LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

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-BP050124.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP020291.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	18	22	30	rVB	19000	30105	1.22%	0.109%
2	3.505	67	71	75	rBV	8549	15122	0.61%	0.055%
3	3.617	87	90	110	rBV	181360	357898	14.47%	1.293%
4	3.911	137	140	143	rVB2	3235	3823	0.15%	0.014%
5	5.128	342	347	350	rBV7	1492	2980	0.12%	0.011%
6	5.517	410	413	419	rBV8	1188	2558	0.10%	0.009%
7	5.999	491	495	502	rBV3	12345	24136	0.98%	0.087%
8	6.805	625	632	653	rBV	255670	551121	22.28%	1.990%
9	6.969	654	660	670	rBV	216876	374706	15.15%	1.353%
10	7.152	684	691	706	rBV	278872	508671	20.56%	1.837%
11	7.611	763	769	779	rBV	230653	420295	16.99%	1.518%
12	8.328	883	891	913	rBV	347255	726539	29.37%	2.624%
13	8.775	959	967	985	rBV	303799	561439	22.70%	2.028%
14	9.116	1022	1025	1030	rVB5	3076	4381	0.18%	0.016%
15	9.363	1061	1067	1076	rBV4	7737	15790	0.64%	0.057%
16	9.493	1080	1089	1104	rBV	202182	405933	16.41%	1.466%
17	9.705	1122	1125	1131	rVB8	1956	2927	0.12%	0.011%
18	10.028	1172	1180	1200	rBV	272886	654595	26.46%	2.364%
19	10.216	1205	1212	1231	rVB4	19956	73087	2.95%	0.264%
20	10.387	1234	1241	1248	rBV	341168	581409	23.50%	2.100%
21	10.552	1261	1269	1299	rBV	291446	716779	28.98%	2.589%
22	10.999	1341	1345	1346	rBV3	2270	3350	0.14%	0.012%
23	11.187	1370	1377	1384	rBV4	7884	23223	0.94%	0.084%
24	11.993	1509	1514	1522	rVV3	5774	14445	0.58%	0.052%
25	12.081	1524	1529	1535	rVB7	3290	5483	0.22%	0.020%
26	12.293	1562	1565	1572	rVB5	2873	5449	0.22%	0.020%
27	13.163	1708	1713	1718	rVB4	7185	12177	0.49%	0.044%
28	13.240	1721	1726	1731	rBV8	2376	4187	0.17%	0.015%
29	13.434	1755	1759	1762	rBV6	2039	2706	0.11%	0.010%
30	13.504	1767	1771	1777	rVB8	2643	4765	0.19%	0.017%
31	13.593	1780	1786	1793	rBV6	4643	11961	0.48%	0.043%
32	13.704	1798	1805	1819	rBV	728746	1096553	44.33%	3.960%
33	13.975	1843	1851	1866	rBV2	801518	1363039	55.10%	4.923%
34	14.181	1882	1886	1891	rVB7	1904	3558	0.14%	0.013%
35	14.287	1896	1904	1911	rVV	539749	809957	32.74%	2.925%
36	14.351	1911	1915	1924	rVB	32312	48126	1.95%	0.174%

```
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On    : 10 May 2024  14:58
Operator  : MA/JU
Sample    : P2370-17
Misc      :
ALS Vial  : 10   Sample Multiplier: 1
```

Instrument :
BNA_P
ClientSampleId :
A43R0

```
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-BP050124.MA.M
Title     : SVOA CALIBRATION
```

37	14.522	1936	1944	1950	rBV5	26113	66004	2.67%	0.238%
38	14.604	1955	1958	1961	rVB5	3691	4781	0.19%	0.017%
39	14.704	1970	1975	1979	rBV3	13291	21548	0.87%	0.078%
40	14.928	2009	2013	2017	rBV7	2919	4954	0.20%	0.018%
41	14.981	2019	2022	2027	rVB7	3269	5407	0.22%	0.020%
42	15.092	2034	2041	2045	rBV10	3899	8484	0.34%	0.031%
43	15.163	2051	2053	2060	rVB8	2821	4698	0.19%	0.017%
44	15.310	2071	2078	2085	rBV	1017136	1634792	66.09%	5.904%
45	15.363	2085	2087	2096	rVV	32796	53440	2.16%	0.193%
46	15.469	2099	2105	2106	rBV6	3094	5375	0.22%	0.019%
47	15.492	2106	2109	2113	rBV6	3107	4467	0.18%	0.016%
48	15.675	2137	2140	2145	rVB3	9191	13135	0.53%	0.047%
49	15.828	2162	2166	2175	rVB3	8114	19073	0.77%	0.069%
50	15.898	2175	2178	2179	rBV3	3252	3020	0.12%	0.011%
51	16.081	2201	2209	2216	rBV3	10970	32836	1.33%	0.119%
52	16.245	2232	2237	2239	rBV6	3650	5918	0.24%	0.021%
53	16.410	2254	2265	2268	rBV10	12430	33447	1.35%	0.121%
54	16.651	2302	2306	2309	rBV5	6582	11572	0.47%	0.042%
55	16.686	2310	2312	2316	rVB4	5739	5629	0.23%	0.020%
56	16.781	2322	2328	2335	rVB3	23766	47972	1.94%	0.173%
57	16.869	2338	2343	2349	rVV8	13087	31983	1.29%	0.116%
58	16.945	2351	2356	2364	rVB	28739	49128	1.99%	0.177%
59	17.122	2380	2386	2390	rBV	611976	1000863	40.46%	3.615%
60	17.169	2390	2394	2399	rVV	506024	780186	31.54%	2.818%
61	17.234	2399	2405	2417	rVB	1037205	1884233	76.17%	6.805%
62	17.575	2456	2463	2470	rBV2	25499	66892	2.70%	0.242%
63	17.675	2476	2480	2485	rBV3	18654	30594	1.24%	0.110%
64	18.045	2538	2543	2546	rBV	51363	78578	3.18%	0.284%
65	18.092	2546	2551	2561	rVV2	221669	401563	16.23%	1.450%
66	18.186	2565	2567	2572	rVV	27046	43292	1.75%	0.156%
67	18.257	2573	2579	2582	rVV3	78330	155386	6.28%	0.561%
68	18.292	2583	2585	2591	rVB2	39215	46065	1.86%	0.166%
69	18.581	2630	2634	2639	rBV	54379	77676	3.14%	0.281%
70	18.639	2640	2644	2649	rVB2	41928	65708	2.66%	0.237%
71	19.010	2703	2707	2711	rVB2	43726	61298	2.48%	0.221%
72	19.169	2729	2734	2740	rBV4	70184	152948	6.18%	0.552%
73	19.292	2749	2755	2765	rVB	844914	1274528	51.52%	4.603%
74	19.445	2777	2781	2786	rBV2	120512	186205	7.53%	0.673%
75	19.639	2809	2814	2817	rBV	1886892	2473668	100.00%	8.934%
76	19.669	2817	2819	2825	rVB	882056	976436	39.47%	3.527%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

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-BP050124.MA.M
Title : SVOA CALIBRATION

77	20.198	2907	2909	2915	rVB	77804	93219	3.77%	0.337%
78	20.310	2925	2928	2932	rBV	51782	69965	2.83%	0.253%
79	20.510	2960	2962	2964	rBV	48104	45988	1.86%	0.166%
80	21.316	3095	3099	3104	rBV3	174809	260104	10.51%	0.939%
81	21.627	3145	3152	3158	rBV2	833272	1832225	74.07%	6.617%
82	21.686	3158	3162	3168	rVB	280024	460908	18.63%	1.665%
83	23.945	3540	3546	3553	rBV	183866	475189	19.21%	1.716%
84	24.780	3679	3688	3694	rBV	828124	2111584	85.36%	7.626%
85	25.004	3718	3726	3733	rBV	399853	1072039	43.34%	3.872%

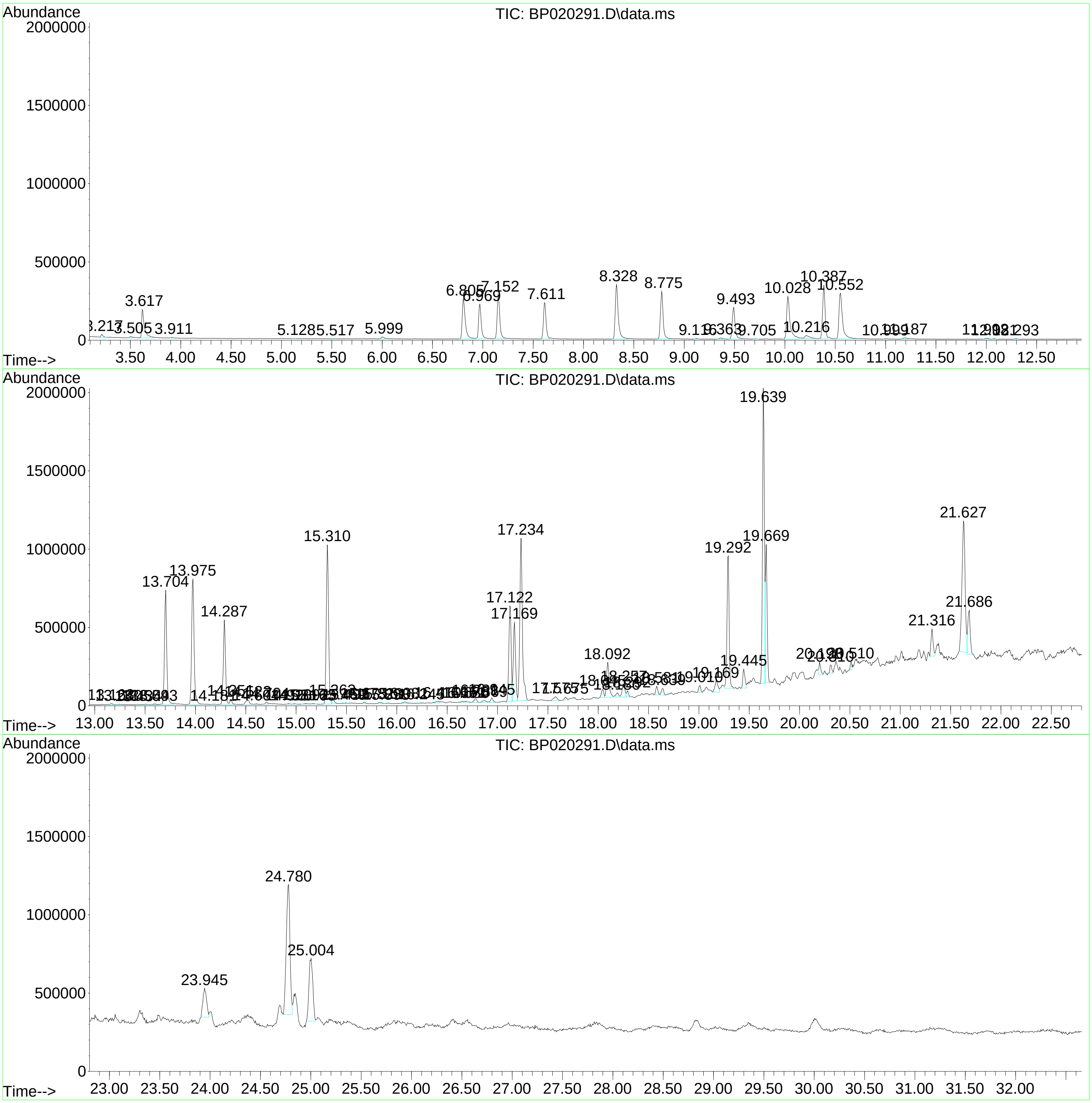
Sum of corrected areas: 27688276

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On    : 10 May 2024  14:58
Operator  : MA/JU
Sample    : P2370-17
Misc      :
ALS Vial  : 10   Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title  : SVOA CALIBRATION

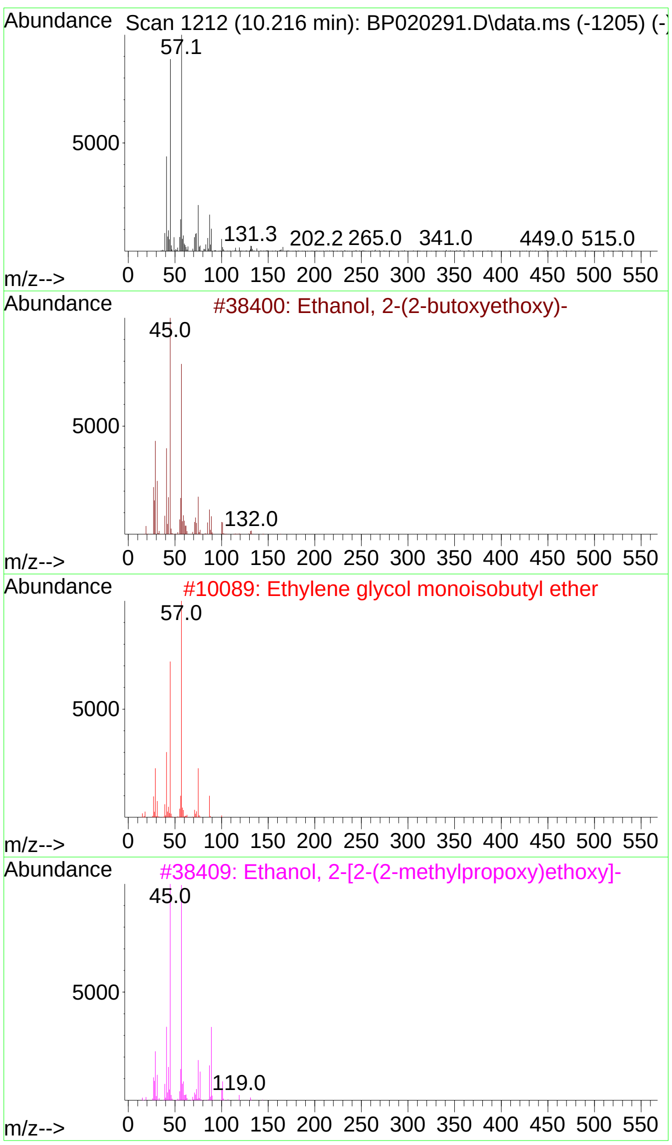
TIC Library    : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P
```

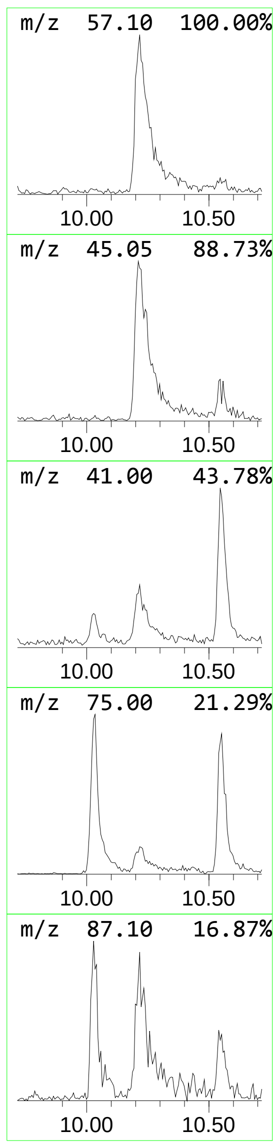
Instrument :
BNA_P
ClientSampleId :
A43R0

```
**********************************************************************
Peak Number  1  Ethanol, 2-(2-butoxyethoxy)-    Concentration Rank  8
```

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.216	2.51 ng/ul	73087	Naphthalene-d8	10.387

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78
2	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	64
4	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	58
5	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	53





Library Search Compound Report

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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On    : 10 May 2024  14:58
Operator  : MA/JU
Sample    : P2370-17
Misc      :
ALS Vial  : 10   Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title  : SVOA CALIBRATION

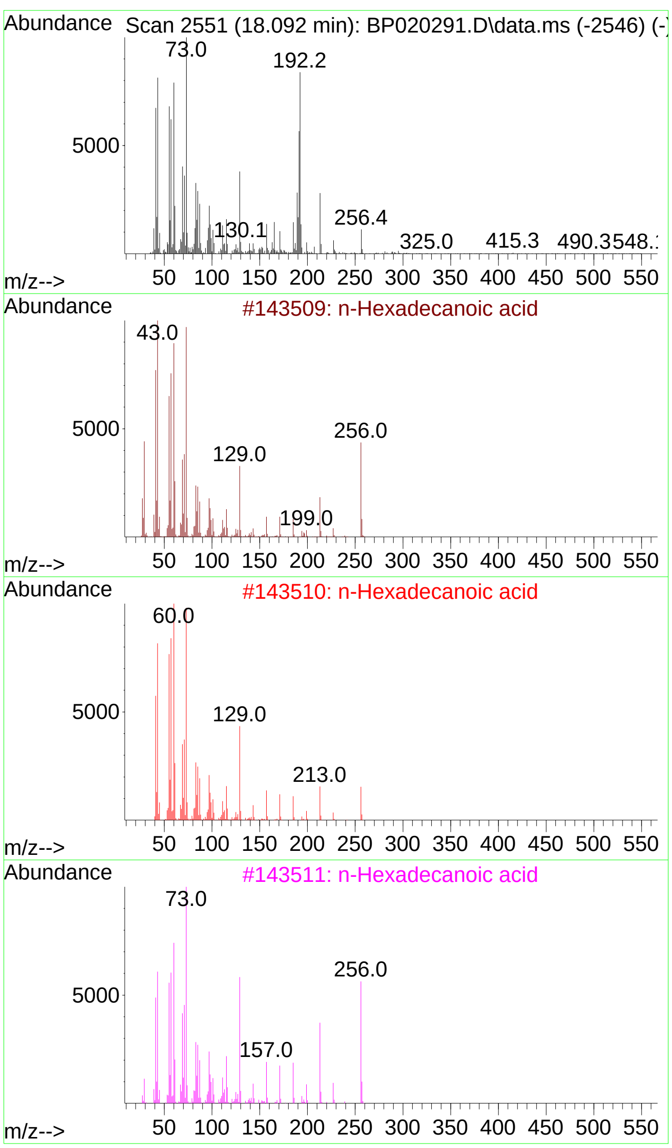
TIC Library    : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P
```

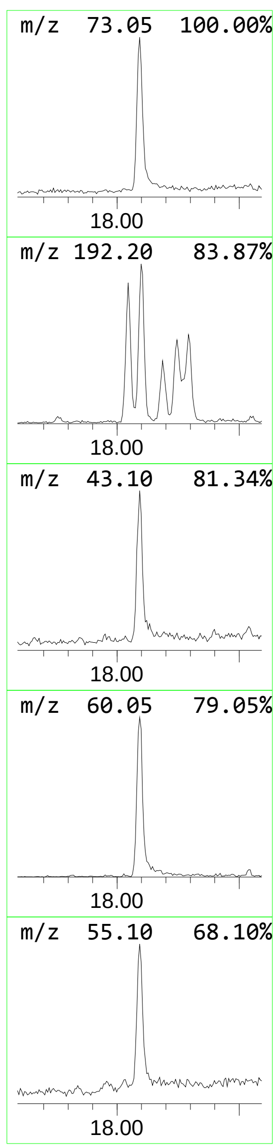
Instrument :
BNA_P
ClientSampleId :
A43R0

```
*********************************************************************
Peak Number  2  n-Hexadecanoic acid            Concentration Rank  3
```

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	8.02 ng/ul	401563	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	64





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

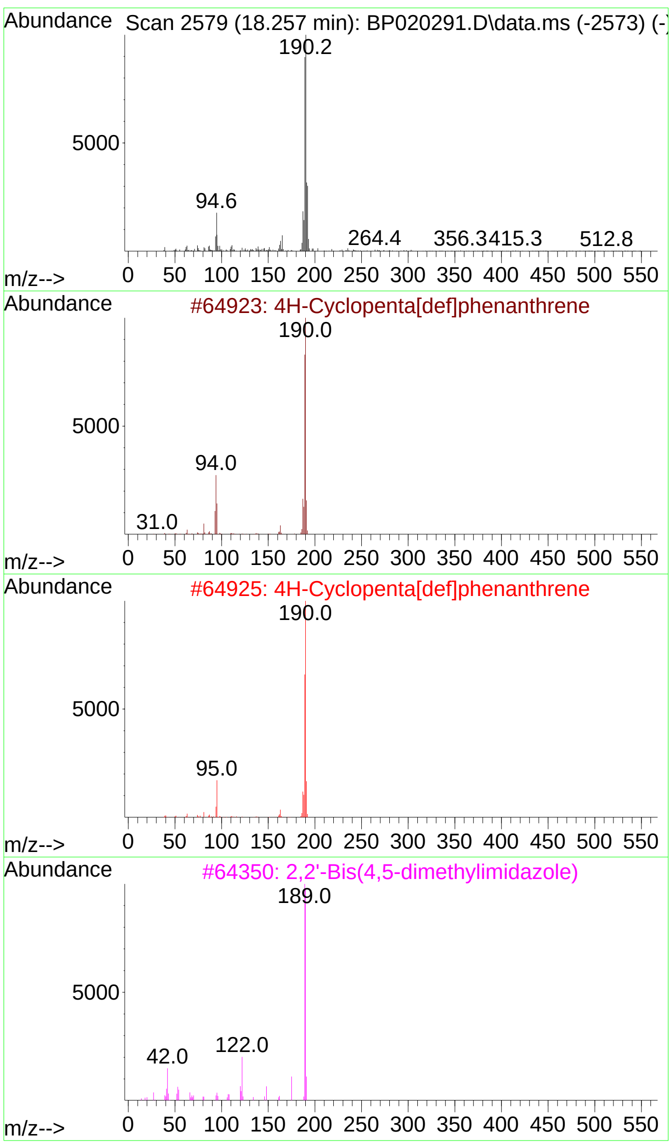
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

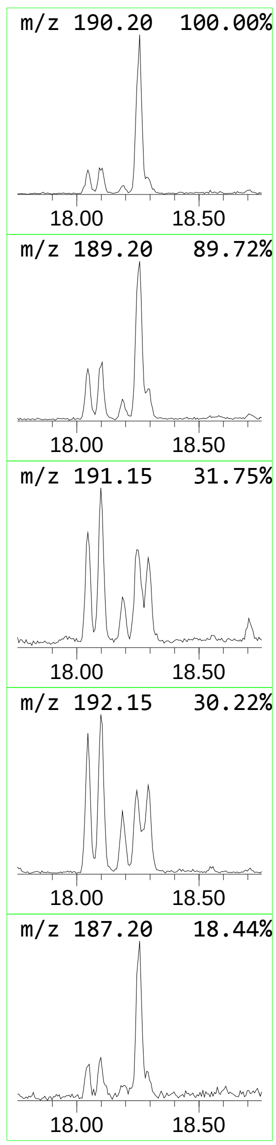
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.257	3.11 ng/ul	155386	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	53
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45





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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On    : 10 May 2024  14:58
Operator  : MA/JU
Sample    : P2370-17
Misc      :
ALS Vial  : 10   Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title  : SVOA CALIBRATION

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

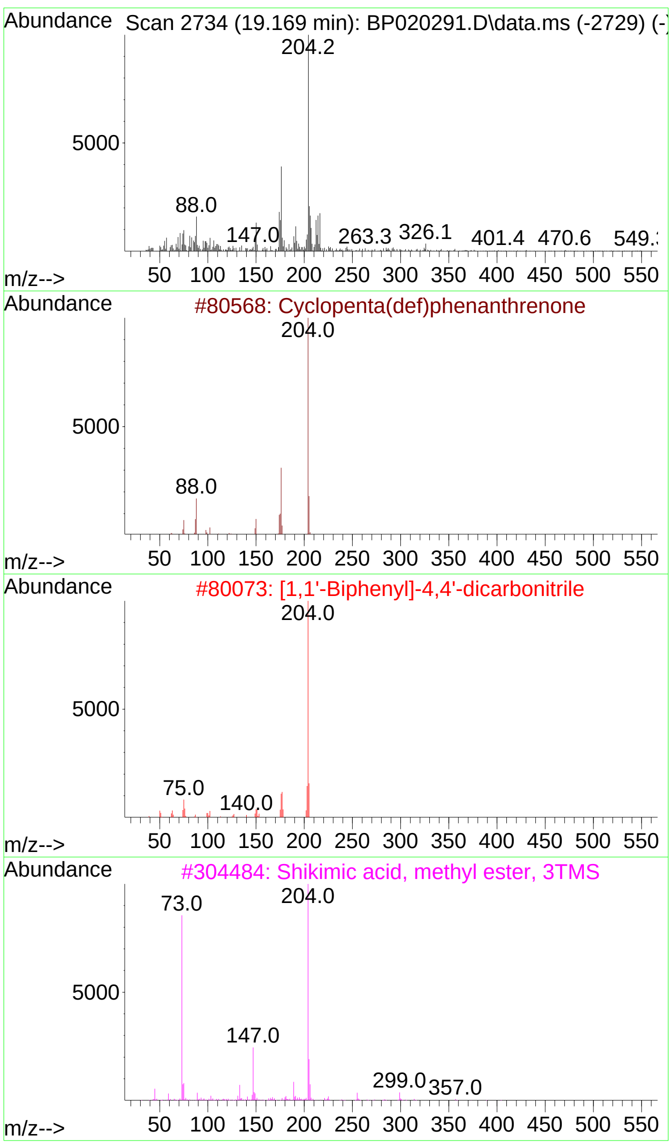
Instrument :
BNA_P
ClientSampleId :
A43R0

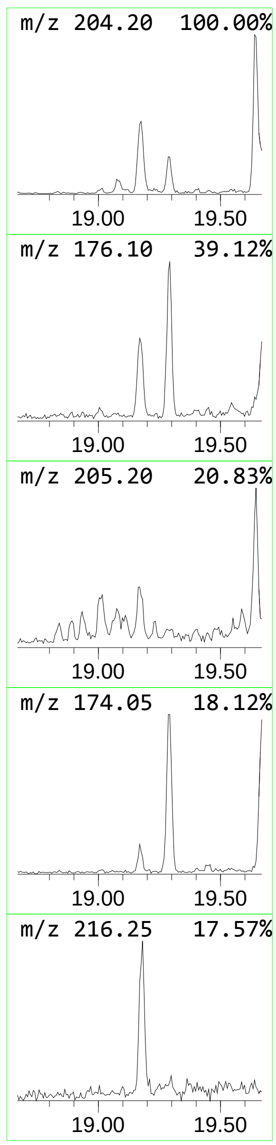
****

Peak Number 4 Cyclopenta(def)phenanthrenone   Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.169	3.06 ng/ul	152948	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	94
2	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	47
3	Shikimic acid, methyl ester, 3TMS	404	C17H36O5Si3	1000510-20-9	46
4	[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	46
5	5H-Dibenzo[a,d]cycloheptene, 5-m...	204	C16H12	002975-79-3	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

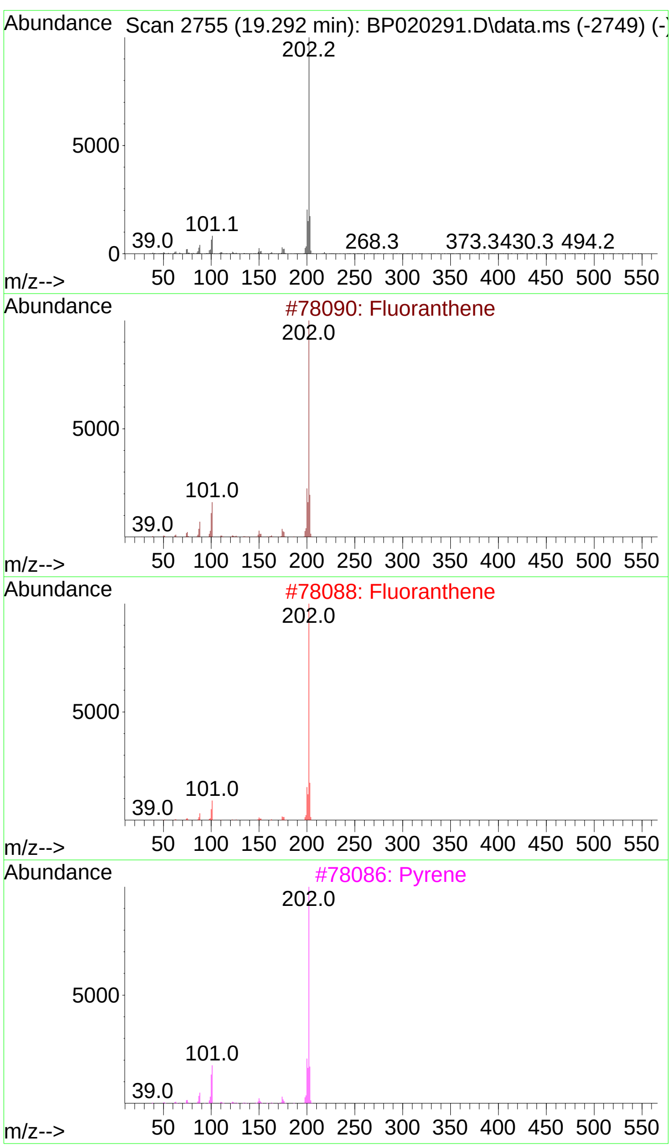
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

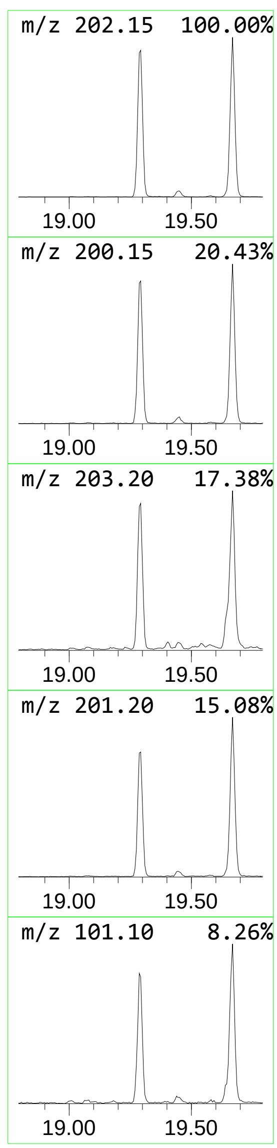
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Fluoran thene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.292	25.47 ng/ul	1274530	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene	202	C16H10	000206-44-0	96
2	Fluoranthene	202	C16H10	000206-44-0	95
3	Pyrene	202	C16H10	000129-00-0	90
4	Pyrene	202	C16H10	000129-00-0	90
5	Pyrene	202	C16H10	000129-00-0	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

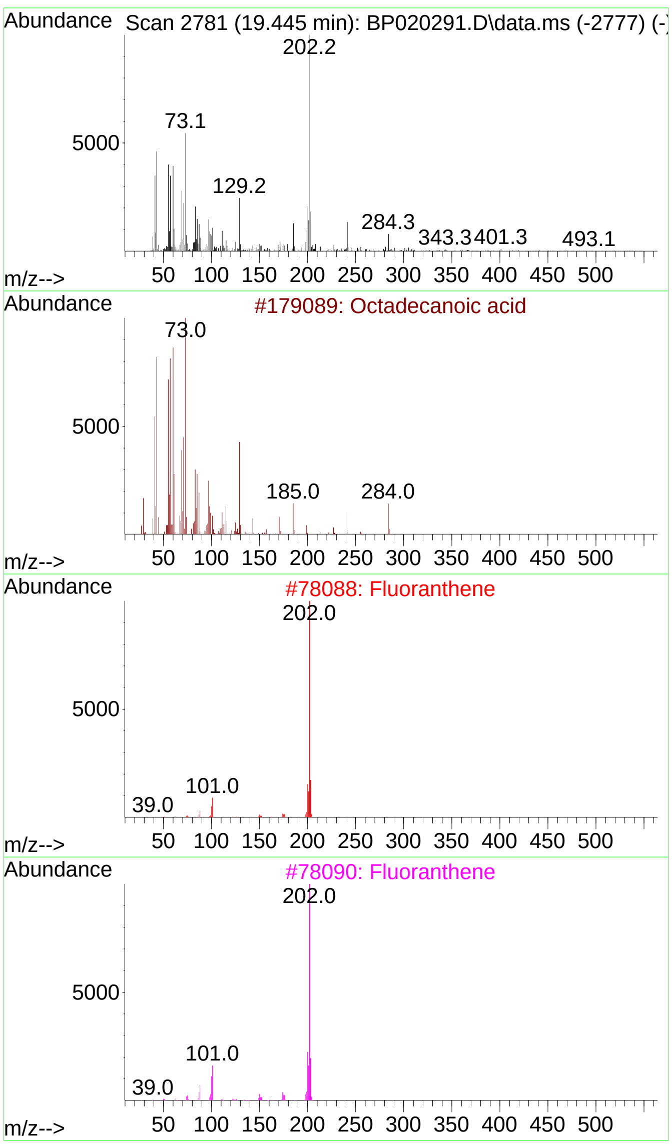
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

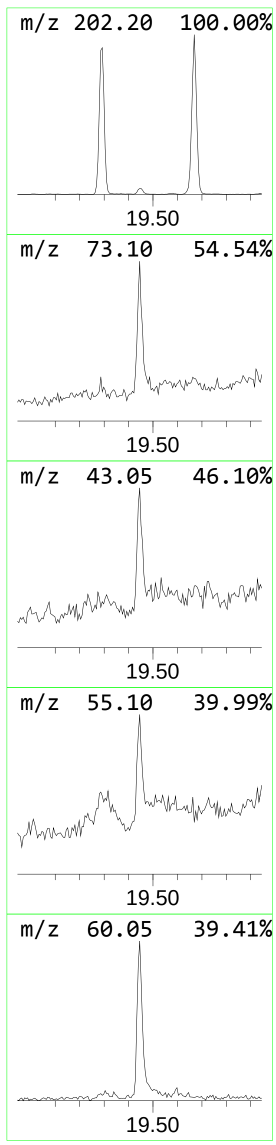
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	2.03 ng/ul	186205	Chrysene-d12	21.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	93
2	Fluoranthene	202	C16H10	000206-44-0	60
3	Fluoranthene	202	C16H10	000206-44-0	50
4	Fluoranthene	202	C16H10	000206-44-0	50
5	Pyrene	202	C16H10	000129-00-0	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

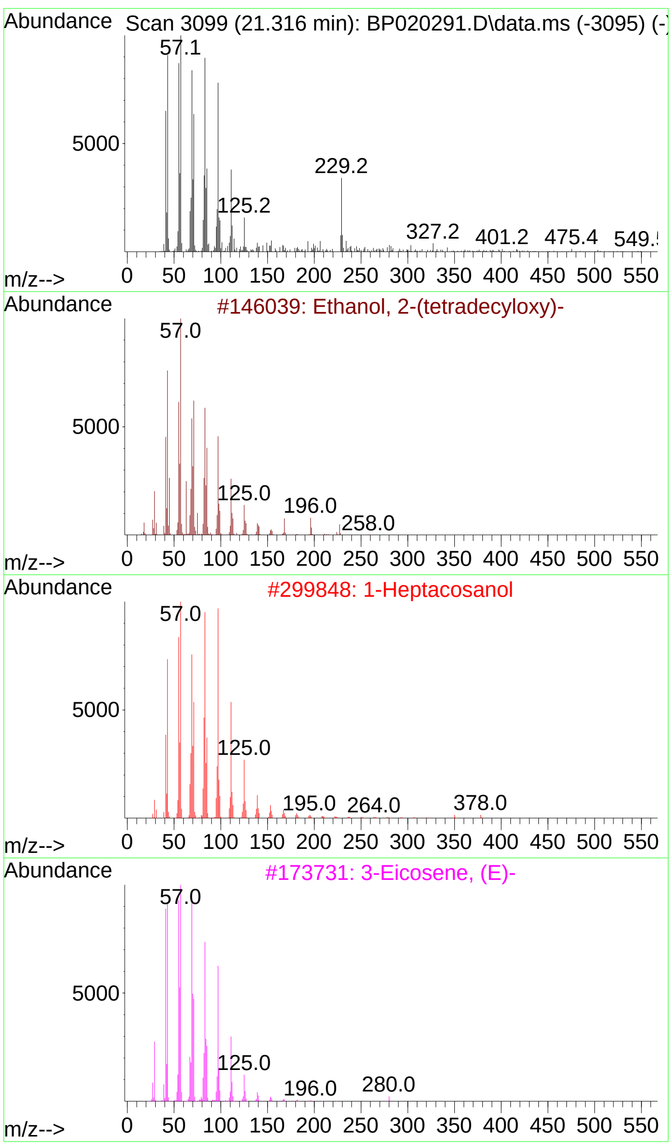
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

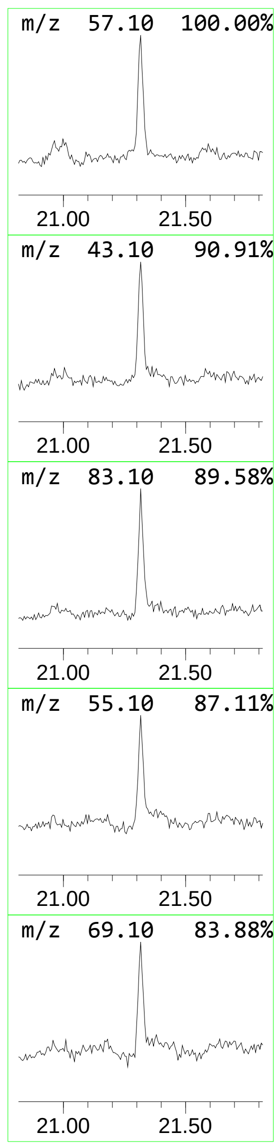
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Ethanol, 2-(tetradecyloxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.316	2.84 ng/ul	260104	Chrysene-d12	21.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	95
2	1-Heptacosanol	396	C27H56O	002004-39-9	94
3	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
5	9-Eicosene, (E)-	280	C20H40	074685-29-3	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

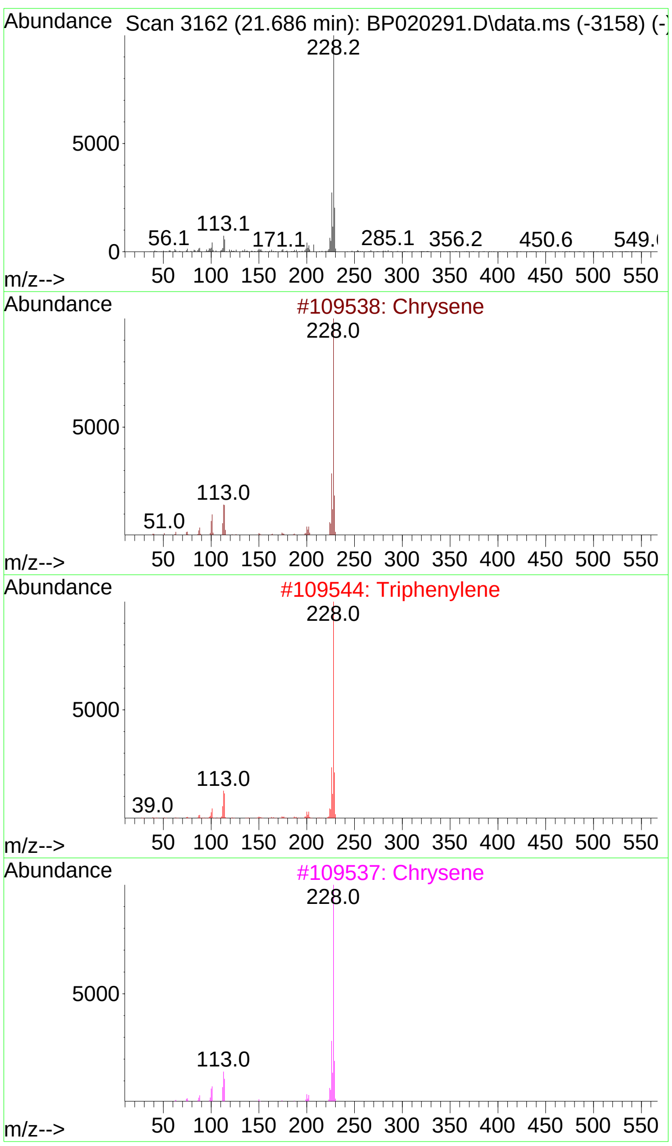
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

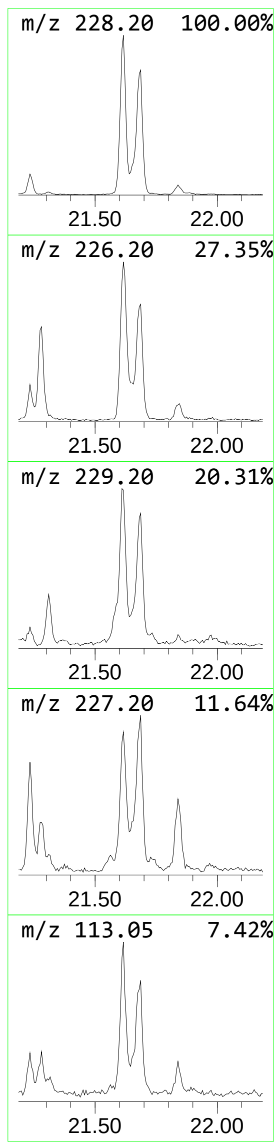
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Chry sene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.686	5.03 ng/ul	460908	Chrysene-d12	21.633

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chrysene	228	C18H12	000218-01-9	94
2	Triphenylene	228	C18H12	000217-59-4	93
3	Chrysene	228	C18H12	000218-01-9	91
4	Chrysene	228	C18H12	000218-01-9	91
5	Triphenylene	228	C18H12	000217-59-4	90





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

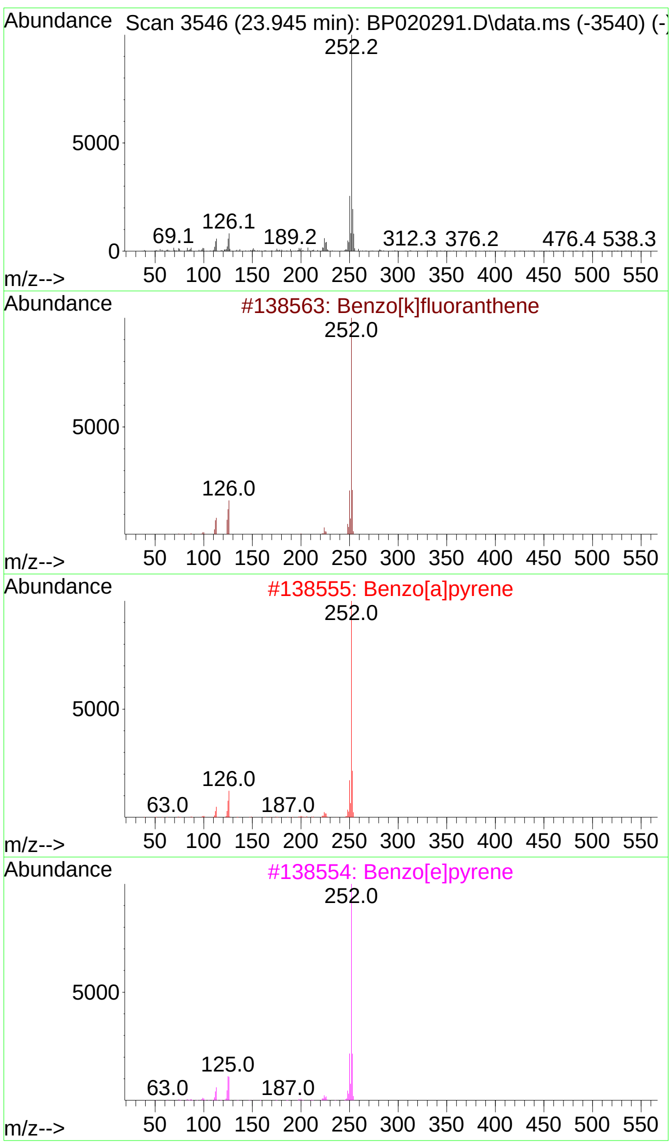
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

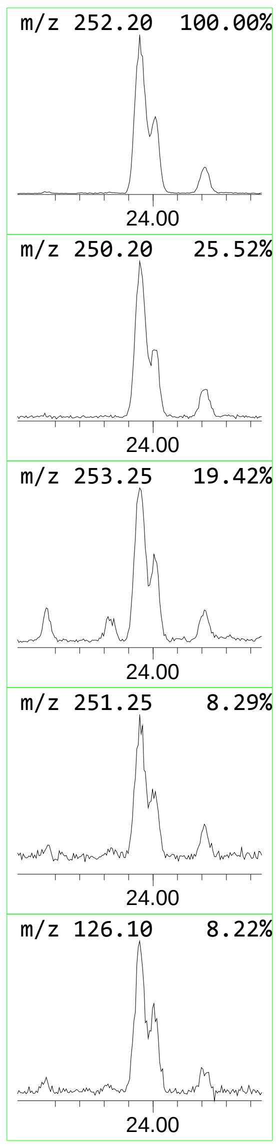
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzo[k]fluoranthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.945	8.87 ng/ul	475189	Perylene-d12	25.004

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
2	Benzo[a]pyrene	252	C20H12	000050-32-8	81
3	Benzo[e]pyrene	252	C20H12	000192-97-2	81
4	Benzo[a]pyrene	252	C20H12	000050-32-8	81
5	Perylene	252	C20H12	000198-55-0	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
Data File : BP020291.D
Acq On : 10 May 2024 14:58
Operator : MA/JU
Sample : P2370-17
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43R0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.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
Ethanol, 2-(2-b...	10.216	2.5	ng/ul	73087	2	10.387	581409	20.0
n-Hexadecanoic ...	18.092	8.0	ng/ul	401563	4	17.128	1000860	20.0
4H-Cyclopenta[d...	18.257	3.1	ng/ul	155386	4	17.128	1000860	20.0
Cyclopenta(def)...	19.169	3.1	ng/ul	152948	4	17.128	1000860	20.0
Fluoran thene	19.292	25.5	ng/ul	1274530	4	17.128	1000860	20.0
Octadecanoic acid	19.445	2.0	ng/ul	186205	5	21.633	1832230	20.0
Ethanol, 2-(tet...	21.316	2.8	ng/ul	260104	5	21.633	1832230	20.0
Chry sene	21.686	5.0	ng/ul	460908	5	21.633	1832230	20.0
Benzo[k]fluoran...	23.945	8.9	ng/ul	475189	6	25.004	1072040	20.0