LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

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-BP040722.M
Title : SVOA CALIBRATION

Signal : TIC: BP009839.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	17	rVB	173821	247616	6.95%	0.620%
2	3.381	46	50	58	rBV	19912	29967	0.84%	0.075%
3	3.799	118	121	131	rBV	280919	464857	13.04%	1.163%
4	4.140	174	179	186	rBV4	19910	32686	0.92%	0.082%
5	4.358	211	216	224	rVB5	13733	26217	0.74%	0.066%
6	4.781	284	288	293	rBV2	9229	13921	0.39%	0.035%
7	4.964	315	319	322	rBV6	5565	8488	0.24%	0.021%
8	5.269	364	371	381	rVB	435806	670921	18.82%	1.679%
9	5.469	399	405	409	rVV8	4163	9197	0.26%	0.023%
10	6.052	498	504	507	rBV8	4812	8841	0.25%	0.022%
11	6.446	566	571	576	rBV6	5847	10474	0.29%	0.026%
12	6.634	596	603	612	rVB4	18124	36609	1.03%	0.092%
13	6.946	650	656	659	rVV6	8346	18034	0.51%	0.045%
14	6.999	659	665	670	rVV2	14310	38228	1.07%	0.096%
15	7.052	670	674	678	rVV5	7277	14072	0.39%	0.035%
16	7.116	678	685	695	rVV	437390	731881	20.53%	1.831%
17	7.287	707	714	723	rVV	325842	536699	15.06%	1.343%
18	7.357	723	726	734	rVB5	7981	15837	0.44%	0.040%
19	7.487	740	748	761	rBV	421472	713282	20.01%	1.785%
20	7.957	821	828	838	rBV	358910	620599	17.41%	1.553%
21	8.034	838	841	848	rVB7	9327	19117	0.54%	0.048%
22	8.181	861	866	872	rBV10	6150	11394	0.32%	0.029%
23	8.257	872	879	885	rVV2	39829	67245	1.89%	0.168%
24	8.581	928	934	938	rBV5	8722	15356	0.43%	0.038%
25	8.663	941	948	955	rVV	584111	963065	27.02%	2.410%
26	8.728	956	959	964	rVV2	17941	29130	0.82%	0.073%
27	8.963	995	999	1009	rVB2	5177	14995	0.42%	0.038%
28	9.116	1012	1025	1033	rBV	448678	772734	21.68%	1.934%
29	9.204	1033	1040	1048	rVV	132055	291152	8.17%	0.729%
30	9.287	1048	1054	1059	rVV3	31205	75930	2.13%	0.190%
31	9.340	1059	1063	1068	rVB4	16565	27417	0.77%	0.069%
32	9.457	1077	1083	1091	rVB	132955	220637	6.19%	0.552%
33	9.575	1098	1103	1112	rVB3	9323	20227	0.57%	0.051%
34	9.663	1112	1118	1121	rBV3	8765	13879	0.39%	0.035%
35	9.775	1132	1137	1142	rBV	41860	75384	2.11%	0.189%
36	9.846	1142	1149	1160	rVB	362804	625517	17.55%	1.565%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

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-BP040722.M
Title : SVOA CALIBRATION

37	10.181	1201	1206	1213	rVB5	9204	16856	0.47%	0.042%
38	10.257	1213	1219	1226	rBV9	8453	17707	0.50%	0.044%
39	10.381	1226	1240	1254	rBV	578190	1014495	28.46%	2.538%
40	10.640	1274	1284	1291	rVV3	20127	44483	1.25%	0.111%
41	10.769	1297	1306	1318	rBV	570011	990808	27.80%	2.479%
42	10.899	1318	1328	1338	rVV	548040	1015797	28.50%	2.542%
43	10.993	1338	1344	1350	rVB7	20786	36418	1.02%	0.091%
44	11.169	1369	1374	1378	rBV5	10001	15422	0.43%	0.039%
45	11.216	1378	1382	1388	rVB8	7498	14106	0.40%	0.035%
46	11.310	1392	1398	1409	rBV2	99983	174891	4.91%	0.438%
47	11.716	1457	1467	1469	rBV7	9010	22996	0.65%	0.058%
48	11.881	1489	1495	1499	rBV7	7639	13997	0.39%	0.035%
49	12.093	1522	1531	1538	rBV5	8063	19974	0.56%	0.050%
50	12.169	1538	1544	1546	rVV6	8499	19622	0.55%	0.049%
51	12.204	1546	1550	1554	rVV5	12718	19730	0.55%	0.049%
52	12.316	1564	1569	1572	rVV8	11937	20172	0.57%	0.050%
53	12.351	1572	1575	1580	rVB2	14724	22903	0.64%	0.057%
54	12.469	1589	1595	1604	rVB	145404	234817	6.59%	0.588%
55	12.640	1613	1624	1637	rVB	80355	150140	4.21%	0.376%
56	12.816	1646	1654	1659	rBV	4366	13184	0.37%	0.033%
57	12.869	1660	1663	1668	rVV6	6500	11732	0.33%	0.029%
58	12.922	1668	1672	1681	rVV4	5242	15572	0.44%	0.039%
59	13.010	1681	1687	1698	rVV4	16669	41609	1.17%	0.104%
60	13.192	1714	1718	1724	rVB6	6780	12576	0.35%	0.031%
61	13.322	1731	1740	1748	rVB3	38782	100755	2.83%	0.252%
62	13.428	1748	1758	1764	rBV2	23565	44924	1.26%	0.112%
63	13.522	1768	1774	1779	rVV	37257	59890	1.68%	0.150%
64	13.645	1789	1795	1802	rBV	108133	158442	4.44%	0.396%
65	13.910	1836	1840	1844	rVV7	6518	12786	0.36%	0.032%
66	13.998	1848	1855	1868	rVV	1153441	1672862	46.93%	4.186%
67	14.134	1874	1878	1886	rVB	4682	8957	0.25%	0.022%
68	14.281	1897	1903	1915	rVV	1224277	1819069	51.03%	4.552%
69	14.504	1938	1941	1949	rVB7	4769	10834	0.30%	0.027%
70	14.592	1949	1956	1967	rBV	938146	1452077	40.74%	3.633%
71	14.769	1980	1986	2003	rBV	549650	867998	24.35%	2.172%
72	14.898	2005	2008	2016	rVB8	6810	12148	0.34%	0.030%
73	14.998	2021	2025	2032	rVB10	9470	16716	0.47%	0.042%
74	15.245	2064	2067	2073	rVB5	11894	19818	0.56%	0.050%
75	15.357	2077	2086	2091	rBV	425570	849368	23.83%	2.125%
76	15.581	2117	2124	2133	rVB	1630889	2410428	67.62%	6.031%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

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-BP040722.M
Title : SVOA CALIBRATION

77	15.692	2133	2143	2154	rBV	482991	706827	19.83%	1.769%
78	15.822	2160	2165	2169	rBV2	18161	28909	0.81%	0.072%
79	15.898	2174	2178	2183	rBV5	7125	10459	0.29%	0.026%
80	16.051	2200	2204	2212	rVB3	9262	15550	0.44%	0.039%
81	16.157	2219	2222	2227	rVB4	5692	8752	0.25%	0.022%
82	16.434	2262	2269	2275	rBV4	10606	19290	0.54%	0.048%
83	16.557	2285	2290	2294	rBV5	17072	25331	0.71%	0.063%
84	17.251	2404	2408	2414	rVB	9154	17205	0.48%	0.043%
85	17.339	2417	2423	2430	rVV	1292098	1853814	52.01%	4.639%
86	17.439	2432	2440	2451	rVV	1930569	2779554	77.98%	6.955%
87	17.851	2506	2510	2514	rBV7	8917	12694	0.36%	0.032%
88	18.016	2534	2538	2543	rBV8	7341	9487	0.27%	0.024%
89	18.222	2569	2573	2580	rBV	121601	169255	4.75%	0.424%
90	18.763	2663	2665	2669	rVB5	8461	9221	0.26%	0.023%
91	19.245	2743	2747	2754	rVB2	28999	42142	1.18%	0.105%
92	19.310	2754	2758	2762	rBV	32293	47756	1.34%	0.119%
93	19.457	2779	2783	2788	rBV6	23303	32984	0.93%	0.083%
94	19.669	2813	2819	2822	rBV	2463056	3381221	94.86%	8.460%
95	19.704	2822	2825	2833	rVB	866463	998998	28.03%	2.500%
96	20.163	2901	2903	2907	rBV3	17548	21246	0.60%	0.053%
97	21.127	3063	3067	3074	rBV	429814	520085	14.59%	1.301%
98	21.421	3111	3117	3122	rBV	1748463	2335796	65.53%	5.845%
99	23.727	3501	3509	3516	rBV	1775830	3564517	100.00%	8.919%
100	23.886	3528	3536	3547	rVB2	1168216	2385167	66.91%	5.968%

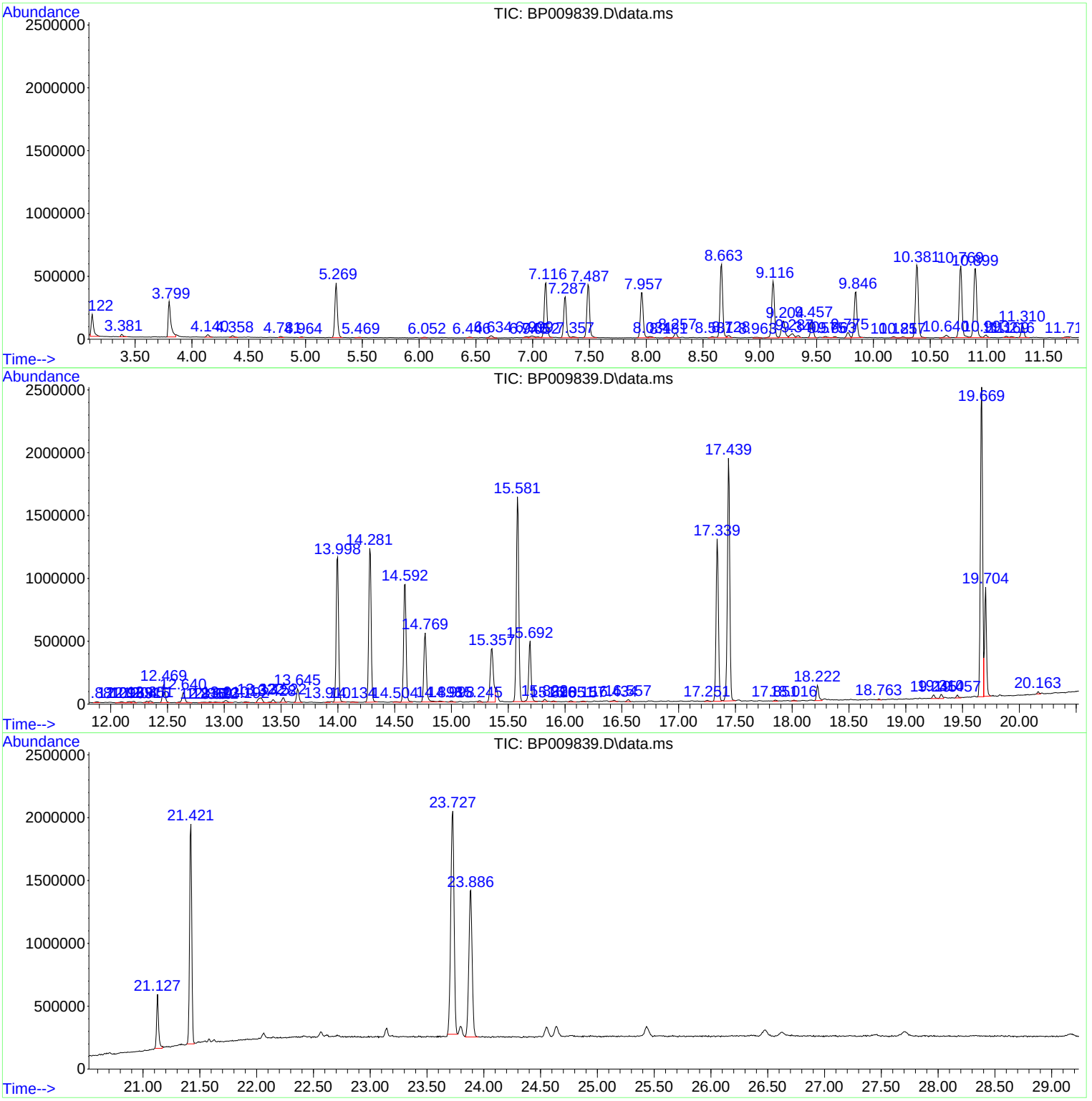
Sum of corrected areas: 39964944

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

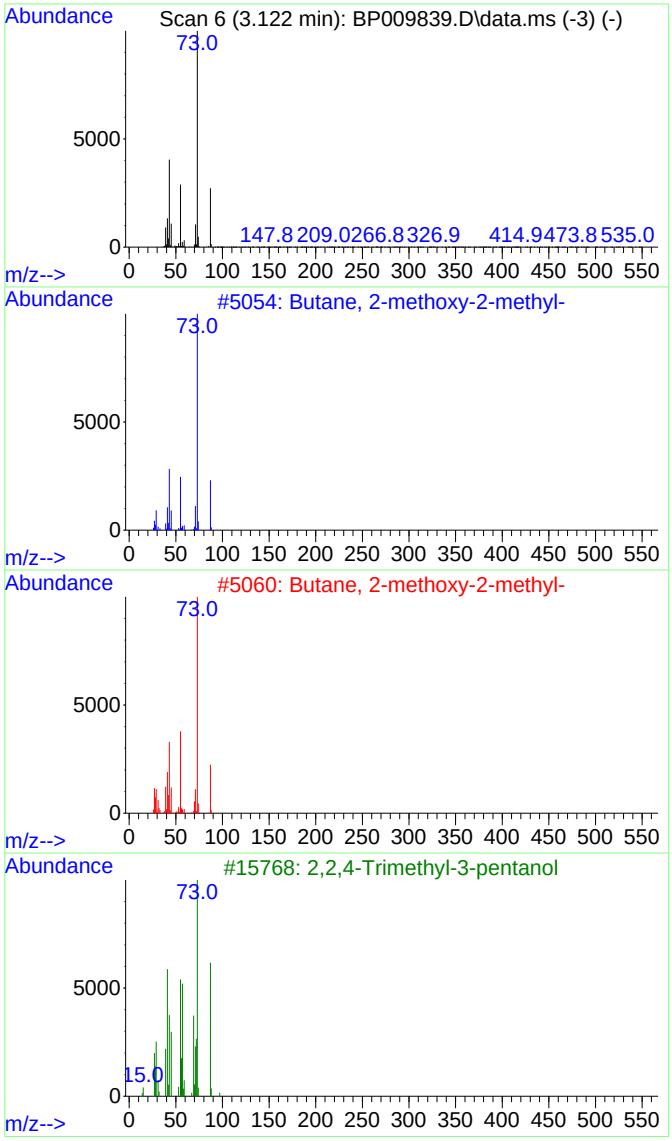
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

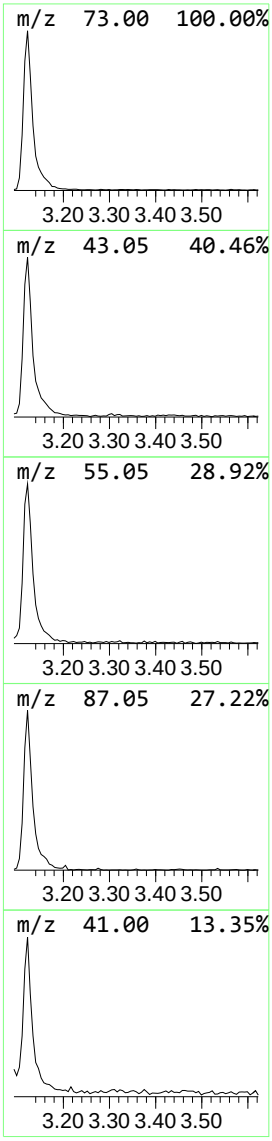
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	7.98 ng/ul	247616	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
4	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

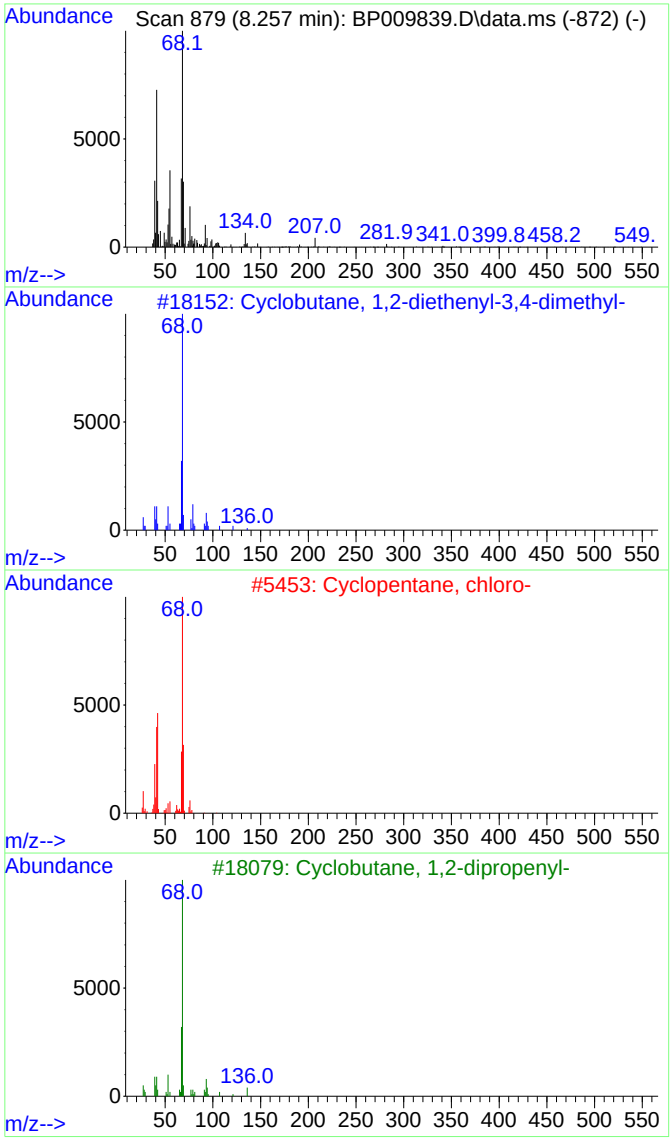
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

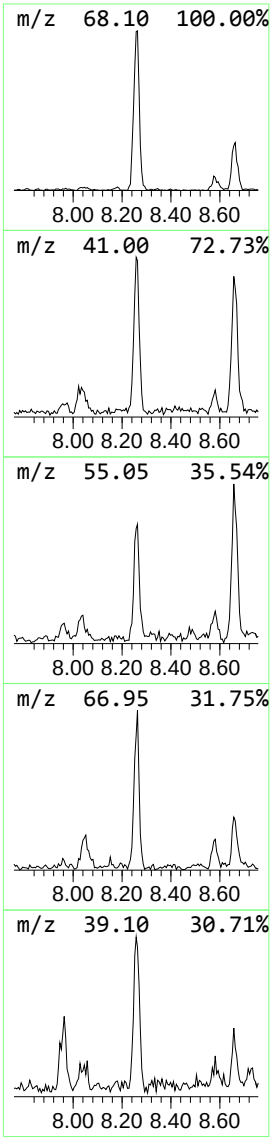
****

Peak Number 3 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.257	2.17 ng/ul	67245	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, 1,2-diethenyl-3,4-d...	136	C10H16	1000034-00-1	47
2	Cyclopentane, chloro-	104	C5H9Cl	000930-28-9	40
3	Cyclobutane, 1,2-dipropenyl-	136	C10H16	022769-00-2	38
4	Cyclopentane, (1-methylethyl)-	112	C8H16	003875-51-2	33
5	Bicyclo[4.1.0]heptane, 7,7-dichl...	164	C7H10Cl2	000823-69-8	33





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

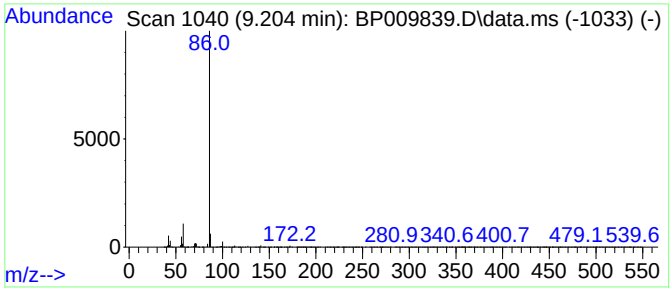
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

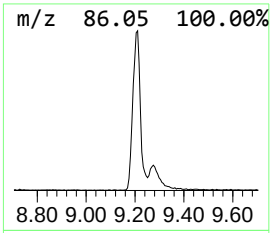
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

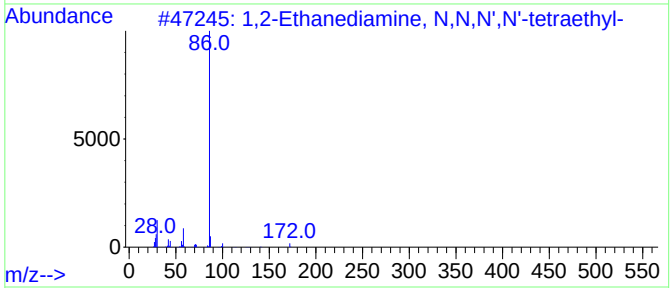
****
Peak Number 4 1,2-Ethanediamine, N,N,N',N... Concentration Rank 3

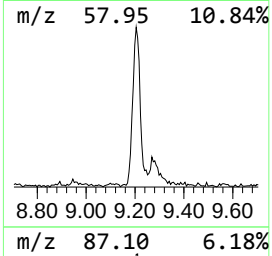
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.204	9.38 ng/ul	291152	1,4-Dichlorobenzene-d4	7.963

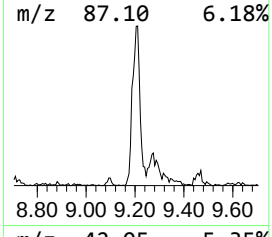
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Ethanediamine, N,N,N',N'-tet...	172	C10H24N2	000150-77-6	78
2	2-Propenoic acid, 2-methyl-, 2-(...	185	C10H19NO2	000105-16-8	78
3	Clofexamide	284	C14H21ClN2O2	001223-36-5	78
4	Ethyl N,N-diethylaminoacetate	159	C8H17NO2	002644-21-5	78
5	3-Fluorobenzoic acid, 2-(diethyl...	239	C13H18FNO2	002994-22-1	78

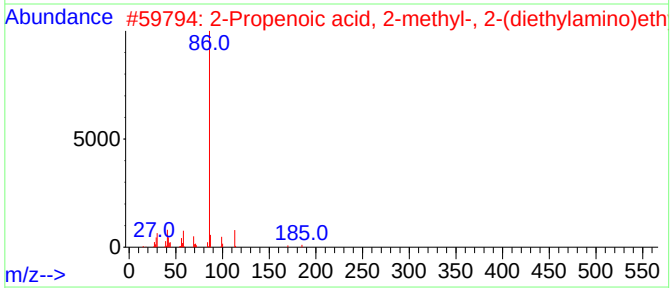


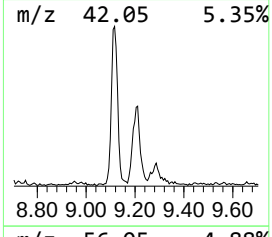


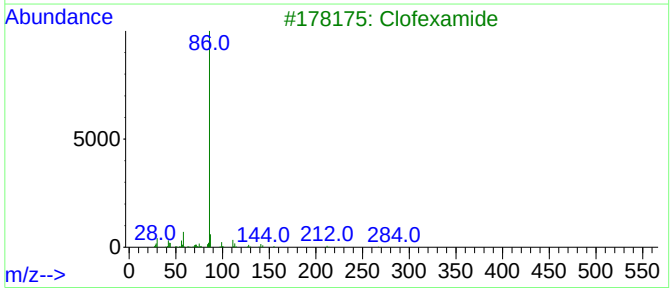


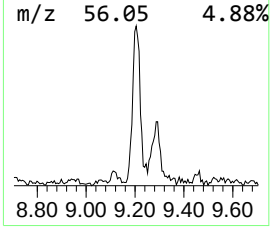












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

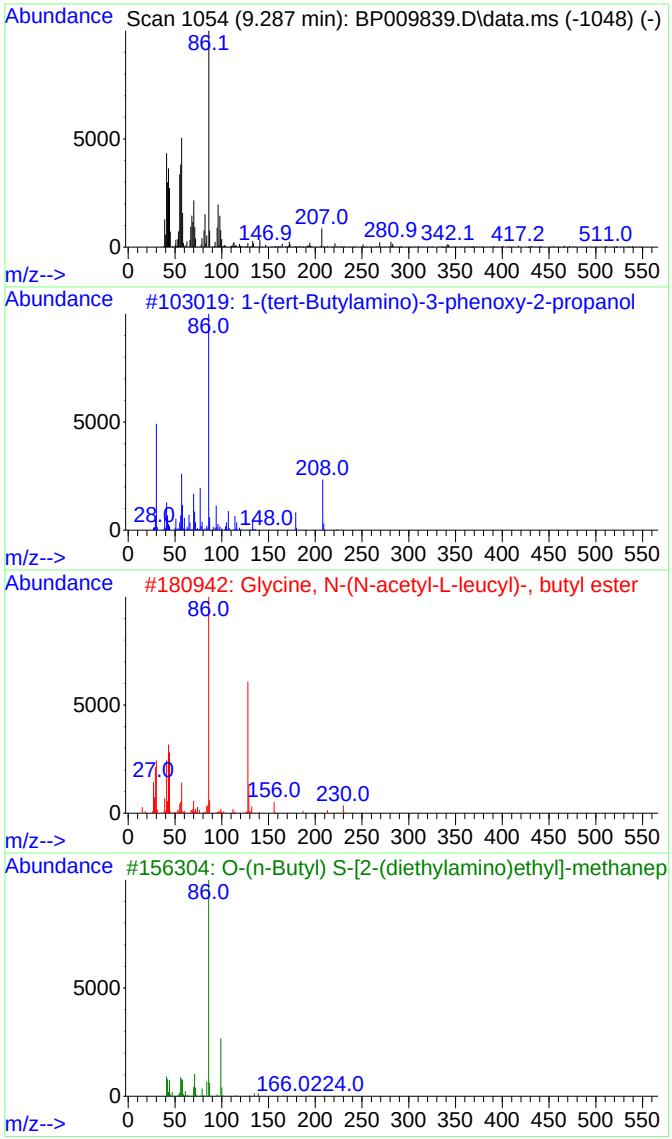
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

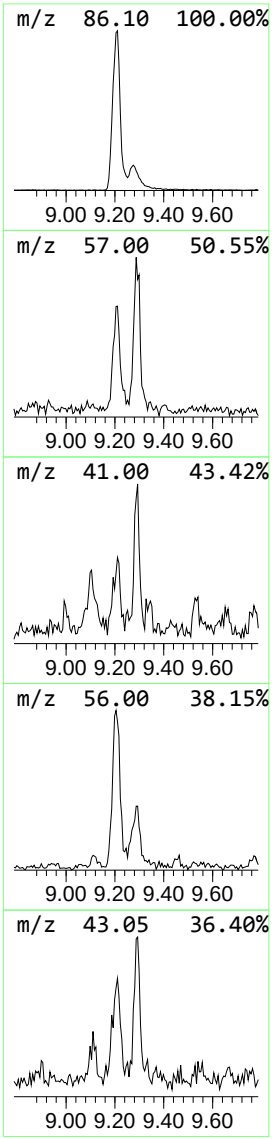
****

Peak Number 5 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.287	2.45 ng/ul	75930	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(tert-Butylamino)-3-phenoxy-2-...	223	C13H21NO2	1000468-85-1	35
2	Glycine, N-(N-acetyl-L-leucyl)-,...	286	C14H26N2O4	055712-44-2	35
3	O-(n-Butyl) S-[2-(diethylamino)e...	267	C11H26NO2PS	468712-10-9	35
4	L-Norleucine	131	C6H13NO2	000327-57-1	35
5	Benzoic acid, 4-amino-2-chloro-,...	270	C13H19ClN2O2	000133-16-4	35





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Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

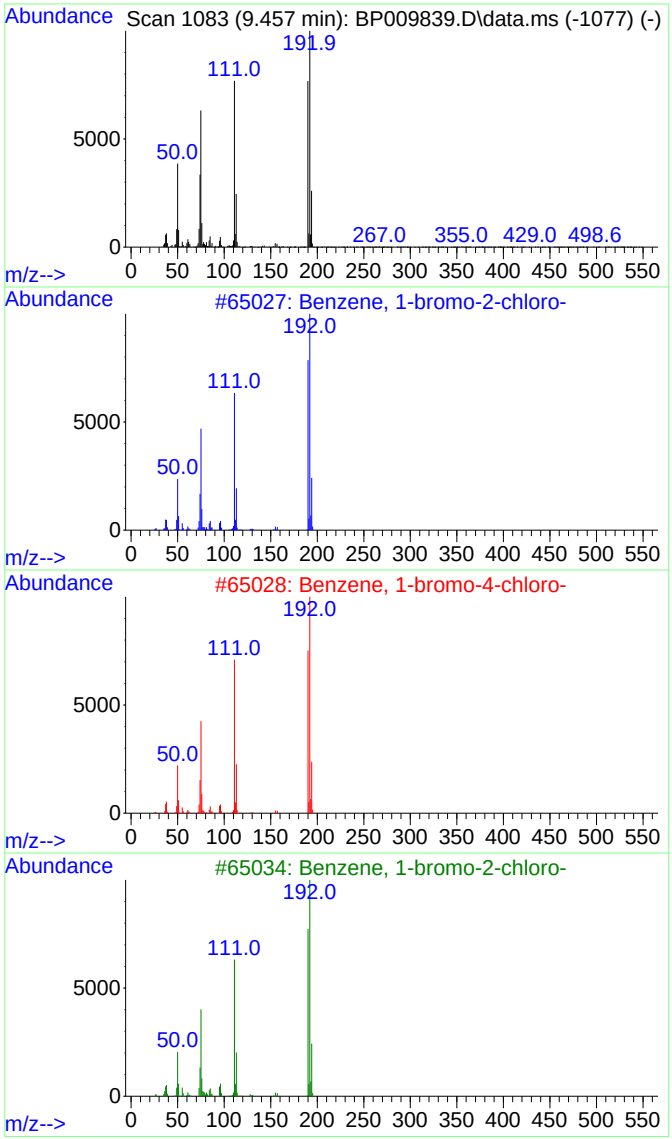
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

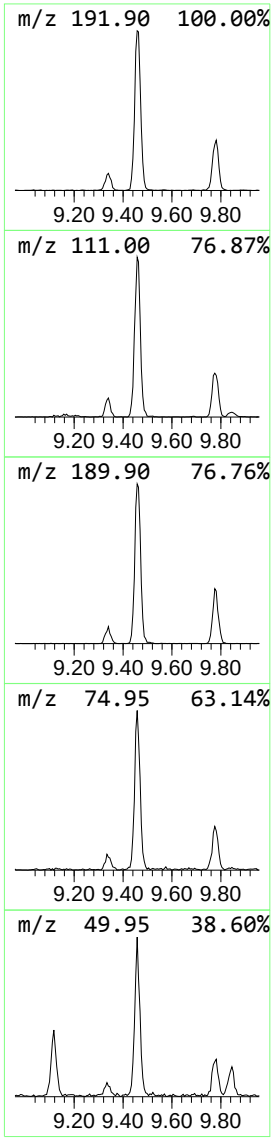
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.457	4.45 ng/ul	220637	Naphthalene-d8	10.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	97
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
3	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	94
4	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94
5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	94





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Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

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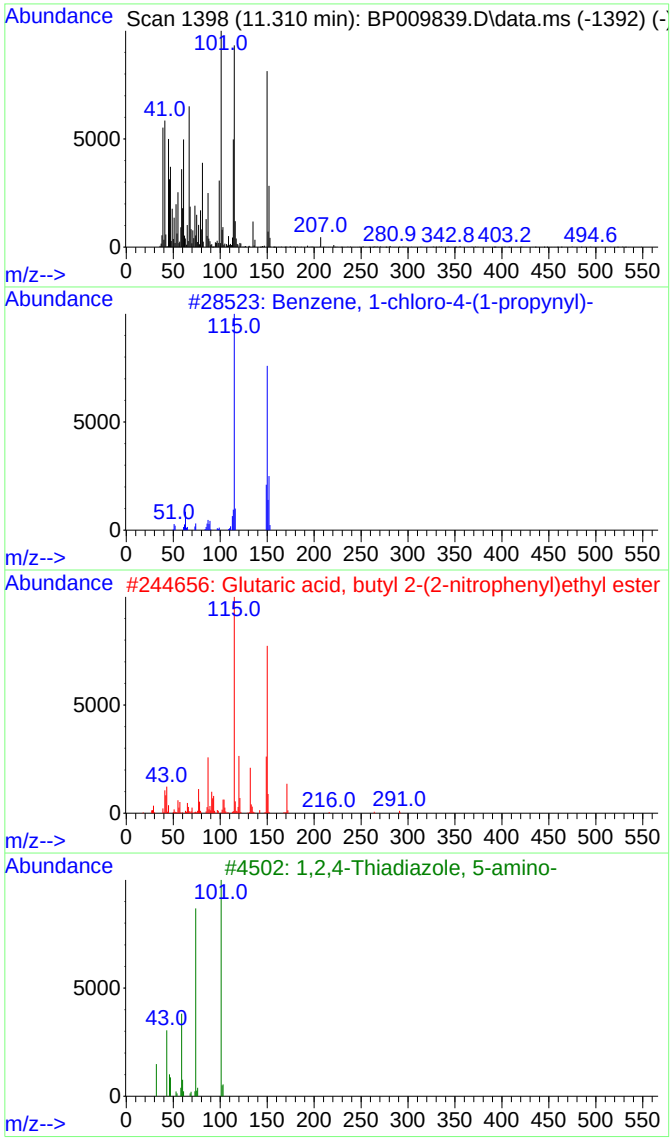
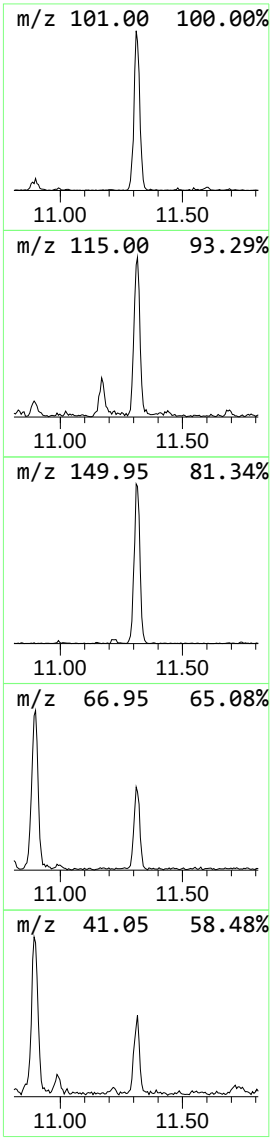
Peak Number 7 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.310	3.53 ng/ul	174891	Naphthalene-d8	10.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-4-(1-propynyl)-	150	C9H7Cl	002809-65-6	38
2	Glutaric acid, butyl 2-(2-nitrop...	337	C17H23NO6	1000377-52-2	27
3	1,2,4-Thiadiazole, 5-amino-	101	C2H3N3S	007552-07-0	14
4	Carbonic acid, monoamide, N-pent...	207	C9H18ClNO2	1010420-19-8	11
5	Benzoic acid, 3,5-dimethyl-	150	C9H10O2	000499-06-9	11





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Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

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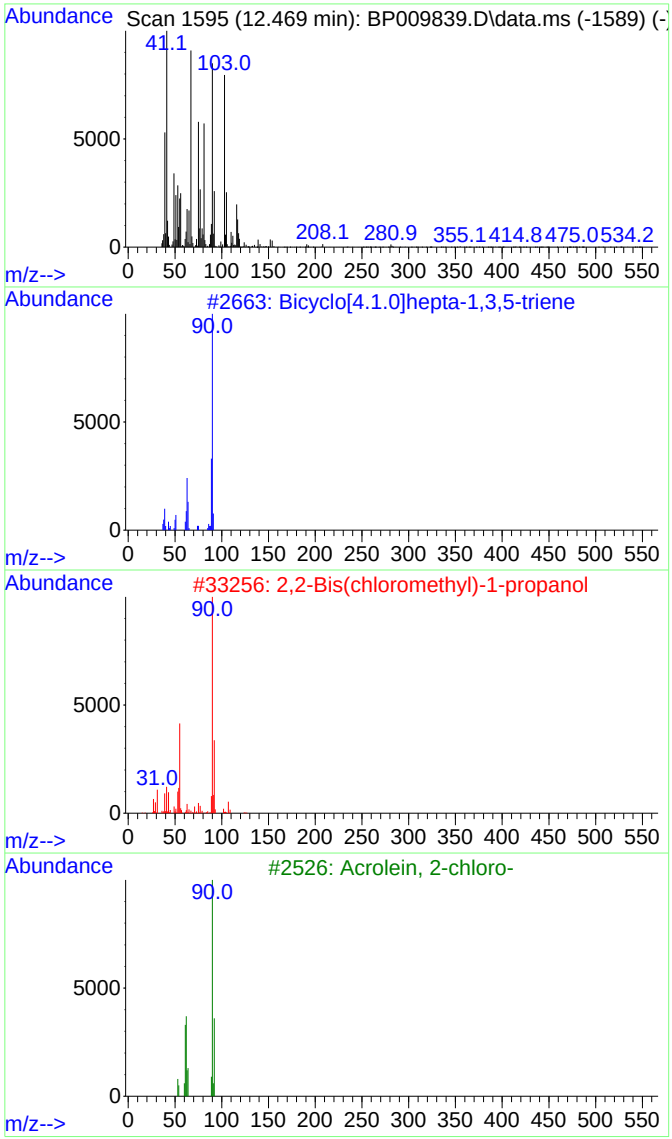
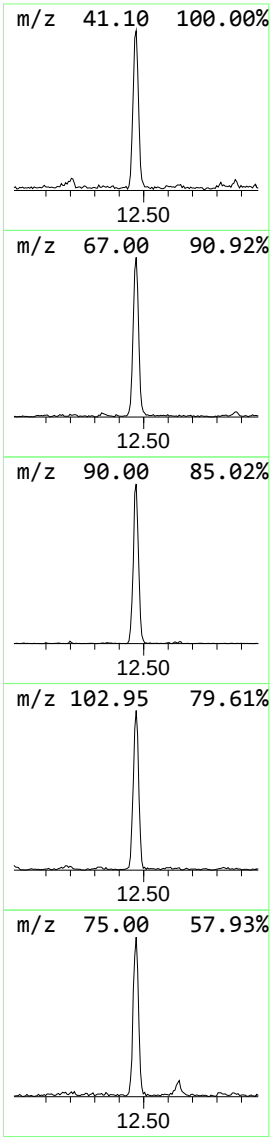
Peak Number 8 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	4.74 ng/ul	234817	Naphthalene-d8	10.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	38
2	2,2-Bis(chloromethyl)-1-propanol	156	C5H10Cl2O	005355-54-4	38
3	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	37
4	1H-1,2,3-Triazol-1-amine, 4-phen...	248	C15H12N4	113261-37-3	33
5	1,1-Dimethyl-1-sila-4-thiacycloh...	146	C6H14SSi	082316-78-7	25





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Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

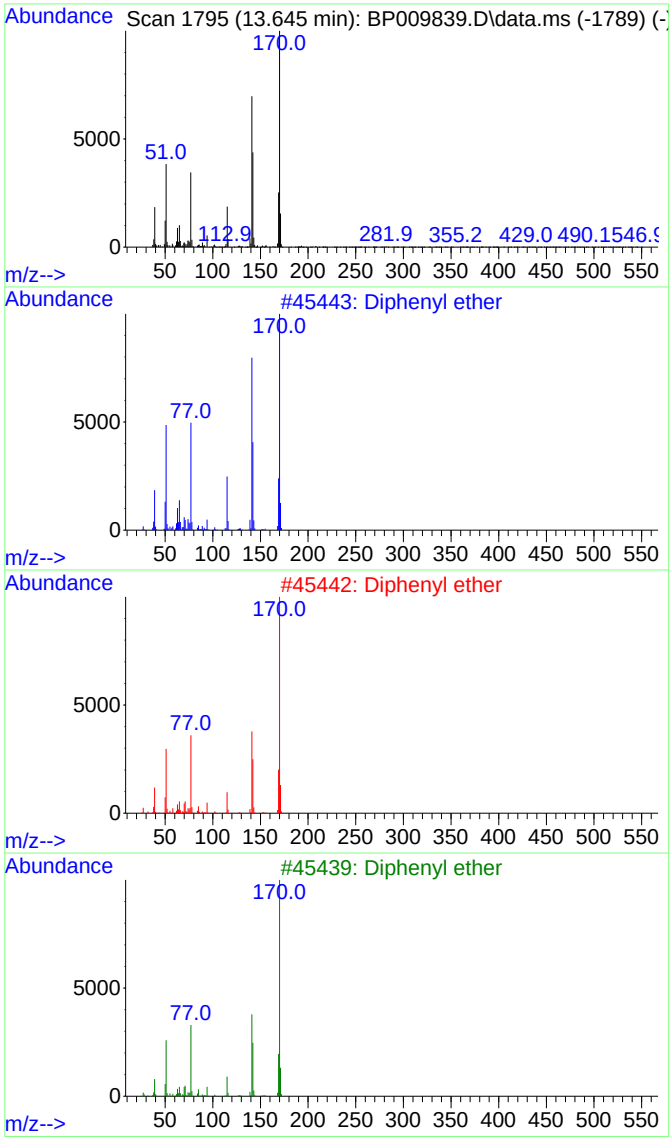
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

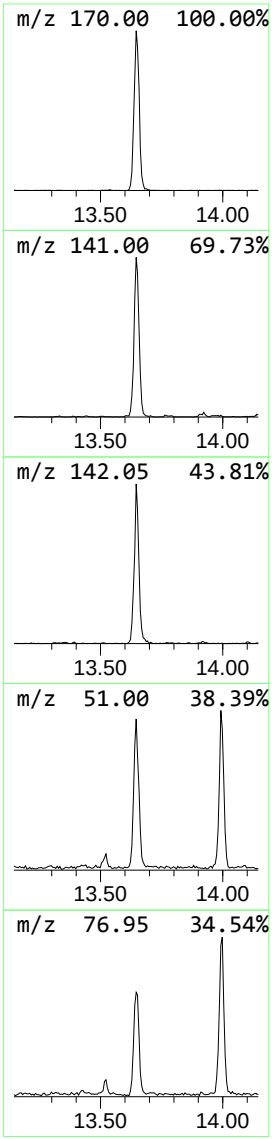
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Diphenyl ether Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.645	2.18 ng/ul	158442	Acenaphthene-d10	14.592

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl ether	170	C12H10O	000101-84-8	96
2	Diphenyl ether	170	C12H10O	000101-84-8	87
3	Diphenyl ether	170	C12H10O	000101-84-8	81
4	Diphenyl ether	170	C12H10O	000101-84-8	74
5	Diphenyl ether	170	C12H10O	000101-84-8	58





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Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

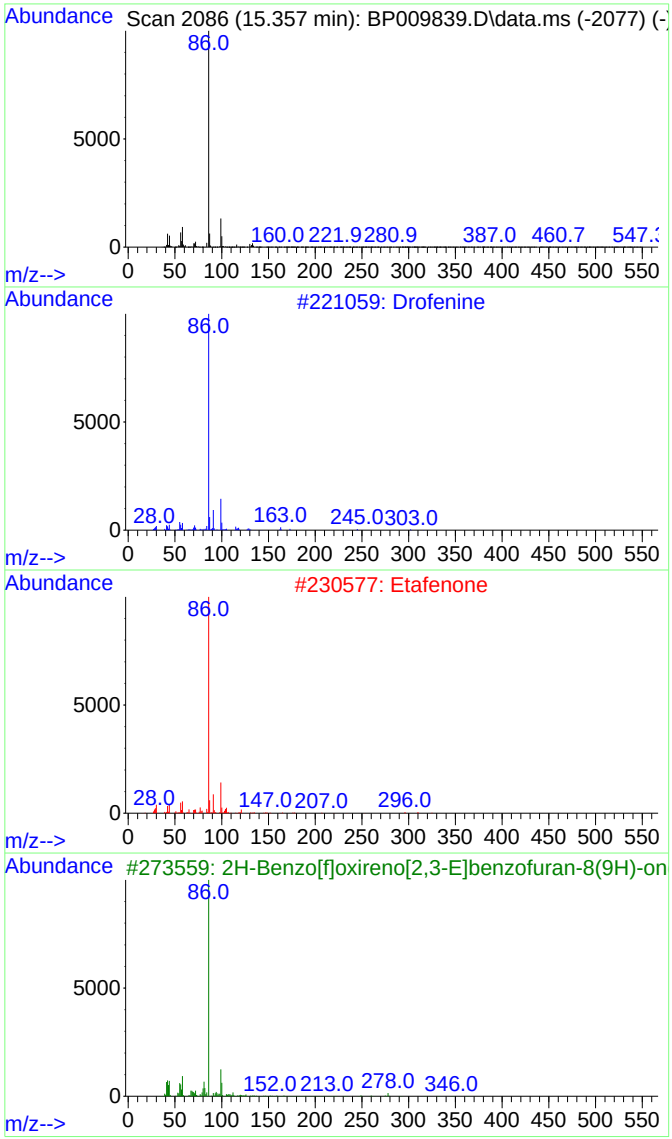
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

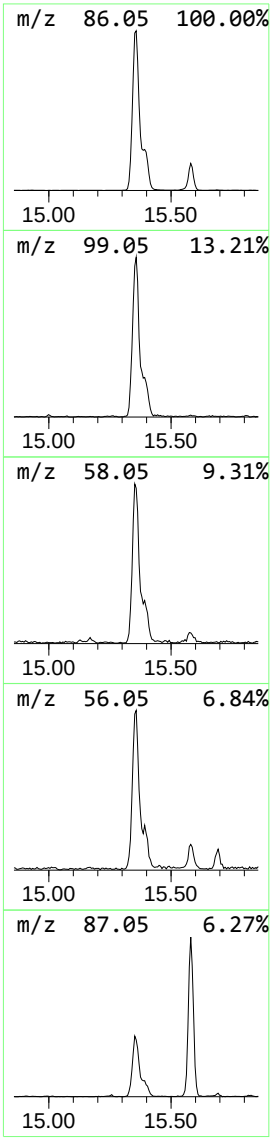
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Drofenine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.357	11.70 ng/ul	849368	Acenaphthene-d10	14.592

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Drofenine	317	C20H31NO2	001679-76-1	78
2	Etafenone	325	C21H27NO2	000090-54-0	78
3	2H-Benzo[f]oxireno[2,3-E]benzofu...	364	C21H36N2O3	1000327-59-9	78
4	2-(Diethylamino)ethyl 3-methyl-2...	291	C18H29NO2	026878-41-1	78
5	O-Ethyl S-2-diethylaminoethyl et...	253	C10H24NO2PS	021738-25-0	78





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Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

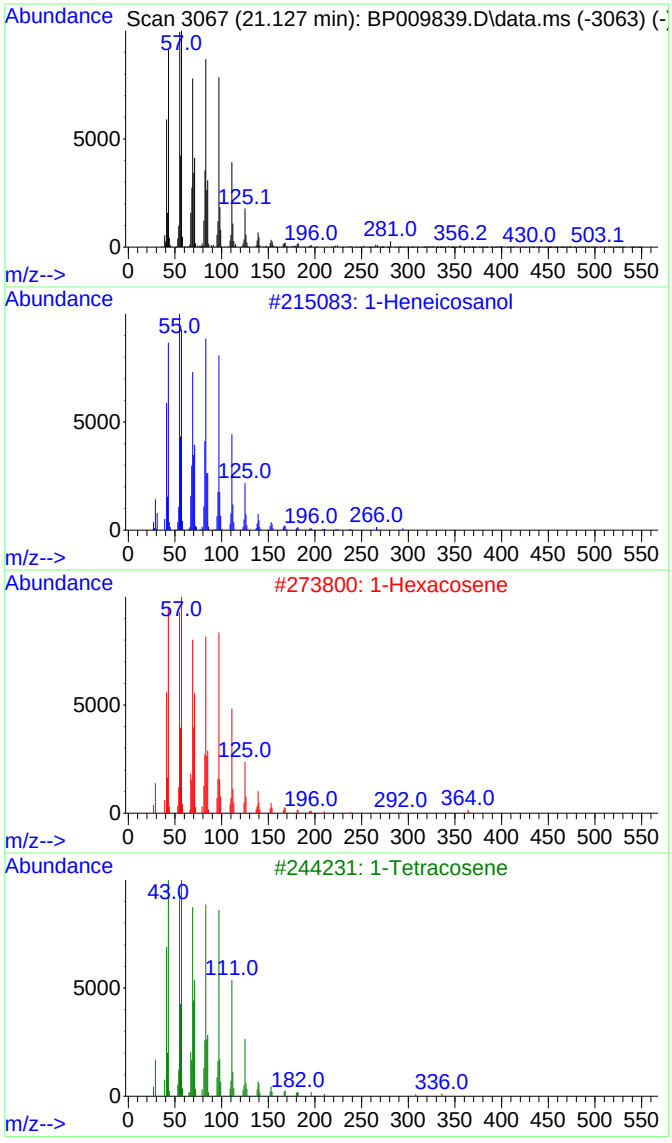
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

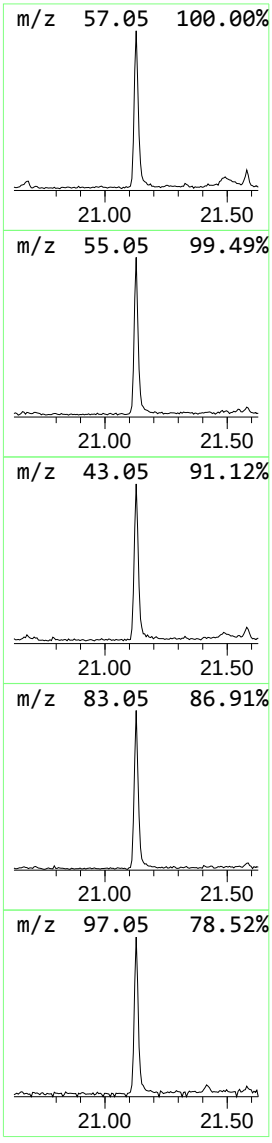
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1-Heneicosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	4.45 ng/ul	520085	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	94
2	1-Hexacosene	364	C26H52	018835-33-1	94
3	1-Tetracosene	336	C24H48	010192-32-2	94
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009839.D
Acq On : 14 Apr 2022 21:28
Operator : CG/JU
Sample : N2293-17
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.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
Butane, 2-metho...	3.122	8.0	ng/ul	247616	1	7.963	620599	20.0
unknown-01	8.257	2.2	ng/ul	67245	1	7.963	620599	20.0
1,2-Ethanediami...	9.204	9.4	ng/ul	291152	1	7.963	620599	20.0
unknown-02	9.287	2.5	ng/ul	75930	1	7.963	620599	20.0
Benzene, 1-brom...	9.457	4.5	ng/ul	220637	2	10.769	990808	20.0
unknown-03	11.310	3.5	ng/ul	174891	2	10.769	990808	20.0
unknown-04	12.469	4.7	ng/ul	234817	2	10.769	990808	20.0
Diphenyl ether	13.645	2.2	ng/ul	158442	3	14.592	1452080	20.0
Drofenine	15.357	11.7	ng/ul	849368	3	14.592	1452080	20.0
1-Heneicosanol	21.127	4.5	ng/ul	520085	5	21.422	2335800	20.0