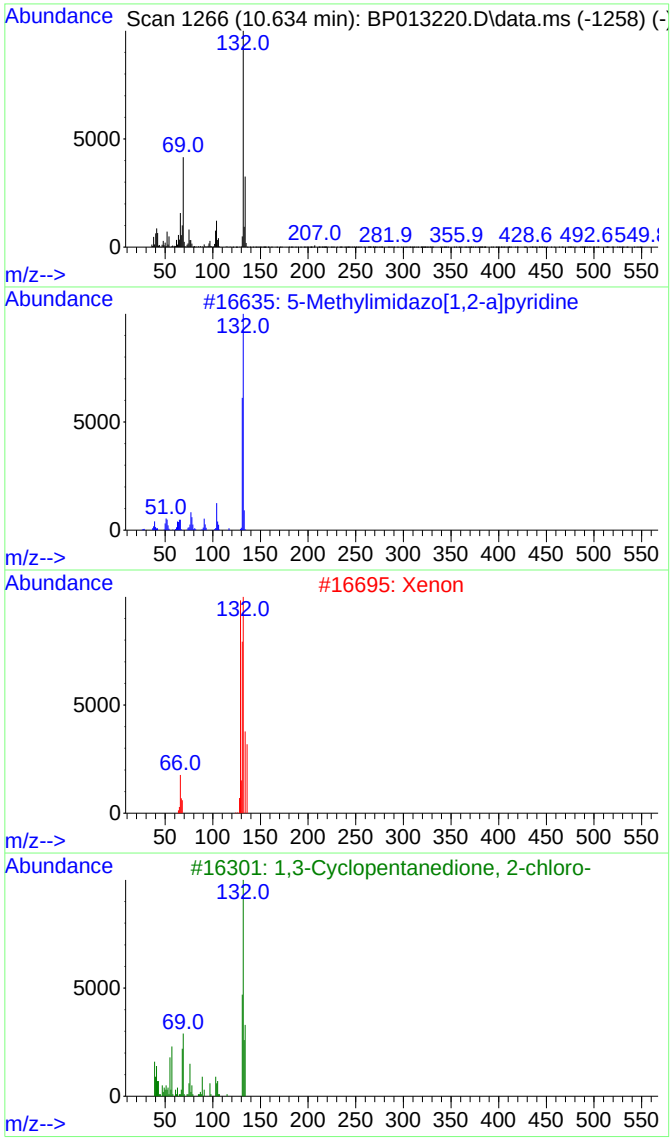
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

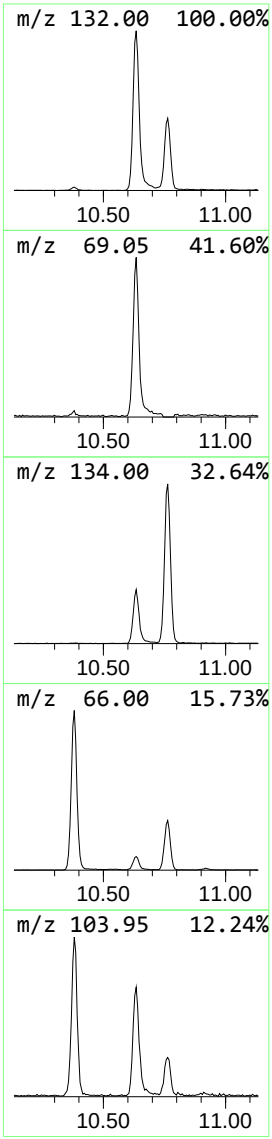
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 5-Methylimidazo[1,2-a]pyridine Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.634	3.71 ng/ul	651620	Naphthalene-d8	10.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methylimidazo[1,2-a]pyridine	132	C8H8N2	000933-69-7	58
2	Xenon	132	Xe	007440-63-3	53
3	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	53
4	7-Methylimidazo[1,2-a]pyridine	132	C8H8N2	000874-39-5	50
5	1H-Indol-3-amine	132	C8H8N2	007250-19-3	50





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Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

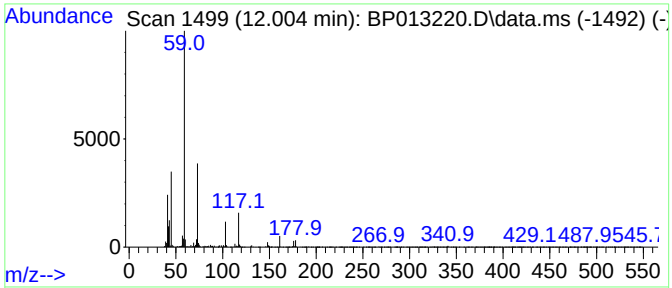
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

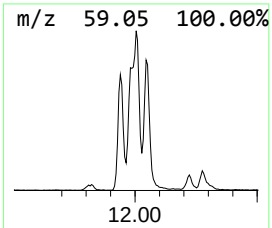
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

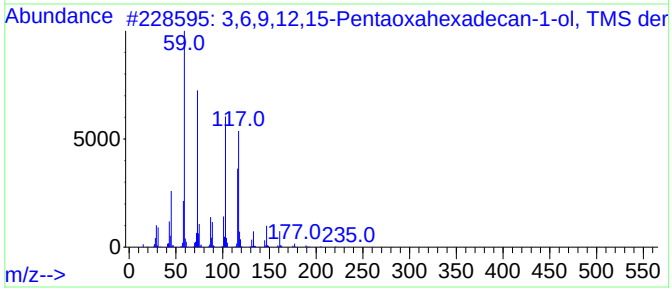
****
Peak Number 9 3,6,9,12,15-Pentaoxahexadec... Concentration Rank 6

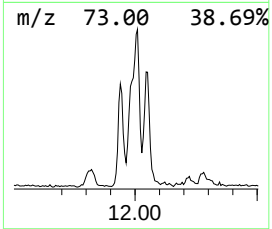
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.004	4.47 ng/ul	786384	Naphthalene-d8	10.763

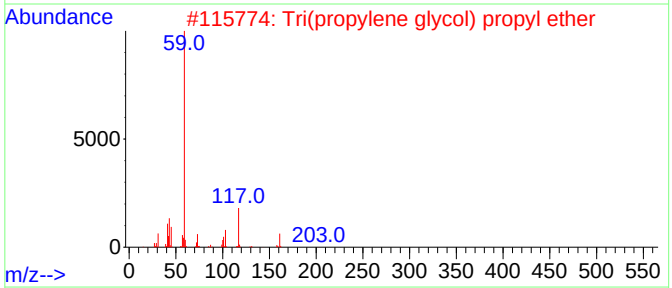
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6,9,12,15-Pentaoxahexadecan-1-...	324	C14H32O6Si	1000352-07-5	78
2	Tri(propylene glycol) propyl ether	234	C12H26O4	096077-04-2	64
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
4	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	64
5	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	64

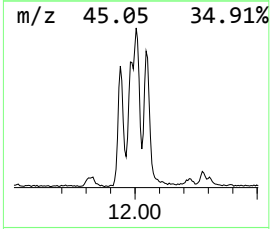


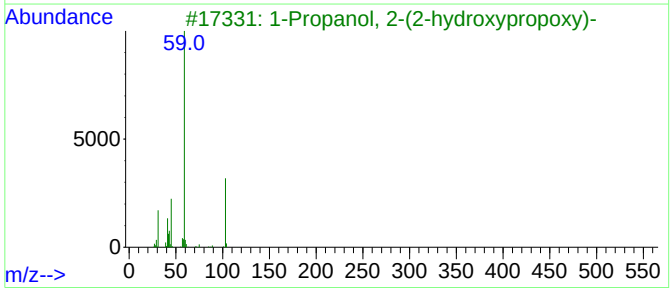


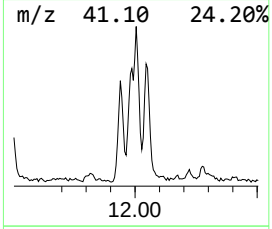


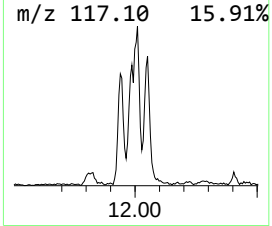












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

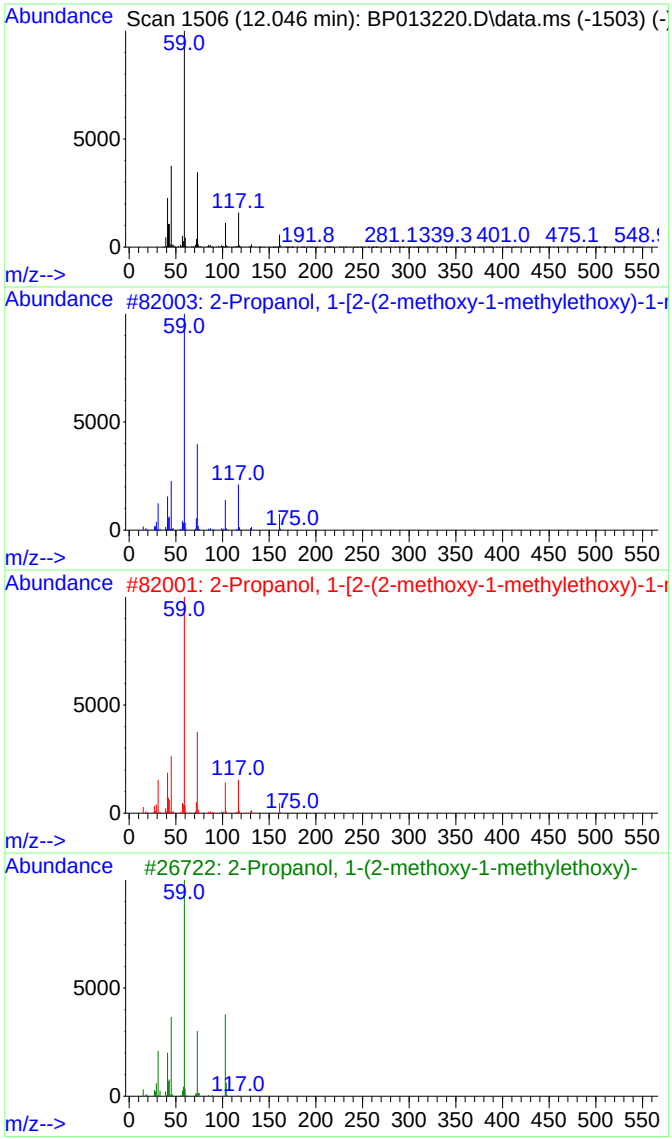
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

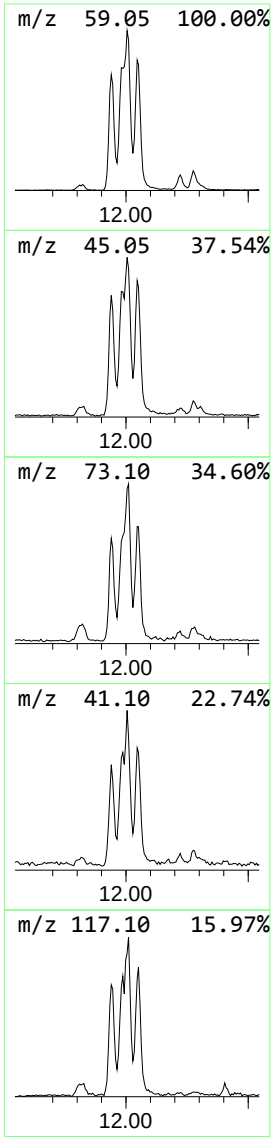
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.046	2.24 ng/ul	393242	Naphthalene-d8	10.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	78
2	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	78
3	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
4	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	56
5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

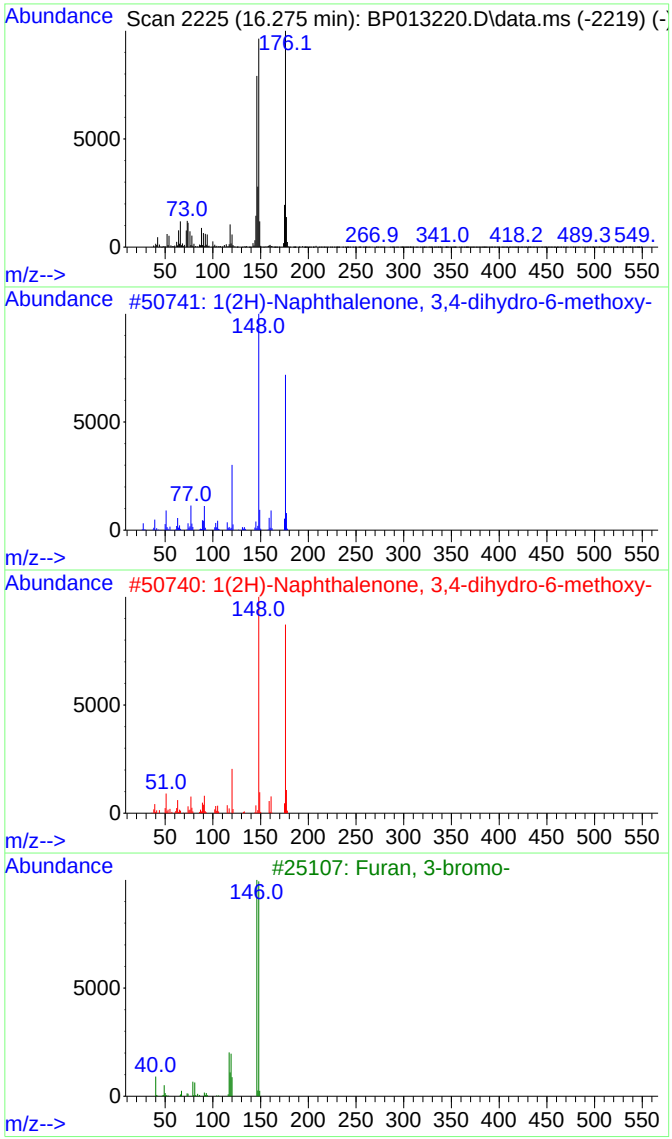
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

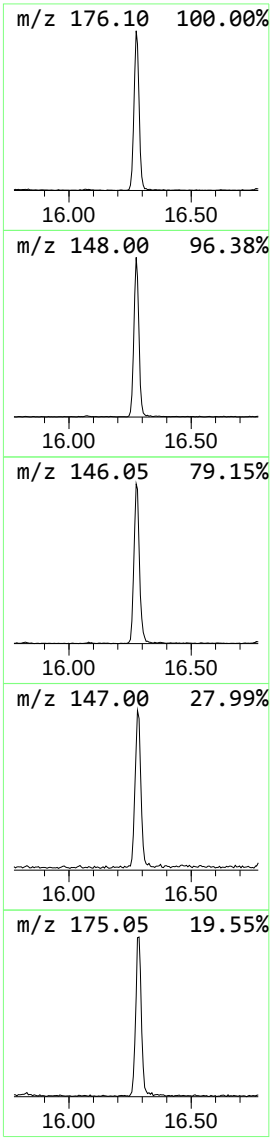
****

Peak Number 11 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	2.72 ng/ul	728043	Phenanthrene-d10	17.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	43
2	1(2H)-Naphthalenone, 3,4-dihydro...	176	C11H12O2	001078-19-9	43
3	Furan, 3-bromo-	146	C4H3BrO	022037-28-1	38
4	6-Fluoro-2-methylquinolin-4-amine	176	C10H9FN2	1000409-97-0	35
5	Pyridine-3-carbonitrile, 1-ethyl...	176	C10H12N2O	1010316-81-3	30





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Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B0AA8

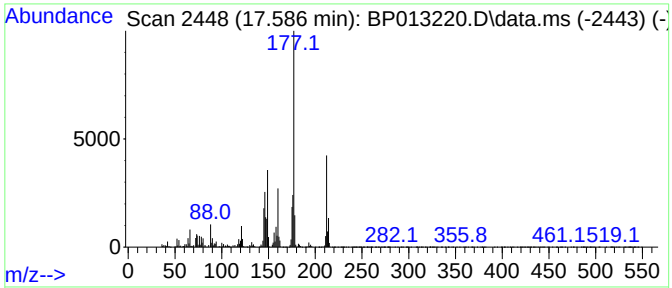
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

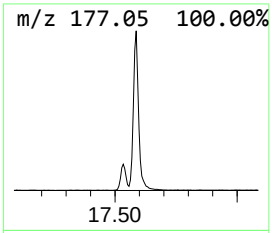
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

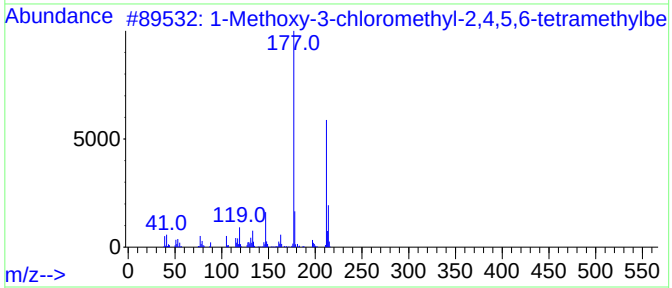
****
Peak Number 12 1-Methoxy-3-chloromethyl-2,... Concentration Rank 5

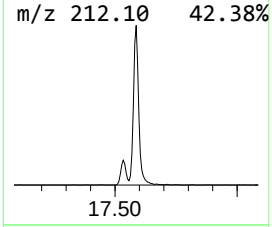
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.586	8.67 ng/ul	2323300	Phenanthrene-d10	17.345

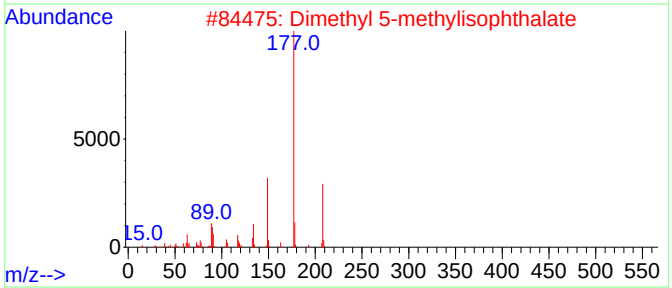
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methoxy-3-chloromethyl-2,4,5,6...	212	C12H17ClO	073942-80-0	53
2	Dimethyl 5-methylisophthalate	208	C11H12O4	017649-58-0	35
3	Terephthalic acid, but-3-enyl et...	248	C14H16O4	1000356-33-3	35
4	Spiro[cyclopenta[d]-1,3,2-dioxab...	206	C12H19BO2	062238-26-0	35
5	trans-.beta.-Ionone	192	C13H20O	000079-77-6	27

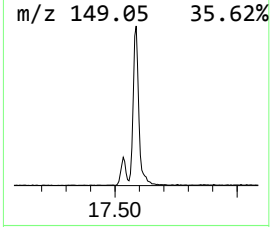


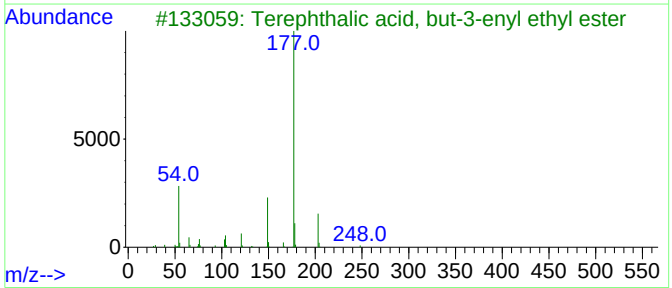


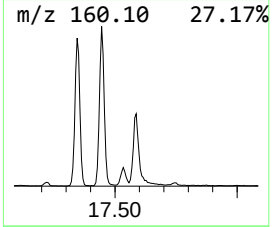


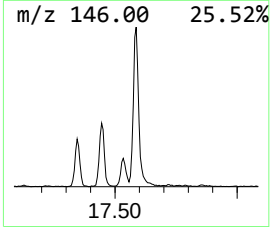












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013220.D
Acq On : 21 Dec 2022 22:02
Operator : CG/JU
Sample : N6050-01
Misc :
ALS Vial : 102 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Butene, 1-chl...	3.228	12.5	ng/ul	1607870	1	7.952	2571940	20.0
Propylene Glycol	3.611	9.0	ng/ul	1159980	1	7.952	2571940	20.0
1,1-Dimethyl-3-...	4.605	41.4	ng/ul	5327410	1	7.952	2571940	20.0
Butane, 2,3-dic...	4.905	4.2	ng/ul	533075	1	7.952	2571940	20.0
unknown-01	5.116	2.1	ng/ul	265851	1	7.952	2571940	20.0
Hexylene glycol	6.258	2.4	ng/ul	313060	1	7.952	2571940	20.0
Benzene,1-chlor...	9.228	14.6	ng/ul	1876790	1	7.952	2571940	20.0
5-Methylimidazo...	10.634	3.7	ng/ul	651620	2	10.763	3517100	20.0
3,6,9,12,15-Pen...	12.004	4.5	ng/ul	786384	2	10.763	3517100	20.0
2-Propanol, 1-[...	12.046	2.2	ng/ul	393242	2	10.763	3517100	20.0
unknown-02	16.275	2.7	ng/ul	728043	4	17.345	5359900	20.0
1-Methoxy-3-chl...	17.586	8.7	ng/ul	2323300	4	17.345	5359900	20.0