LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

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-BP071024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP020944.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.211	34	38	42	rBV	59658	69477	1.32%	0.090%
2	3.252	42	45	48	rVV	36162	45026	0.86%	0.058%
3	3.287	48	51	59	rVB	46454	61847	1.18%	0.080%
4	3.528	88	92	101	rBV	91936	130226	2.48%	0.168%
5	3.670	113	116	127	rBV	253630	383795	7.31%	0.495%
6	3.975	163	168	175	rVB	151097	204793	3.90%	0.264%
7	5.393	404	409	411	rBV	22179	36190	0.69%	0.047%
8	5.752	463	470	476	rBV	61788	100091	1.91%	0.129%
9	5.834	476	484	496	rVB3	17173	49366	0.94%	0.064%
10	6.222	542	550	555	rBV2	22899	41406	0.79%	0.053%
11	6.699	626	631	637	rVV	52053	80663	1.54%	0.104%
12	6.805	644	649	655	rVV	73861	124054	2.36%	0.160%
13	6.869	655	660	663	rVV	60576	98281	1.87%	0.127%
14	6.911	663	667	678	rVV2	186405	402975	7.68%	0.519%
15	7.058	683	692	696	rBV	529695	809198	15.42%	1.043%
16	7.105	696	700	707	rVB	136390	223881	4.27%	0.289%
17	7.252	719	725	733	rBV	633770	1043142	19.87%	1.345%
18	7.352	733	742	754	rVB2	211996	446198	8.50%	0.575%
19	7.599	776	784	790	rVB10	17215	42791	0.82%	0.055%
20	7.705	790	802	813	rBV	659722	1316866	25.09%	1.697%
21	7.816	816	821	832	rVB4	40091	95761	1.82%	0.123%
22	8.058	856	862	871	rVB	358299	613324	11.68%	0.791%
23	8.422	917	924	932	rVV	512914	930521	17.73%	1.199%
24	8.493	932	936	943	rVB3	15521	34674	0.66%	0.045%
25	8.758	975	981	991	rVB4	24299	76672	1.46%	0.099%
26	8.858	991	998	1005	rBV	616118	1055369	20.11%	1.360%
27	8.928	1005	1010	1017	rVB	161160	268872	5.12%	0.347%
28	9.116	1036	1042	1050	rVB3	86927	140010	2.67%	0.180%
29	9.410	1088	1092	1100	rVB3	55511	98882	1.88%	0.127%
30	9.575	1113	1120	1131	rBV	518630	995344	18.96%	1.283%
31	9.675	1131	1137	1143	rVV	176172	295269	5.63%	0.381%
32	9.899	1161	1175	1185	rVB4	301039	640627	12.20%	0.826%
33	10.028	1190	1197	1204	rVB3	75885	142734	2.72%	0.184%
34	10.116	1204	1212	1221	rBV	830910	1636682	31.18%	2.110%
35	10.263	1233	1237	1244	rVB7	18966	44107	0.84%	0.057%
36	10.340	1244	1250	1253	rBV7	15882	35803	0.68%	0.046%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

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-BP071024.MA.M
Title : SVOA CALIBRATION

37	10.475	1265	1273	1278	rBV	978107	1560408	29.73%	2.011%
38	10.522	1278	1281	1287	rVV	172222	278091	5.30%	0.358%
39	10.628	1292	1299	1312	rVB	553544	1115727	21.26%	1.438%
40	10.763	1317	1322	1327	rVB4	28523	60969	1.16%	0.079%
41	10.834	1329	1334	1338	rBV4	23345	47739	0.91%	0.062%
42	11.046	1361	1370	1376	rBV8	52942	186780	3.56%	0.241%
43	11.228	1395	1401	1407	rVB	131602	237235	4.52%	0.306%
44	11.381	1422	1427	1433	rVB4	58073	114420	2.18%	0.147%
45	11.516	1445	1450	1454	rBV3	39275	82229	1.57%	0.106%
46	11.569	1455	1459	1463	rBV	53376	84869	1.62%	0.109%
47	11.657	1469	1474	1481	rVB3	38971	65952	1.26%	0.085%
48	11.722	1481	1485	1491	rVB2	26754	53366	1.02%	0.069%
49	11.799	1492	1498	1501	rBV	88095	169301	3.23%	0.218%
50	11.840	1501	1505	1516	rVB2	177588	386136	7.36%	0.498%
51	11.940	1517	1522	1527	rVB5	30565	58540	1.12%	0.075%
52	12.010	1528	1534	1540	rBV4	81988	160720	3.06%	0.207%
53	12.134	1549	1555	1563	rVB	783311	1364743	26.00%	1.759%
54	12.351	1586	1592	1598	rBV	1693024	2657042	50.62%	3.425%
55	12.410	1598	1602	1608	rVV5	91667	197304	3.76%	0.254%
56	12.475	1609	1613	1619	rVB5	57040	114376	2.18%	0.147%
57	12.634	1637	1640	1643	rBV4	25776	37796	0.72%	0.049%
58	12.757	1654	1661	1664	rBV5	56574	130715	2.49%	0.168%
59	12.916	1685	1688	1693	rVB4	36722	54625	1.04%	0.070%
60	13.251	1741	1745	1751	rBV	53717	86129	1.64%	0.111%
61	13.510	1781	1789	1794	rBV3	107285	276264	5.26%	0.356%
62	13.651	1807	1813	1818	rBV	258295	494517	9.42%	0.637%
63	13.704	1818	1822	1824	rVV	158566	251391	4.79%	0.324%
64	13.746	1824	1829	1840	rVB	1862112	2495609	47.54%	3.217%
65	13.887	1849	1853	1857	rBV	92450	161368	3.07%	0.208%
66	14.028	1872	1877	1887	rBV2	1876982	3086146	58.79%	3.978%
67	14.163	1897	1900	1906	rVB4	70965	88465	1.69%	0.114%
68	14.257	1912	1916	1921	rBV4	74274	143241	2.73%	0.185%
69	14.340	1924	1930	1936	rVV	1531048	2250219	42.87%	2.901%
70	14.516	1958	1960	1963	rVB	67403	61022	1.16%	0.079%
71	14.598	1970	1974	1983	rVB2	271044	480357	9.15%	0.619%
72	14.822	2008	2012	2018	rVB	226521	345556	6.58%	0.445%
73	15.022	2042	2046	2050	rVB2	97287	137101	2.61%	0.177%
74	15.104	2055	2060	2071	rVV	326812	653343	12.45%	0.842%
75	15.193	2071	2075	2081	rVB7	110530	181137	3.45%	0.233%
76	15.351	2096	2102	2108	rVB2	2534952	3751845	71.47%	4.836%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

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-BP071024.MA.M
Title : SVOA CALIBRATION

77	15.498	2121	2127	2134	rBV	603739	1125028	21.43%	1.450%
78	15.763	2165	2172	2178	rVB4	243234	498122	9.49%	0.642%
79	16.134	2229	2235	2244	rVB	1330918	2300039	43.82%	2.965%
80	16.210	2245	2248	2252	rBV	114382	187186	3.57%	0.241%
81	16.257	2253	2256	2257	rBV3	105091	119167	2.27%	0.154%
82	16.334	2266	2269	2274	rVB2	171476	227328	4.33%	0.293%
83	16.557	2299	2307	2314	rBV4	466209	1017712	19.39%	1.312%
84	16.628	2315	2319	2326	rVB3	223078	483801	9.22%	0.624%
85	16.934	2366	2371	2376	rBV	690634	1114965	21.24%	1.437%
86	17.151	2402	2408	2416	rVB	1818737	2769397	52.76%	3.570%
87	17.257	2420	2426	2436	rVV2	2166984	4052058	77.19%	5.223%
88	17.351	2439	2442	2449	rVB	290923	457868	8.72%	0.590%
89	17.963	2542	2546	2551	rBV2	342807	534118	10.18%	0.689%
90	18.098	2565	2569	2574	rBV	1582545	2121492	40.42%	2.735%
91	19.292	2769	2772	2781	rVB3	384803	674480	12.85%	0.869%
92	19.428	2792	2795	2806	rVB2	394648	608167	11.59%	0.784%
93	19.639	2825	2831	2838	rVB	3296056	4714999	89.82%	6.078%
94	19.904	2873	2876	2881	rBV	208961	334223	6.37%	0.431%
95	20.233	2929	2932	2941	rVB	341189	627106	11.95%	0.808%
96	20.433	2963	2966	2976	rVB	279107	520790	9.92%	0.671%
97	21.604	3159	3165	3173	rVB2	1771669	2852509	54.34%	3.677%
98	23.369	3459	3465	3474	rVB2	2265796	5249192	100.00%	6.766%
99	24.716	3685	3694	3706	rVB	1777033	4962600	94.54%	6.397%
100	24.939	3724	3732	3742	rVB	1138563	3046922	58.05%	3.928%

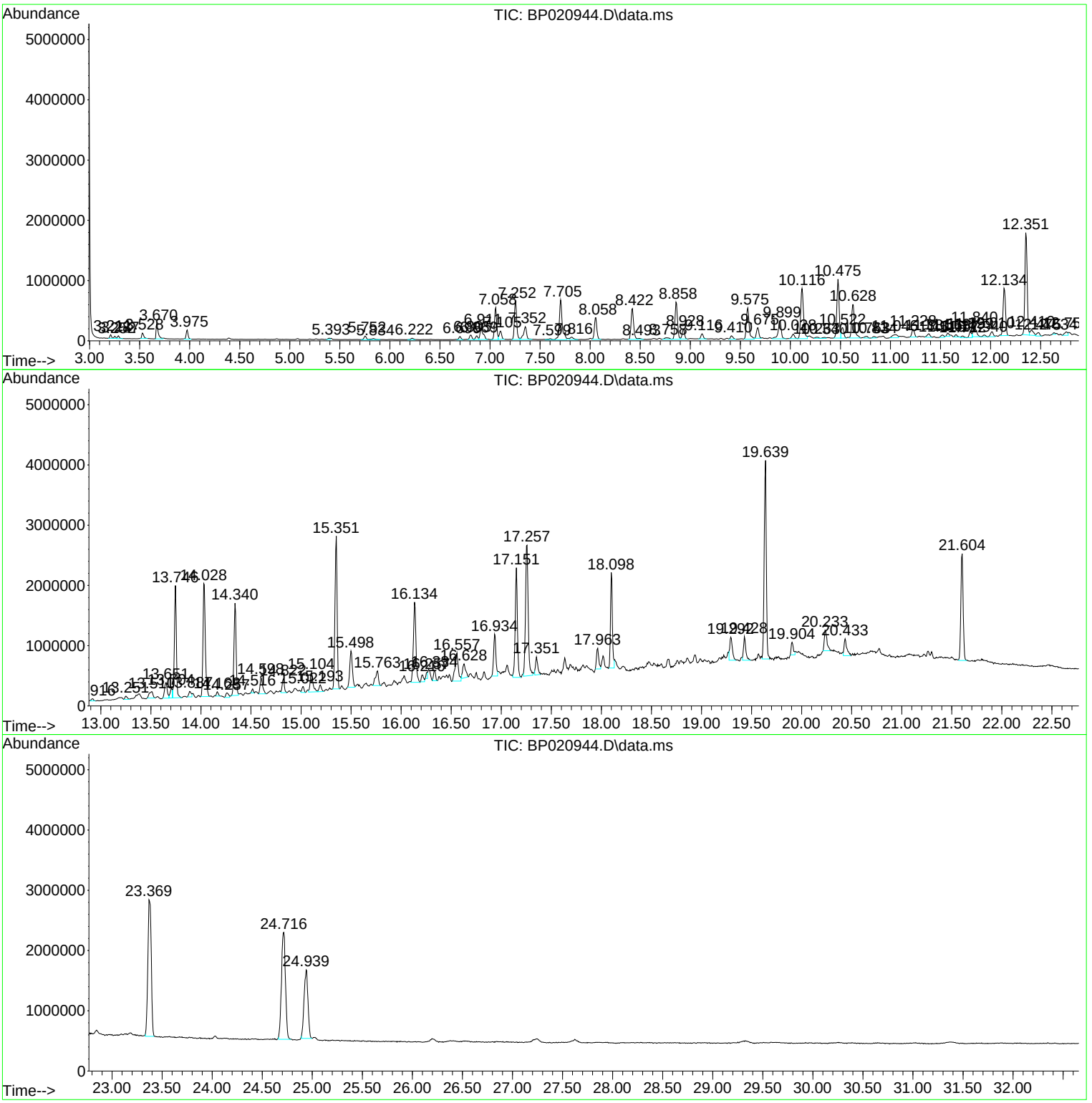
Sum of corrected areas: 77576979

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

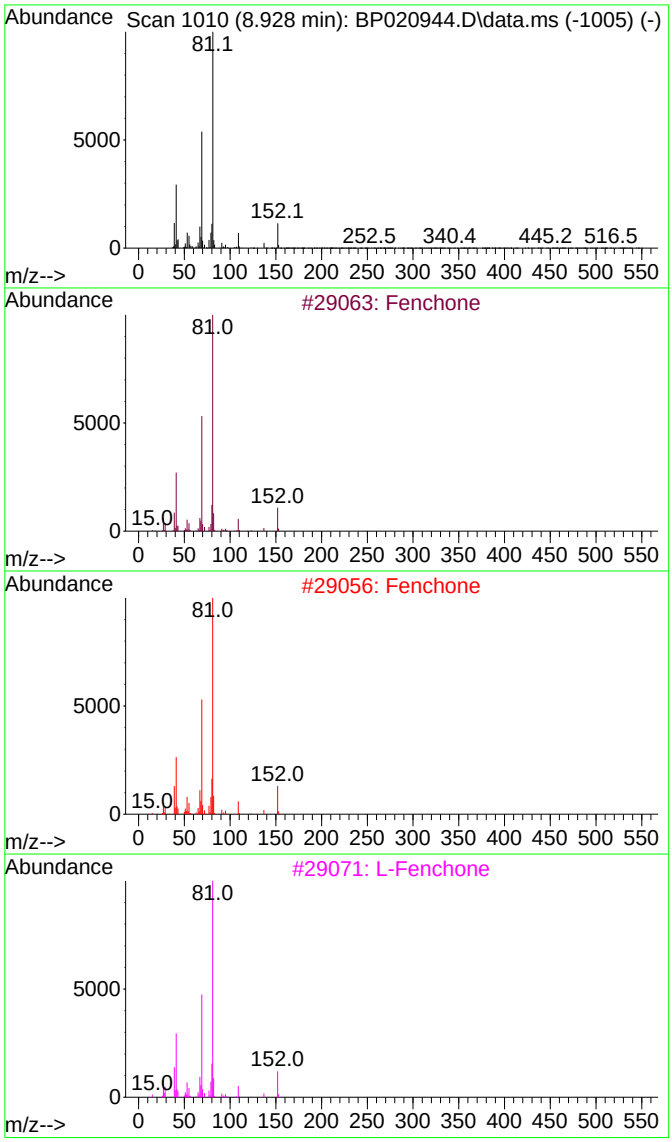
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

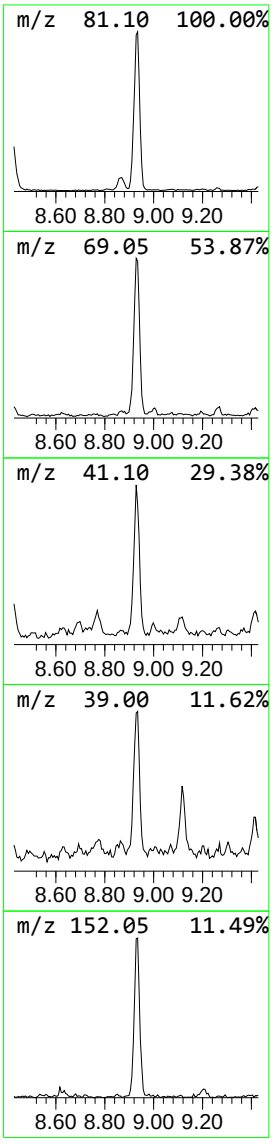
****

Peak Number 4 Fenchone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	4.08 ng/ul	268872	1,4-Dichlorobenzene-d4	7.705

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fenchone	152	C10H16O	001195-79-5	94
2	Fenchone	152	C10H16O	001195-79-5	91
3	L-Fenchone	152	C10H16O	007787-20-4	90
4	L-Fenchone	152	C10H16O	007787-20-4	90
5	D-Fenchone	152	C10H16O	004695-62-9	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

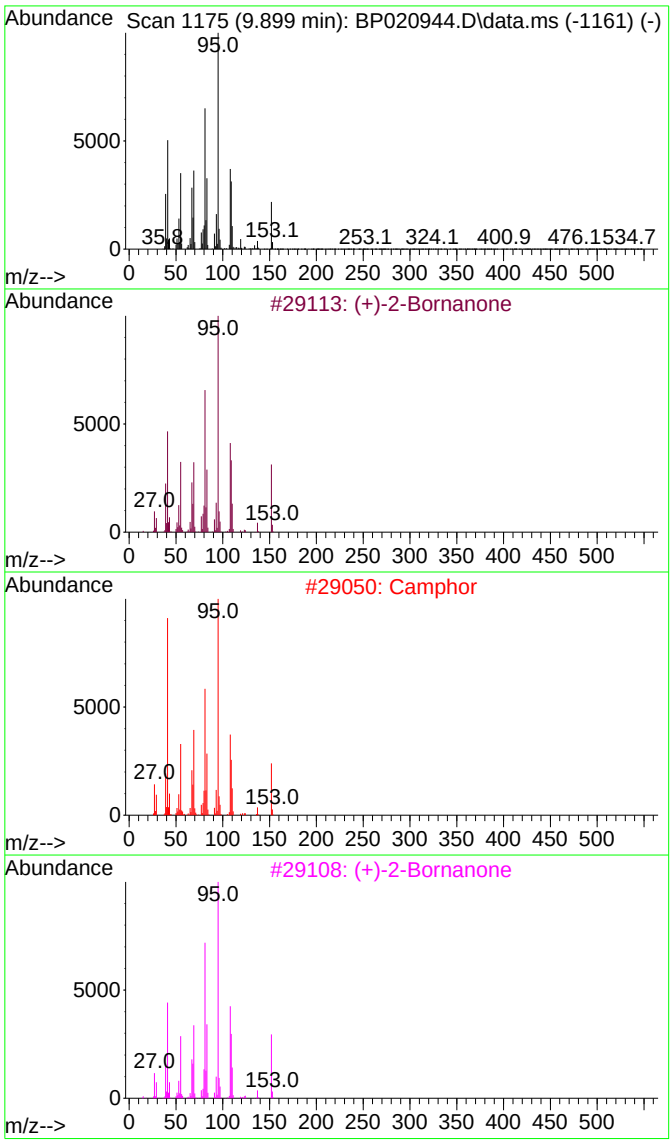
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

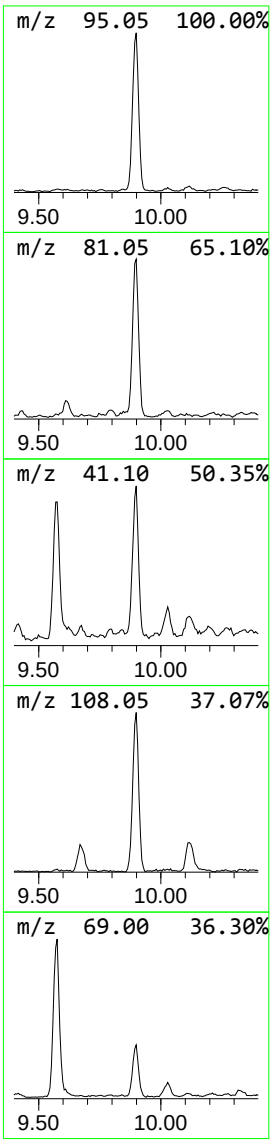
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (+)-2-Bornanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.899	8.21 ng/ul	640627	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
2	Camphor	152	C10H16O	000076-22-2	96
3	(+)-2-Bornanone	152	C10H16O	000464-49-3	96
4	(+)-2-Bornanone	152	C10H16O	000464-49-3	96
5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

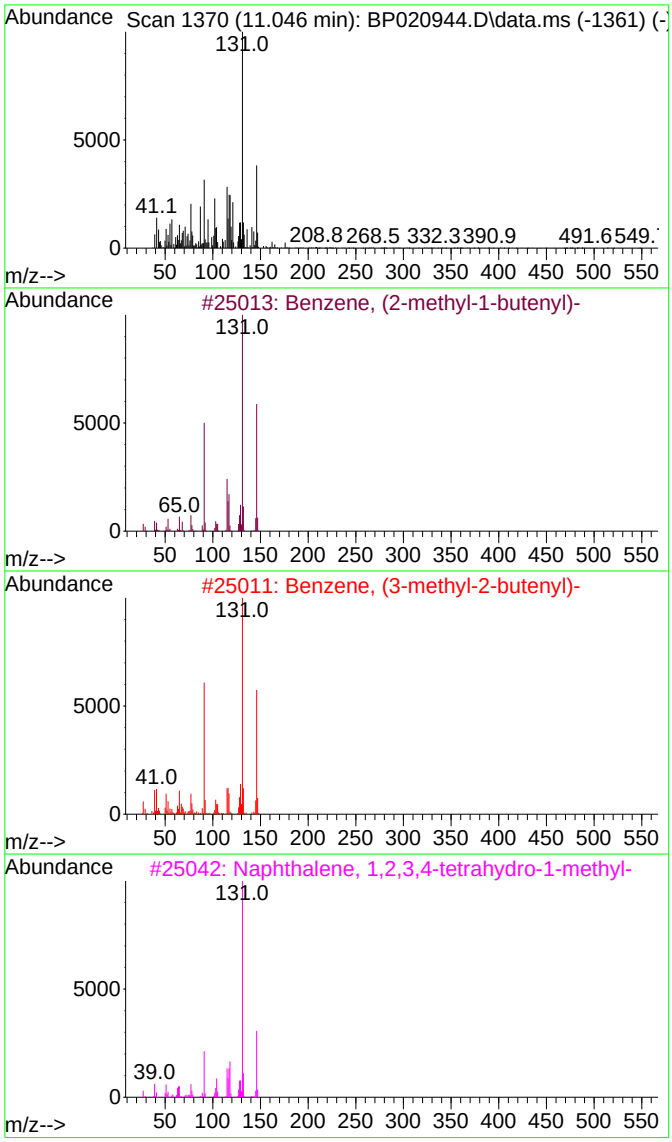
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

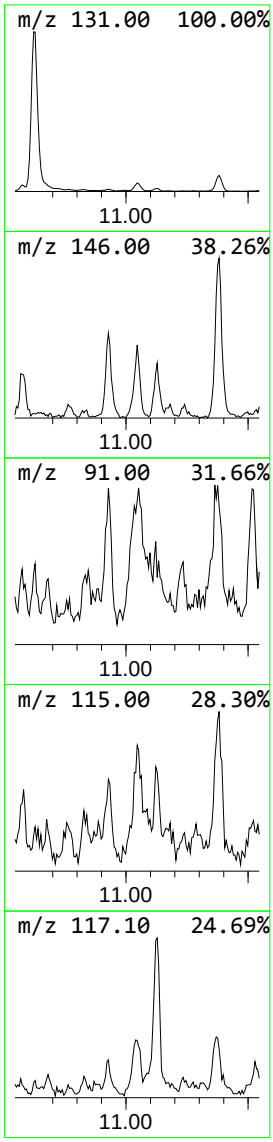
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (2-methyl-1-butenyl)- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.046	2.39 ng/ul	186780	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	89
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	89
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

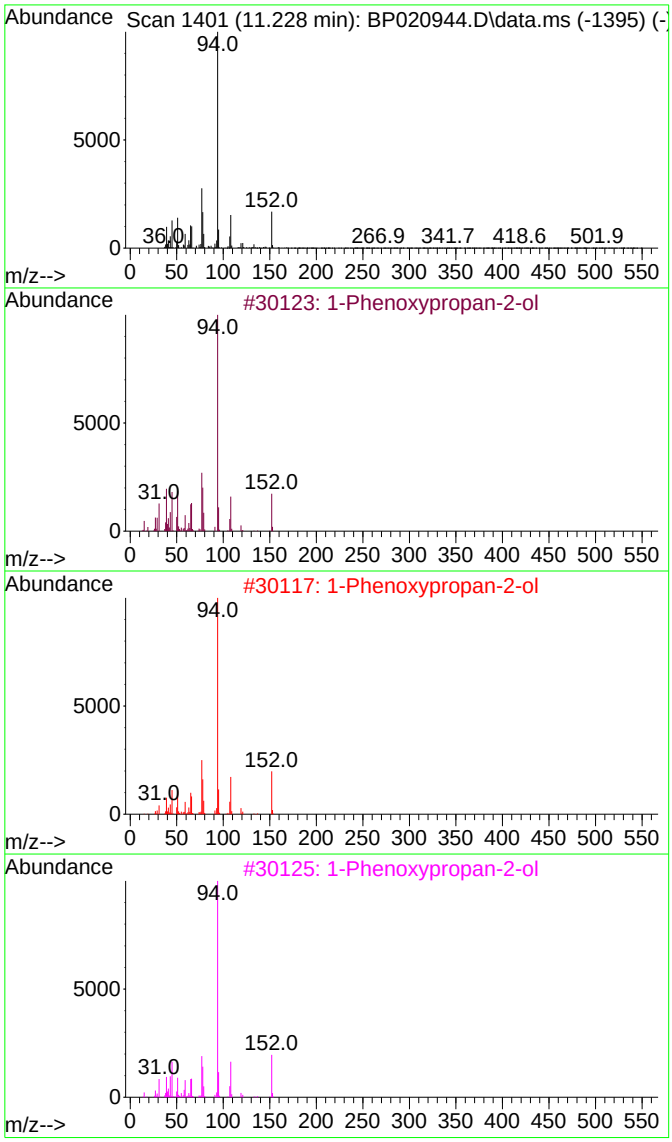
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

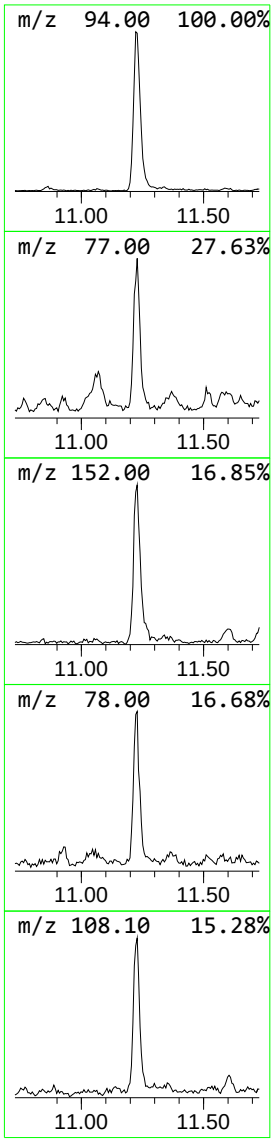
****

Peak Number 7 1-Phenoxypropan-2-ol Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.228	3.04 ng/ul	237235	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5	Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

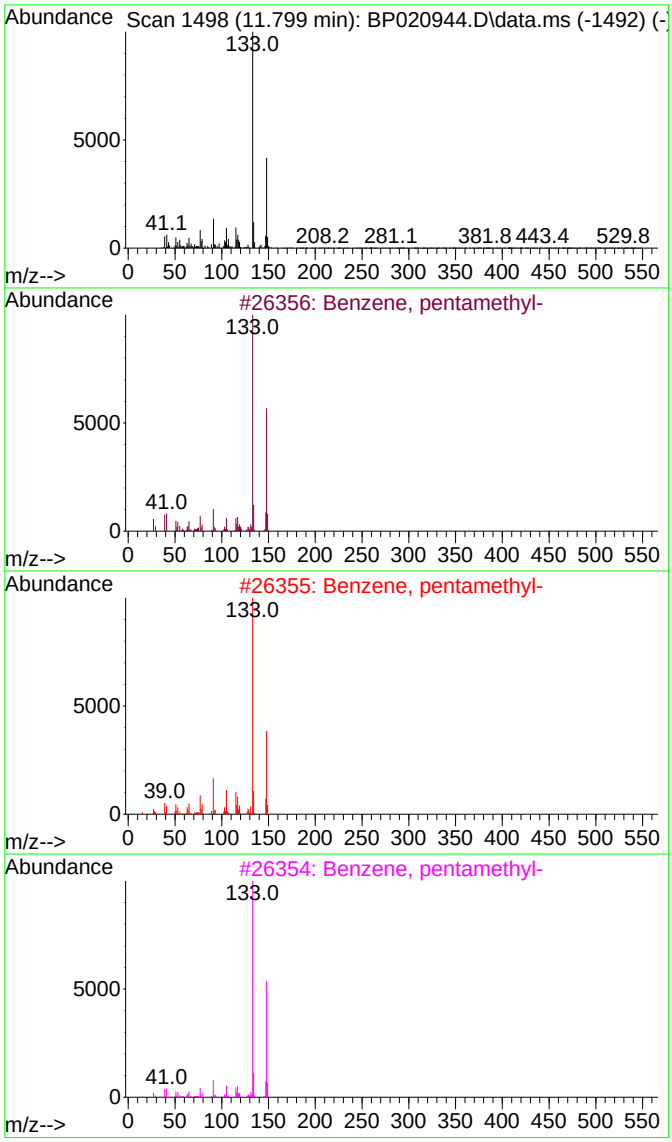
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

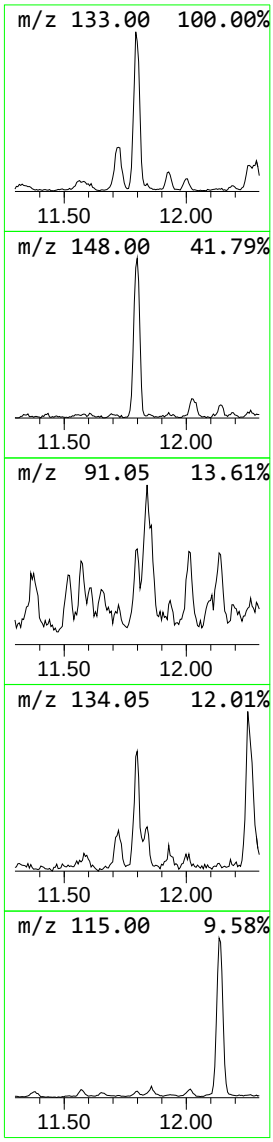
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, pentamethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.799	2.17 ng/ul	169301	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	91
2	Benzene, pentamethyl-	148	C11H16	000700-12-9	91
3	Benzene, pentamethyl-	148	C11H16	000700-12-9	91
4	3,4-Dimethylcumene	148	C11H16	1000370-34-1	91
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	90





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Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

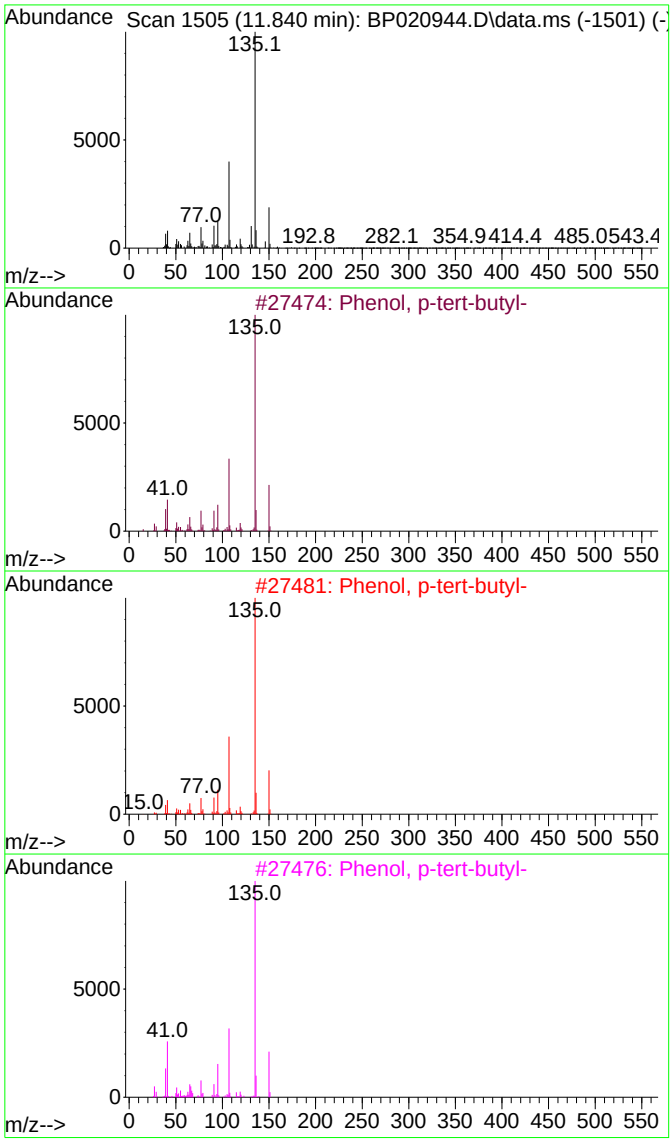
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

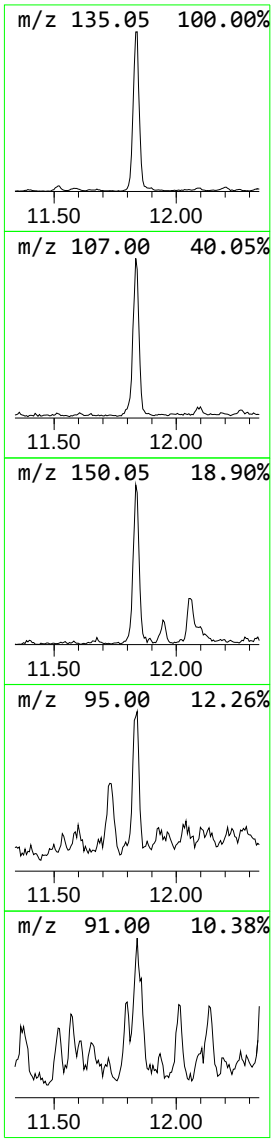
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, p-tert-butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.840	4.95 ng/ul	386136	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	95
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	81
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	81
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	76





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Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

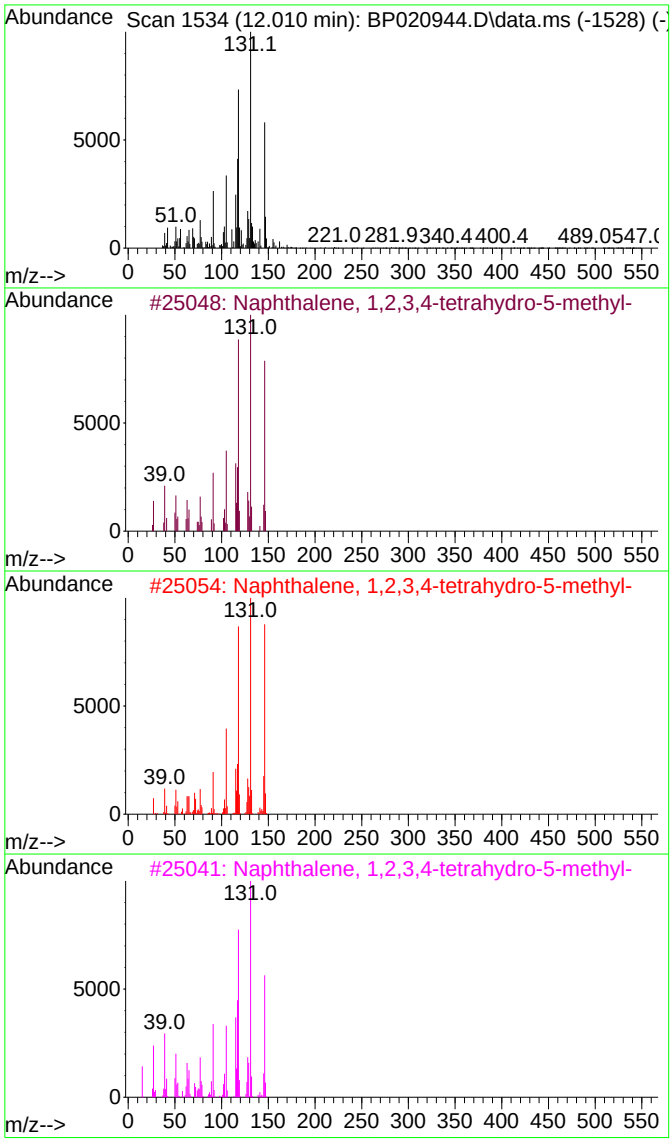
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

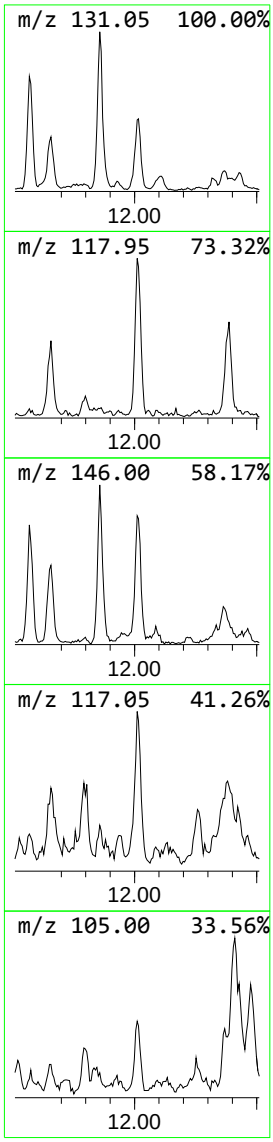
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.010	2.06 ng/ul	160720	Naphthalene-d8	10.475

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	91
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	90
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	89
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	83





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Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

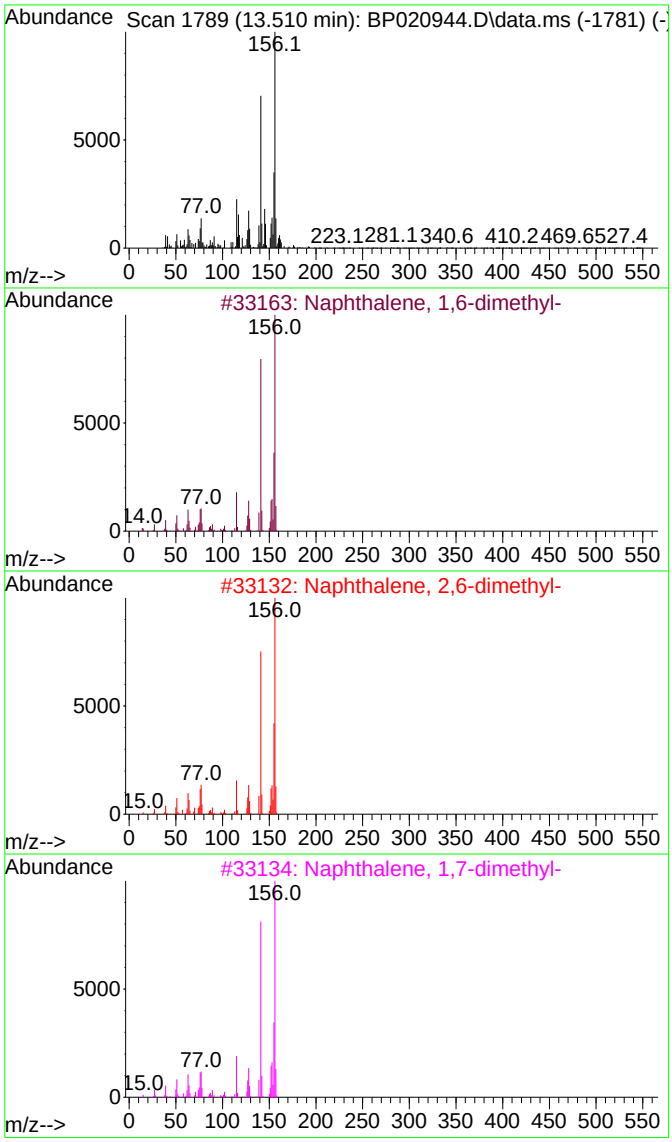
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

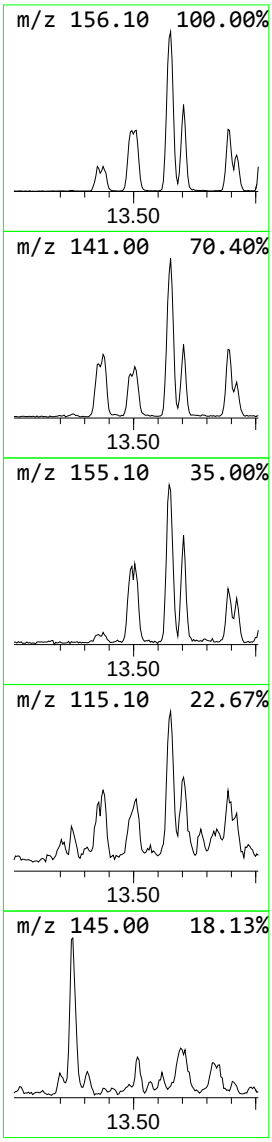
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Naphthalene, 1,6-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.510	2.46 ng/ul	276264	Acenaphthene-d10	14.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96





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Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

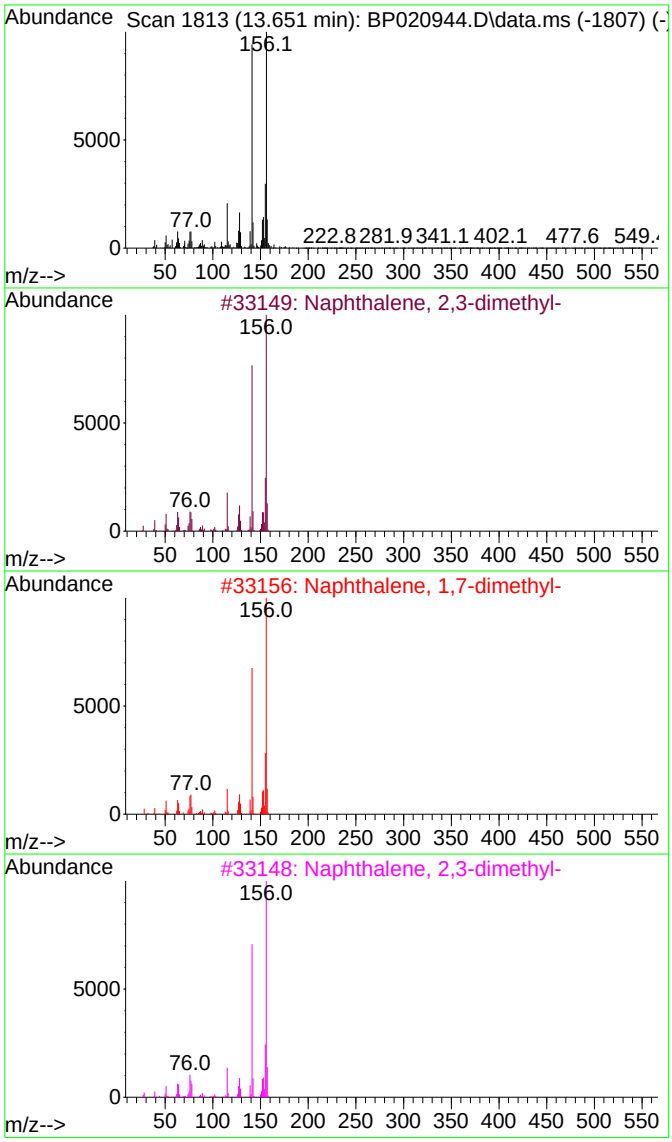
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

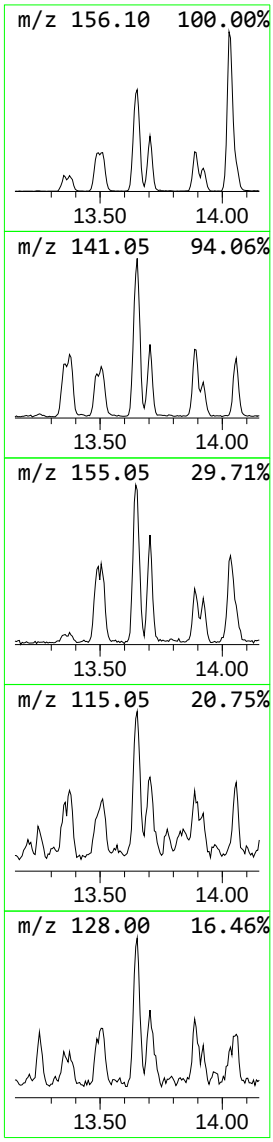
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Naphthalene, 2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.651	4.40 ng/ul	494517	Acenaphthene-d10	14.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97





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Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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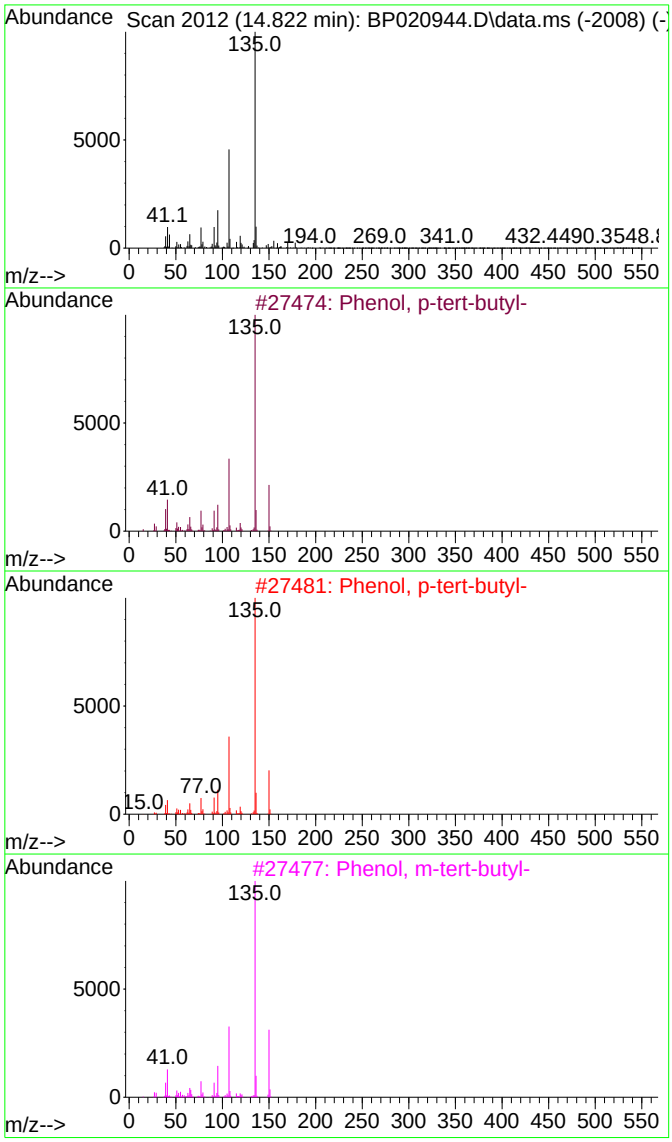
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

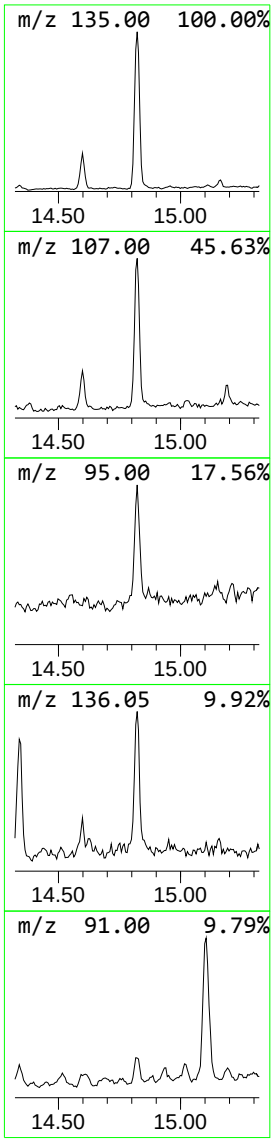
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenol, m-tert-butyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.822	3.07 ng/ul	345556	Acenaphthene-d10	14.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	90
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	83
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	83
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	83
5	4,4'-(1,2-diethylethylene)diphenol	270	C18H22O2	005635-50-7	80





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Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

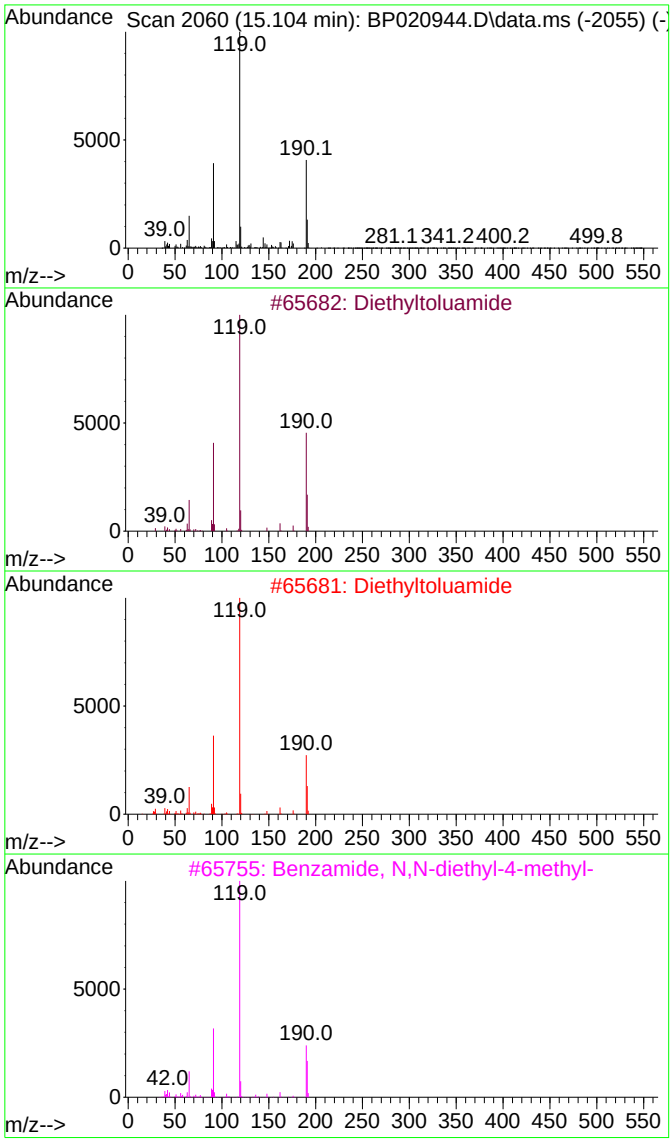
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

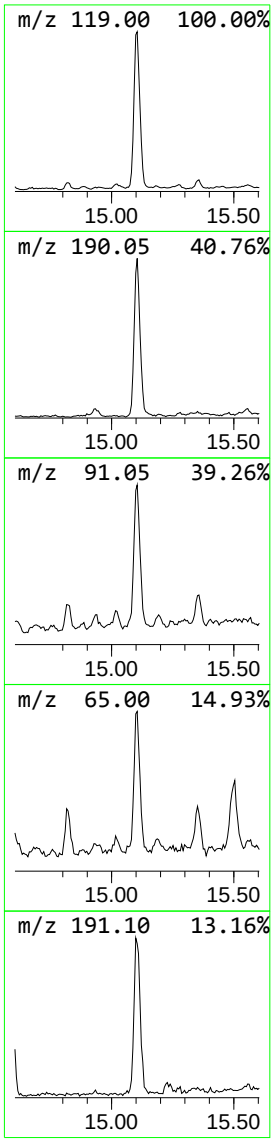
****

Peak Number 14 Diethyltoluamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.104	5.81 ng/ul	653343	Acenaphthene-d10	14.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	94
3	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	76
4	Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	70
5	Benzamide, 4-methyl-N-butyl-	191	C12H17NO	1000407-46-6	70





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Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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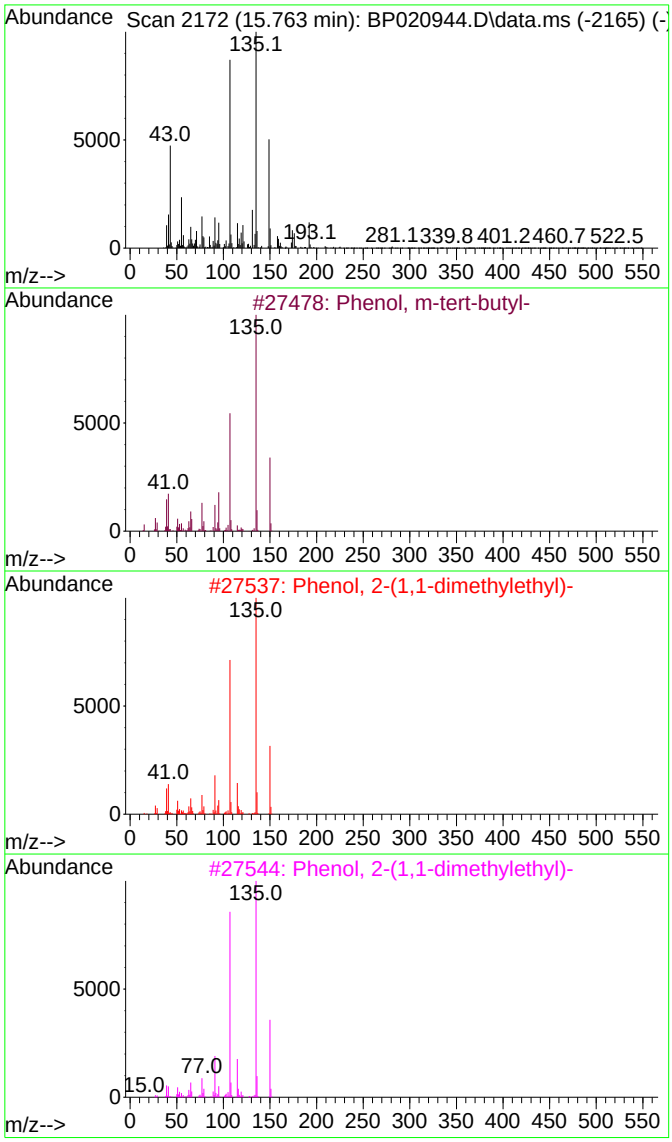
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

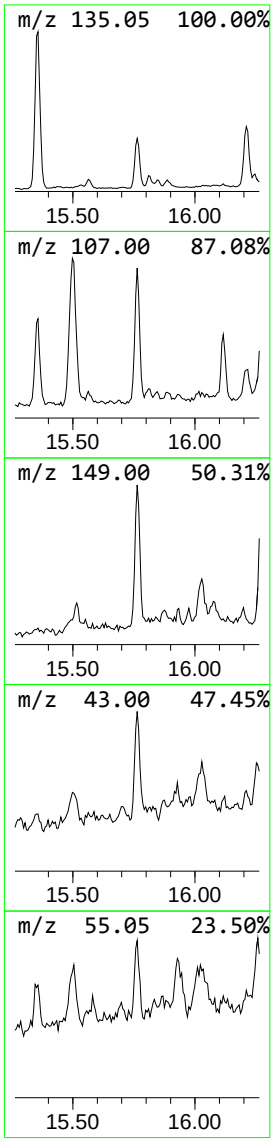
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Phenol, 2-(1,1-dimethylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.763	3.60 ng/ul	498122	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	58
2	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	58
3	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	50
4	1,5,6,7-Tetrahydro-4-indolone	135	C8H9NO	013754-86-4	49
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38





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Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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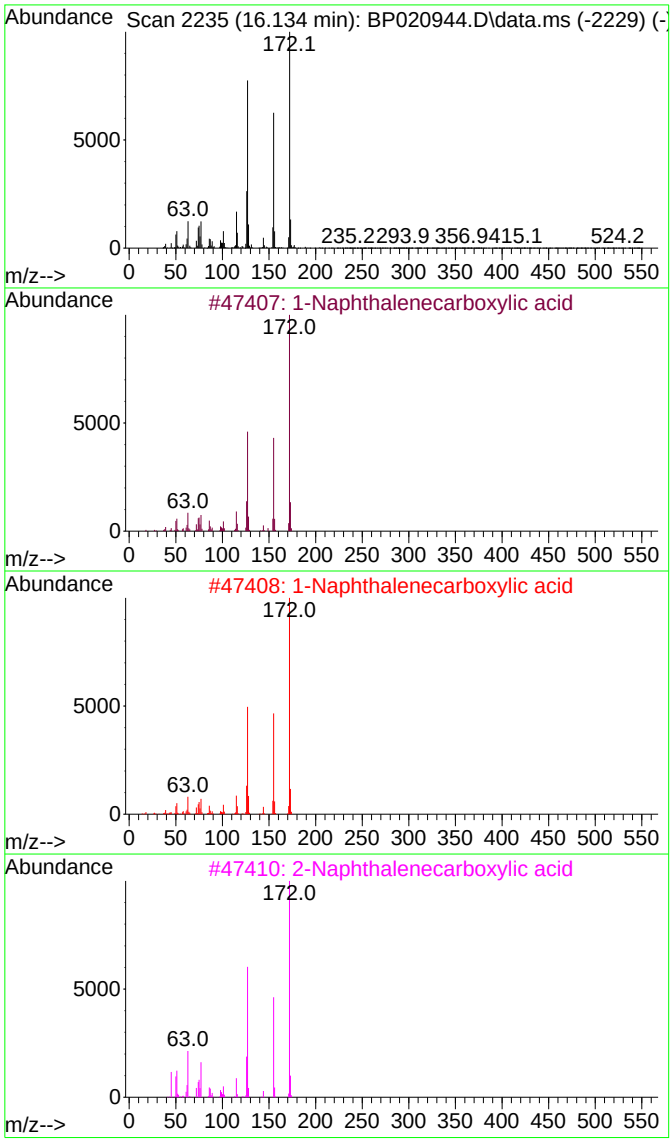
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

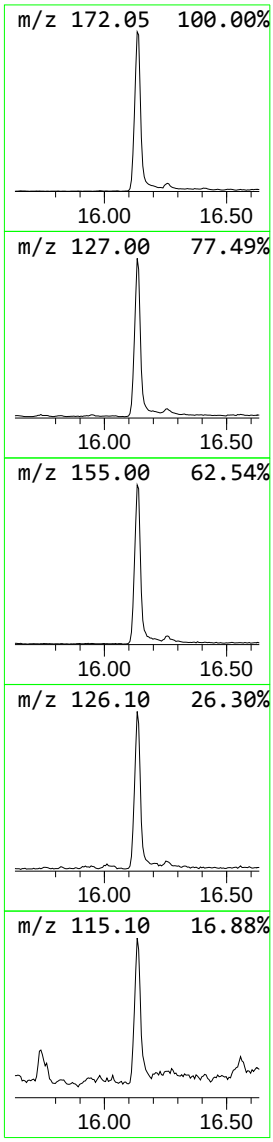
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1-Naphthalenecarboxylic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.134	16.61 ng/ul	2300040	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	95
2	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	95
3	2-Naphthalenecarboxylic acid	172	C11H8O2	000093-09-4	94
4	1-Naphthalenecarboxylic acid	172	C11H8O2	000086-55-5	93
5	2-Naphthalenecarboxylic acid	172	C11H8O2	000093-09-4	93





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Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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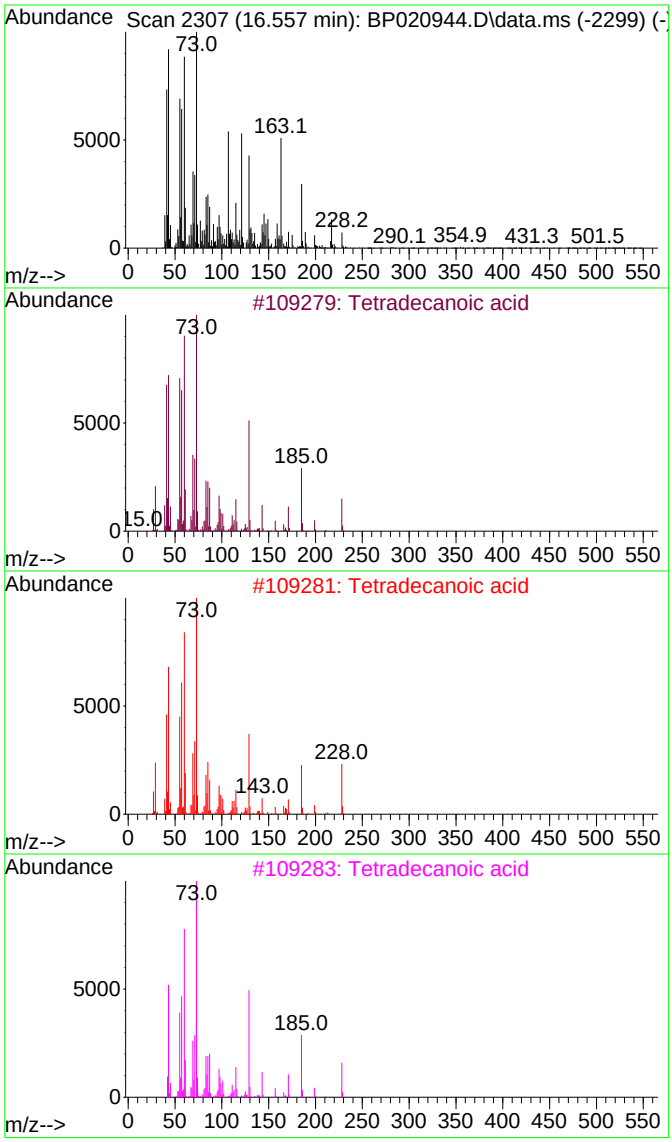
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

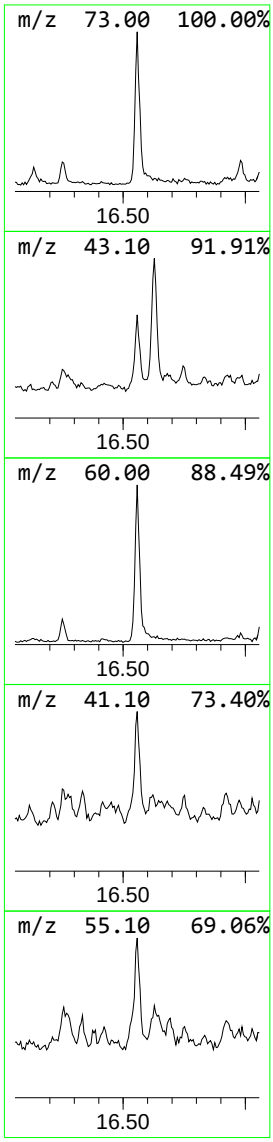
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Tetradecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.557	7.35 ng/ul	1017710	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	60
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
5	Tridecanoic acid	214	C13H26O2	000638-53-9	53





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Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

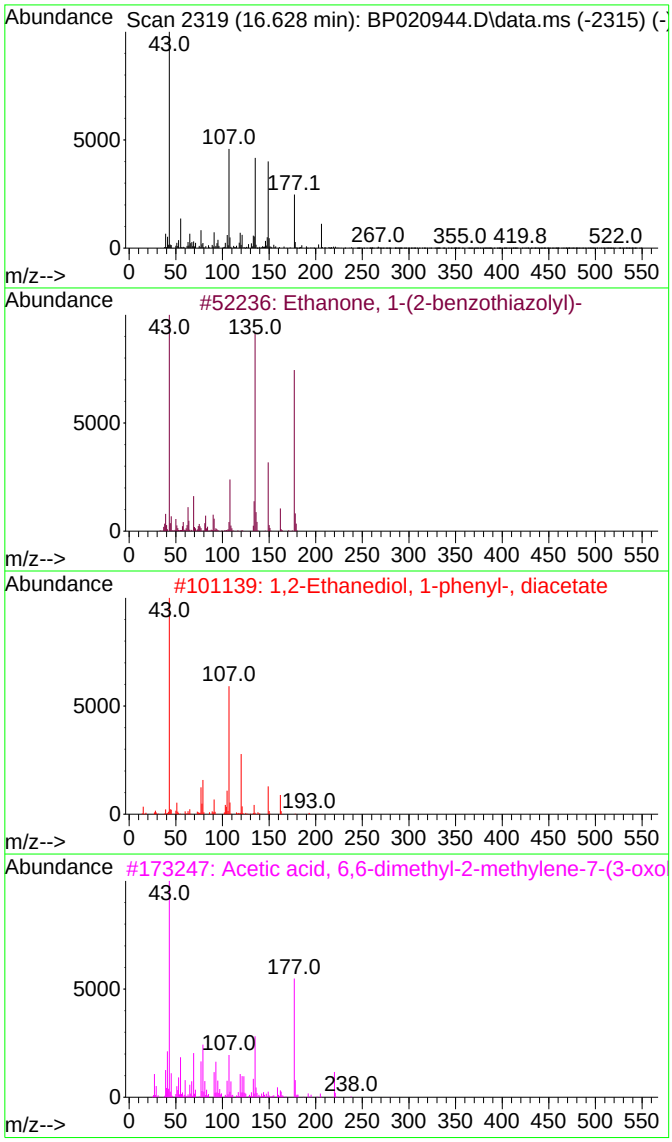
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

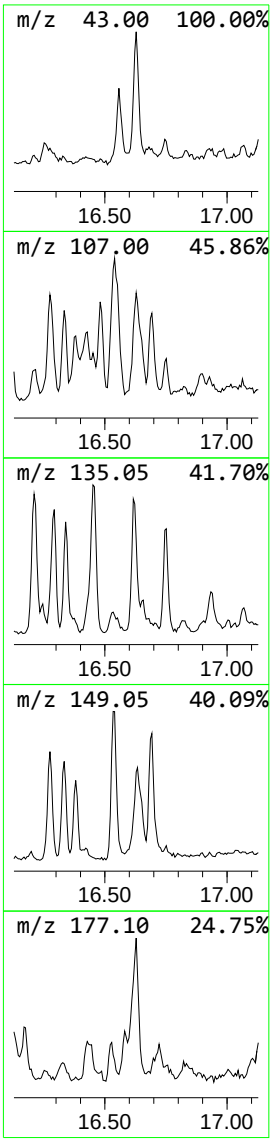
****

Peak Number 18 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.628	3.49 ng/ul	483801	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(2-benzothiazolyl)-	177	C9H7NOS	001629-78-3	25
2	1,2-Ethanediol, 1-phenyl-, diace...	222	C12H14O4	006270-03-7	12
3	Acetic acid, 6,6-dimethyl-2-meth...	280	C16H24O4	1000186-15-5	11
4	Benzenemethanol, .alpha.,.alpha....	150	C10H14O	001197-01-9	9
5	Benzenemethanol, .alpha.,.alpha....	150	C10H14O	001197-01-9	9





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Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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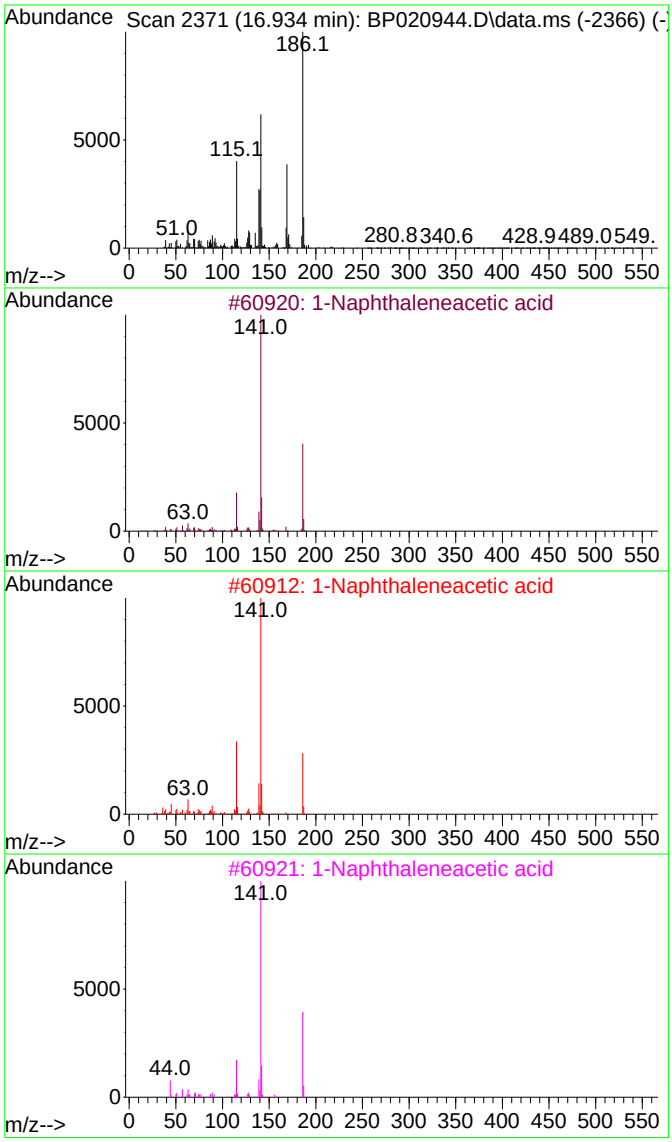
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

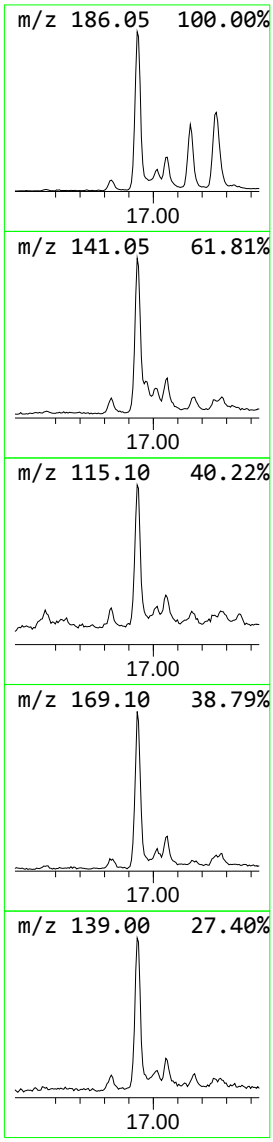
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 1-Naphthaleneacetic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.934	8.05 ng/ul	1114970	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthaleneacetic acid	186	C12H10O2	000086-87-3	60
2	1-Naphthaleneacetic acid	186	C12H10O2	000086-87-3	52
3	1-Naphthaleneacetic acid	186	C12H10O2	000086-87-3	50
4	2-Naphthaleneacetic acid	186	C12H10O2	000581-96-4	49
5	1-Naphthaleneacetic acid	186	C12H10O2	000086-87-3	49





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Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

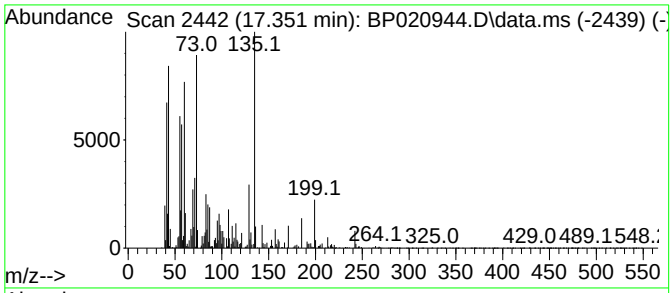
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

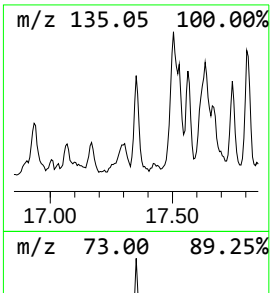
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

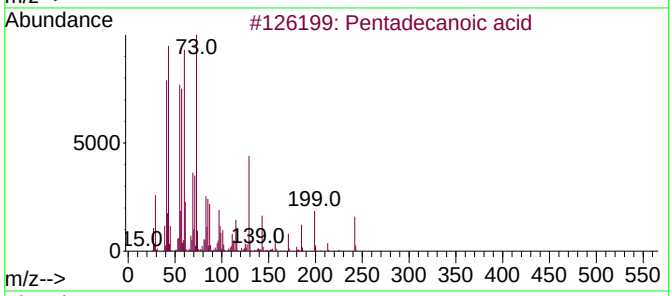
****
Peak Number 20 Pentadecanoic acid Concentration Rank 20

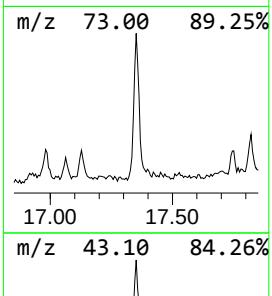
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.351	3.31 ng/ul	457868	Phenanthrene-d10	17.151

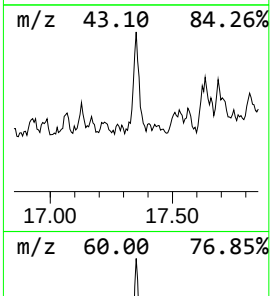
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	97
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	96
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	55
5	Dodecanoic acid	200	C12H24O2	000143-07-7	55

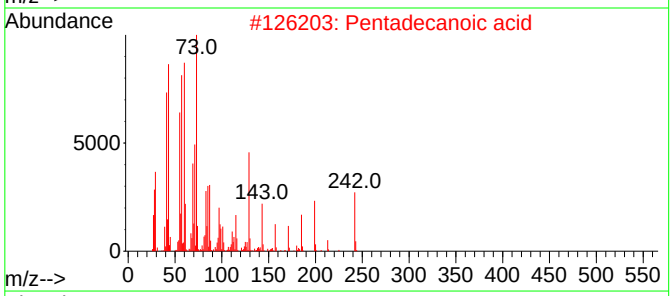


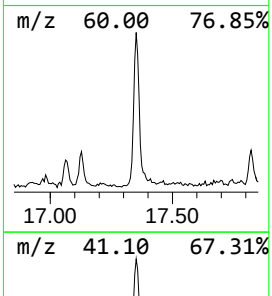


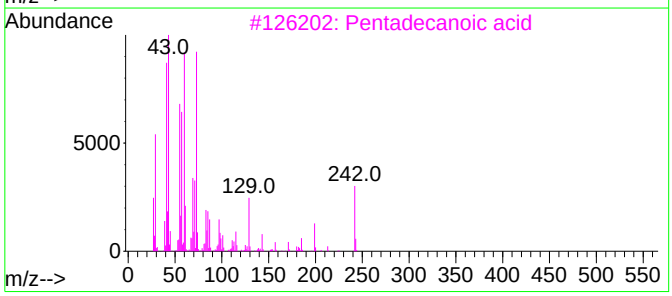


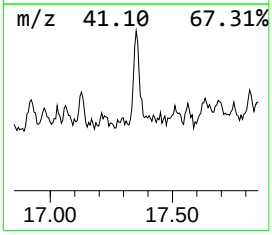












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

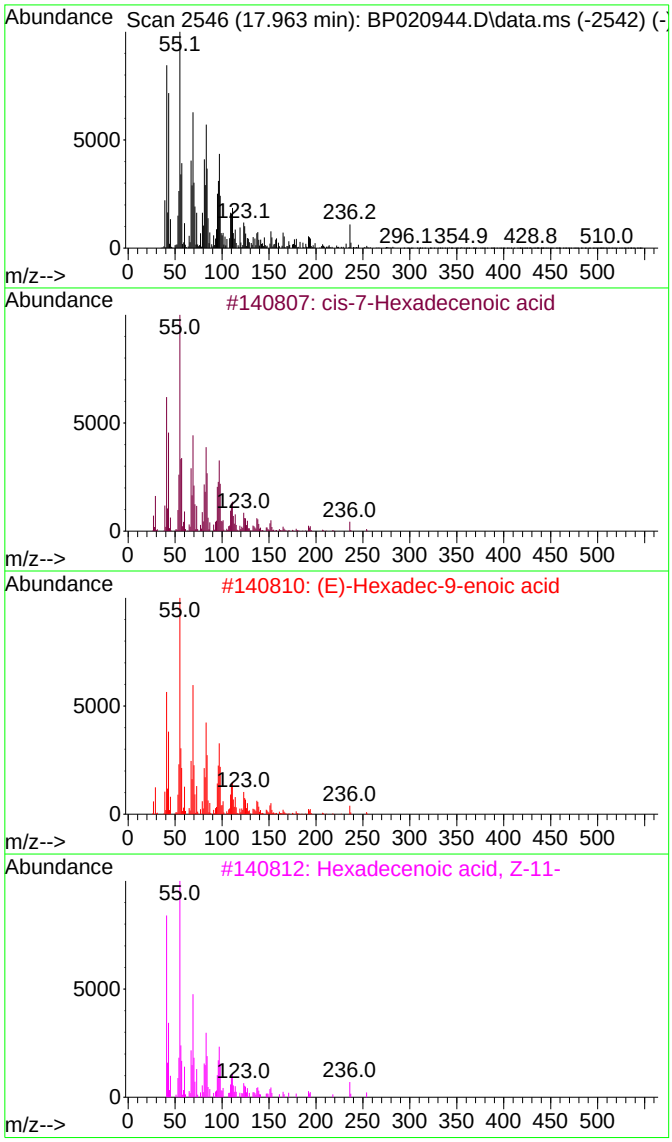
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

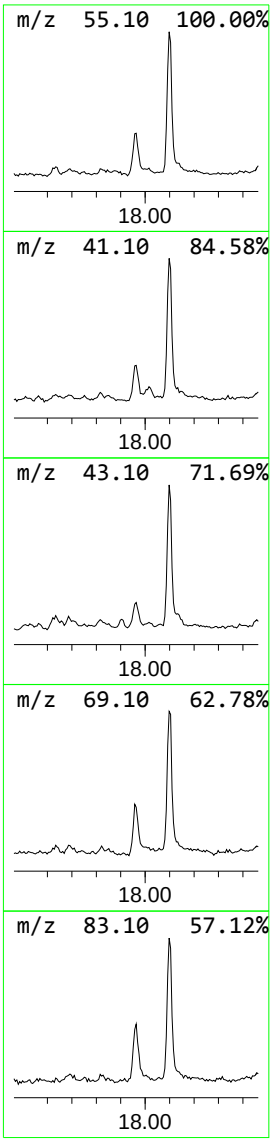
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 cis-7-Hexadecenoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.963	3.86 ng/ul	534118	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	99
2	(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
3	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	94
4	Palmitoleic acid	254	C16H30O2	000373-49-9	90
5	cis-Vaccenic acid	282	C18H34O2	000506-17-2	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

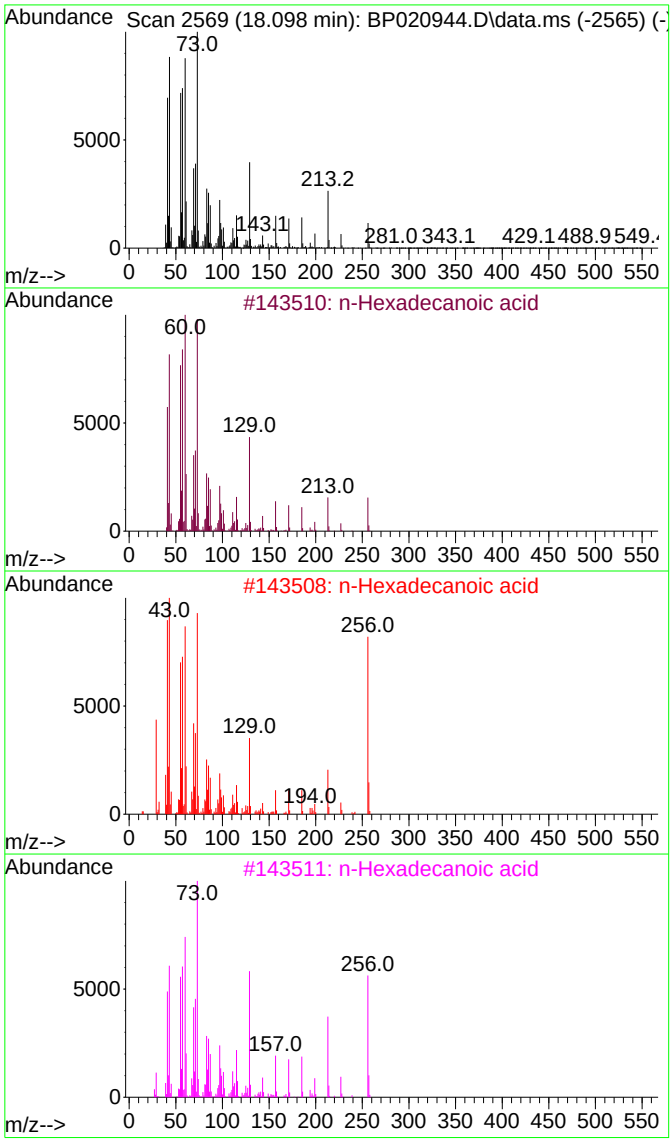
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

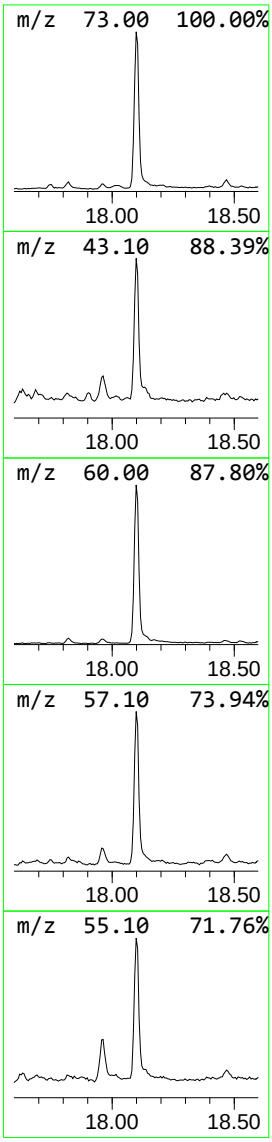
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.098	15.32 ng/ul	2121490	Phenanthrene-d10	17.151

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	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	93





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Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

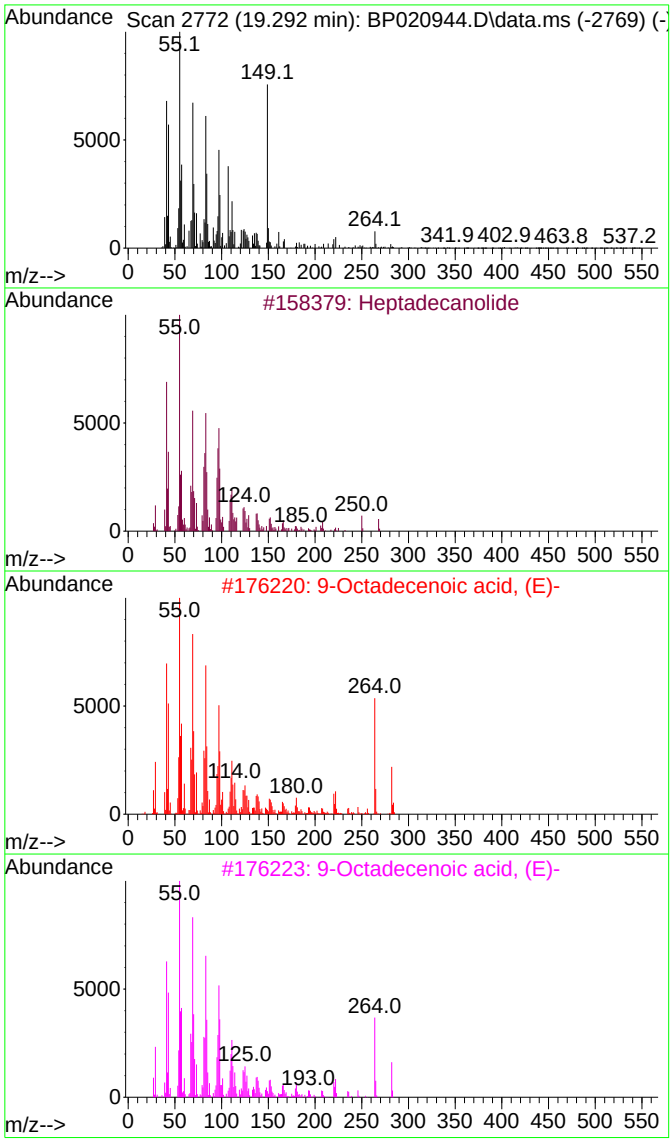
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

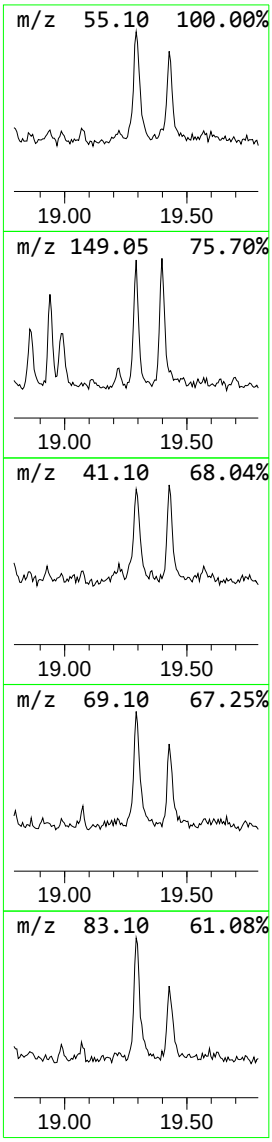
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Heptadecanolide Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.292	4.87 ng/ul	674480	Phenanthrene-d10	17.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecanolide	268	C17H32O2	005637-97-8	55
2	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	49
3	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	46
4	Cyclohexadecane	224	C16H32	000295-65-8	38
5	1-Hexadecanol	242	C16H34O	036653-82-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
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Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

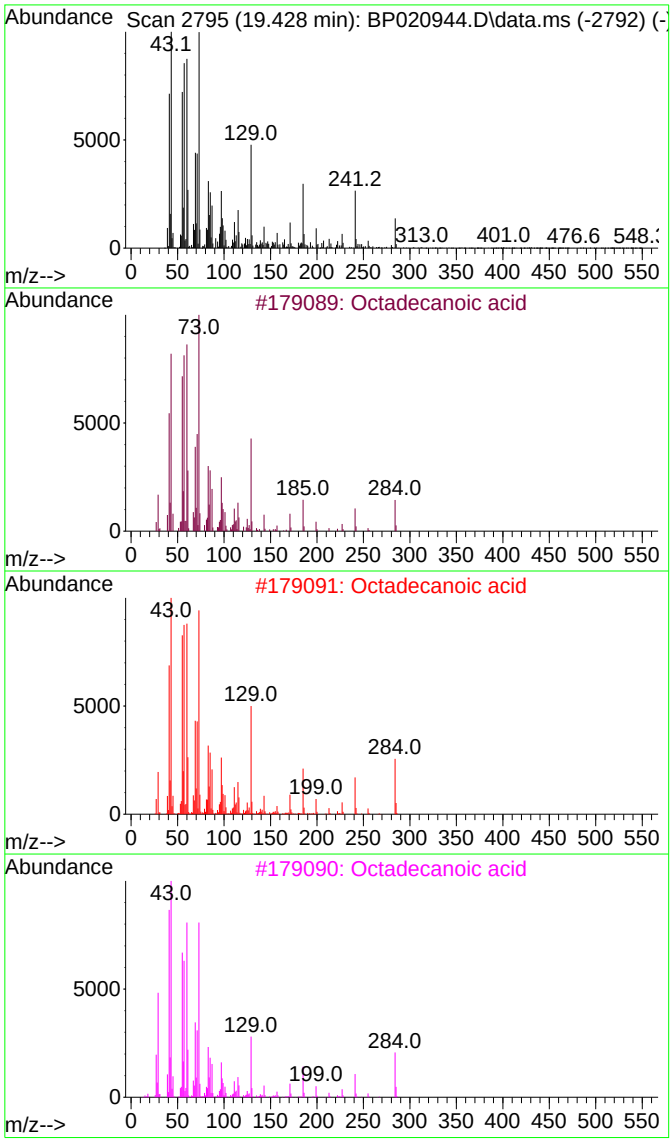
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

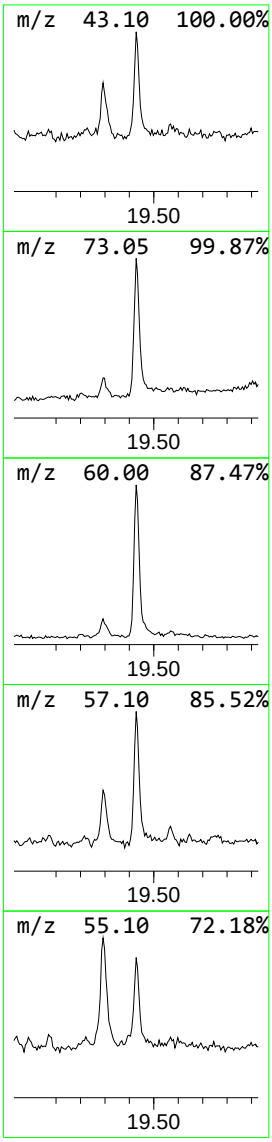
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.428	4.26 ng/ul	608167	Chrysene-d12	21.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	95
4	Octadecanoic acid	284	C18H36O2	000057-11-4	94
5	Octadecanoic acid	284	C18H36O2	000057-11-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

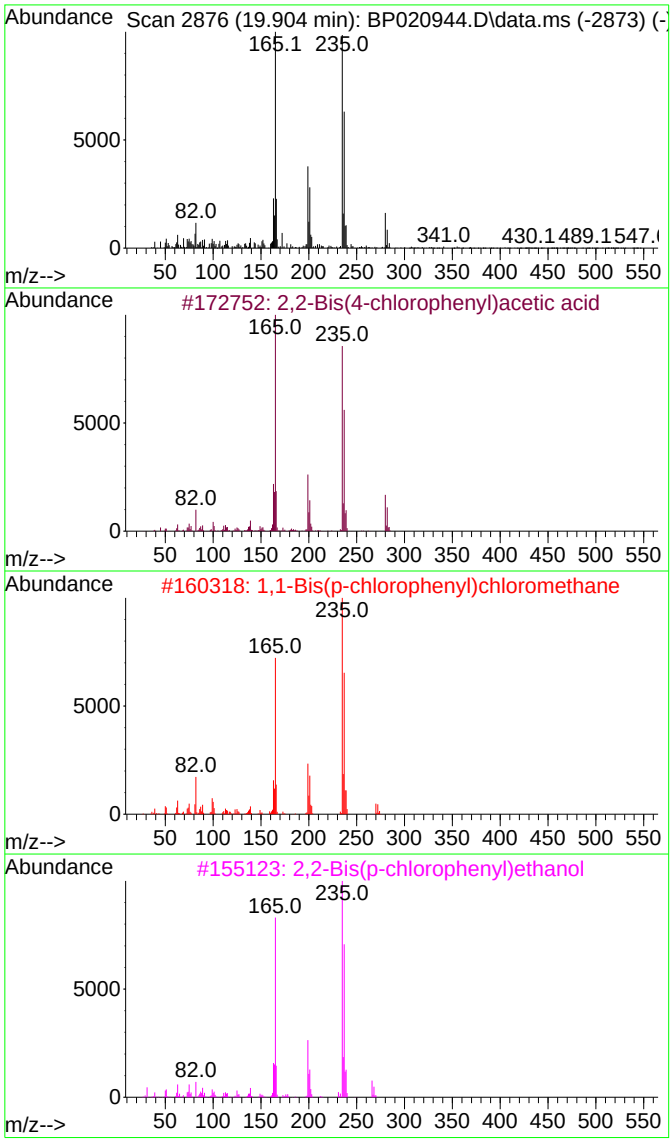
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

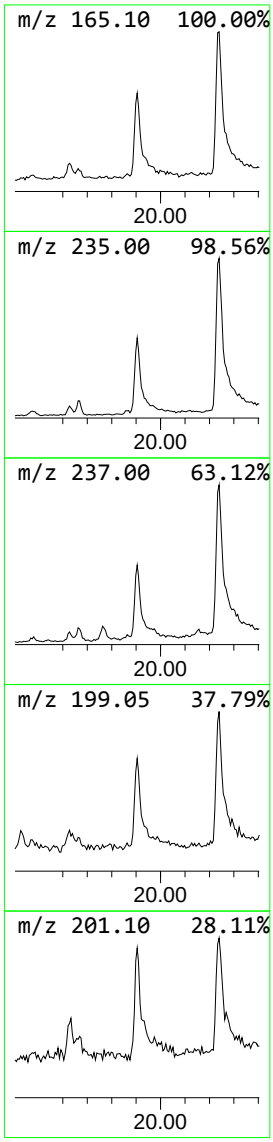
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 2,2-Bis(4-chlorophenyl)acet... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.904	2.34 ng/ul	334223	Chrysene-d12	21.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Bis(4-chlorophenyl)acetic acid	280	C14H10Cl2O2	000083-05-6	94
2	1,1-Bis(p-chlorophenyl)chloromet...	270	C13H9Cl3	000782-08-1	91
3	2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	72
4	m,p'-DDD	318	C14H10Cl4	004329-12-8	72
5	3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	72





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Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

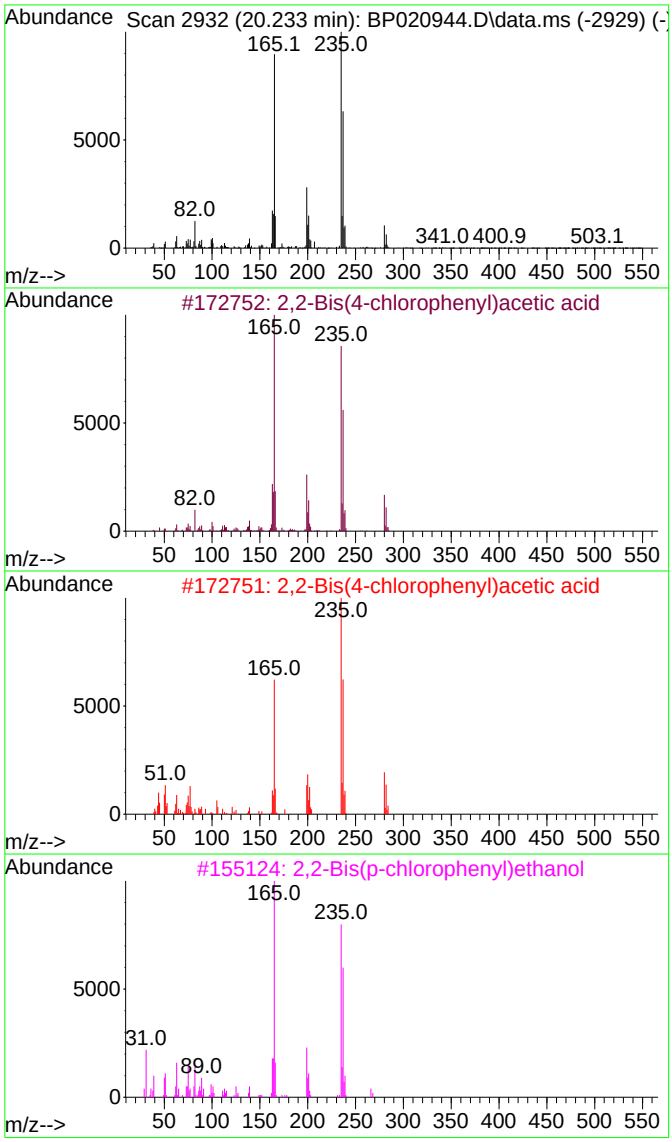
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

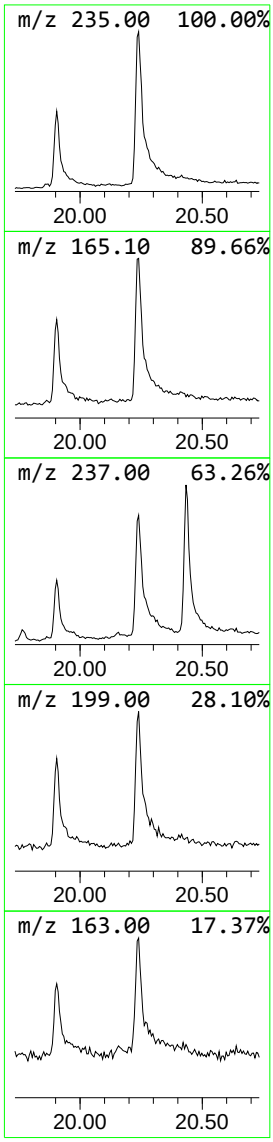
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 2,2-Bis(p-chlorophenyl)ethanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.233	4.40 ng/ul	627106	Chrysene-d12	21.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Bis(4-chlorophenyl)acetic acid	280	C14H10Cl2O2	000083-05-6	96
2	2,2-Bis(4-chlorophenyl)acetic acid	280	C14H10Cl2O2	000083-05-6	94
3	2,2-Bis(p-chlorophenyl)ethanol	266	C14H12Cl2O	002642-82-2	90
4	Mitotane	318	C14H10Cl4	000053-19-0	78
5	1-Chloro-2,2-Bis(p-chlorophenyl)...	284	C14H11Cl3	002642-80-0	78





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Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

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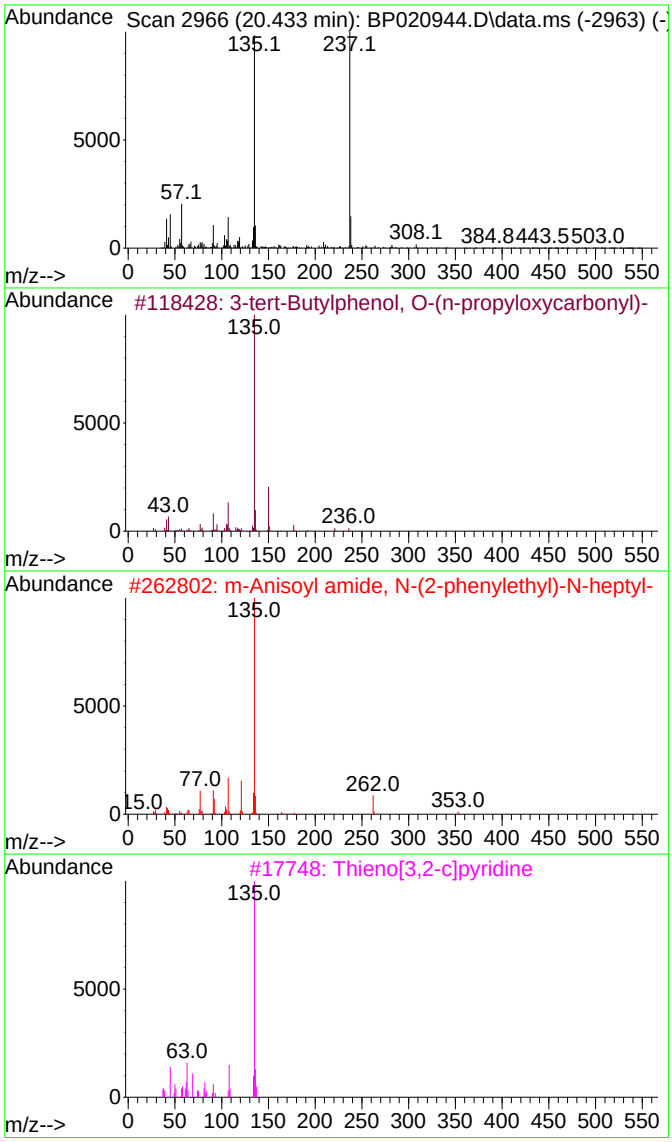
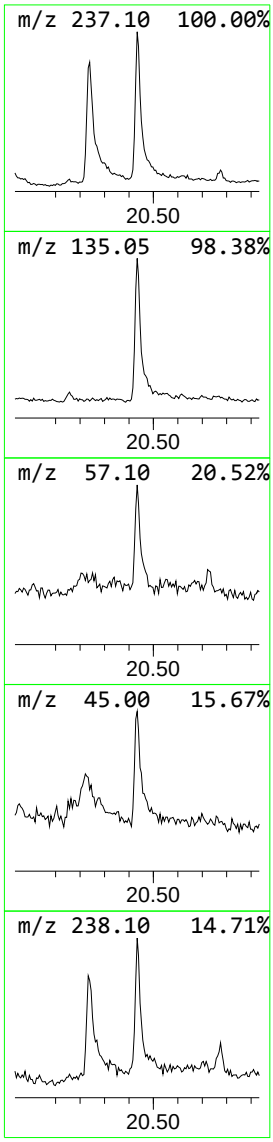
Peak Number 27 unknown-02 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.433	3.65 ng/ul	520790	Chrysene-d12	21.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-tert-Butylphenol, O-(n-propylo...	236	C14H20O3	1000487-38-2	41
2	m-Anisoyl amide, N-(2-phenylethy...	353	C23H31NO2	1000451-74-5	38
3	Thieno[3,2-c]pyridine	135	C7H5NS	000272-14-0	35
4	Trichloroacetic acid, 1-adamanty...	310	C13H17Cl3O2	1000282-92-0	35
5	1-Adamantanecarboxylic acid, 4-c...	290	C17H19ClO2	1000307-66-0	35





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Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

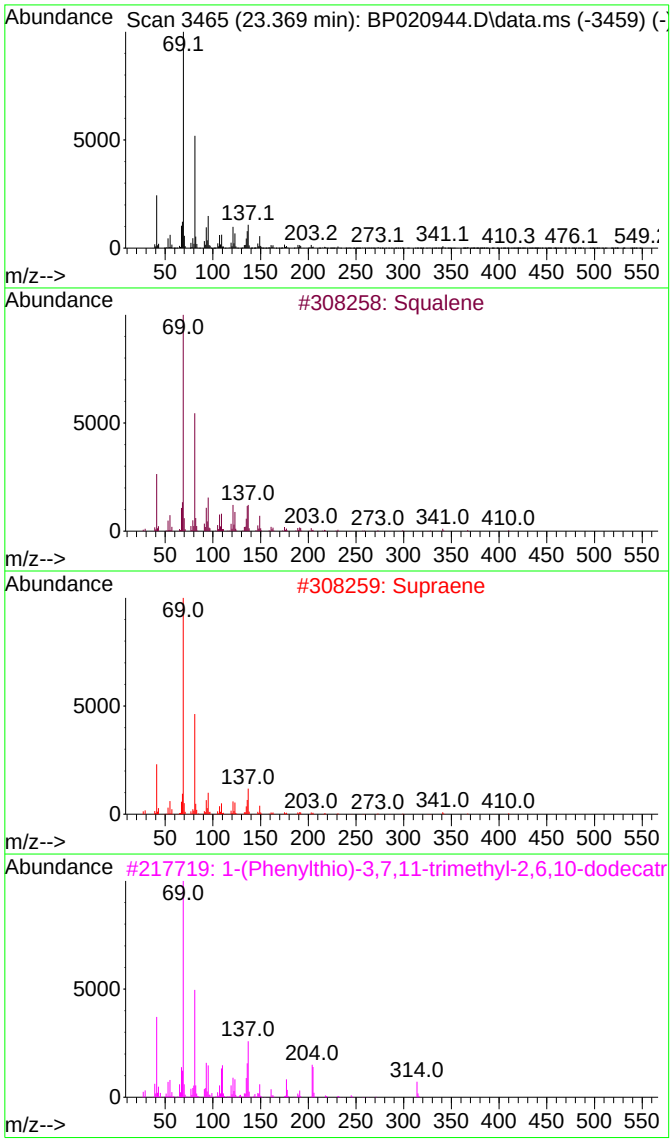
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

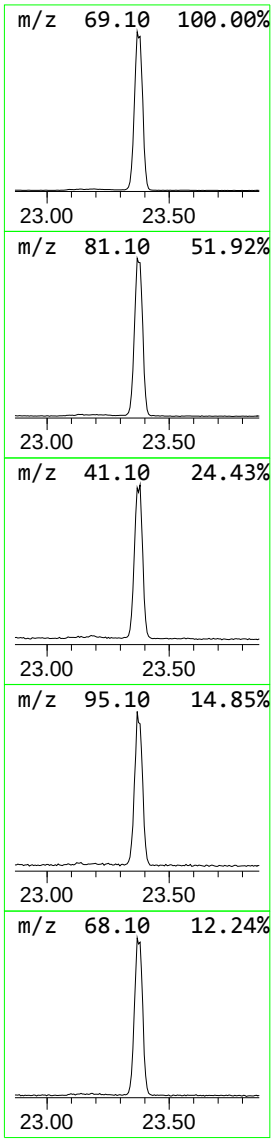
****

Peak Number 28 Squalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.369	34.46 ng/ul	5249190	Perylene-d12	24.939

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	95
2	Supraene	410	C30H50	007683-64-9	91
3	1-(Phenylthio)-3,7,11-trimethyl-...	314	C21H30S	028413-57-2	90
4	Umbelliprenin	366	C24H30O3	023838-17-7	78
5	2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020944.D
Acq On : 12 Jul 2024 22:23
Operator : MA/JU
Sample : P3217-14
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
Fenchone	8.928	4.1	ng/ul	268872	1	7.705	1316870	20.0
(+)-2-Bornanone	9.899	8.2	ng/ul	640627	2	10.475	1560410	20.0
Benzene, (2-met...	11.046	2.4	ng/ul	186780	2	10.475	1560410	20.0
1-Phenoxypropan...	11.228	3.0	ng/ul	237235	2	10.475	1560410	20.0
Benzene, pentam...	11.799	2.2	ng/ul	169301	2	10.475	1560410	20.0
Phenol, p-tert-...	11.840	5.0	ng/ul	386136	2	10.475	1560410	20.0
Naphthalene, 1,...	12.010	2.1	ng/ul	160720	2	10.475	1560410	20.0
Naphthalene, 1,...	13.510	2.5	ng/ul	276264	3	14.340	2250220	20.0
Naphthalene, 2,...	13.651	4.4	ng/ul	494517	3	14.340	2250220	20.0
Phenol, m-tert-...	14.822	3.1	ng/ul	345556	3	14.340	2250220	20.0
Diethyltoluamide	15.104	5.8	ng/ul	653343	3	14.340	2250220	20.0
Phenol, 2-(1,1-...	15.763	3.6	ng/ul	498122	4	17.151	2769400	20.0
1-Naphthaleneca...	16.134	16.6	ng/ul	2300040	4	17.151	2769400	20.0
Tetradecanoic acid	16.557	7.3	ng/ul	1017710	4	17.151	2769400	20.0
unknown-01	16.628	3.5	ng/ul	483801	4	17.151	2769400	20.0
1-Naphthaleneac...	16.934	8.1	ng/ul	1114970	4	17.151	2769400	20.0
Pentadecanoic acid	17.351	3.3	ng/ul	457868	4	17.151	2769400	20.0
cis-7-Hexadecen...	17.963	3.9	ng/ul	534118	4	17.151	2769400	20.0
n-Hexadecanoic ...	18.098	15.3	ng/ul	2121490	4	17.151	2769400	20.0
Heptadecanolide	19.292	4.9	ng/ul	674480	4	17.151	2769400	20.0
Octadecanoic acid	19.428	4.3	ng/ul	608167	5	21.604	2852510	20.0
2,2-Bis(4-chlor...	19.904	2.3	ng/ul	334223	5	21.604	2852510	20.0
2,2-Bis(p-chlor...	20.233	4.4	ng/ul	627106	5	21.604	2852510	20.0
unknown-02	20.433	3.6	ng/ul	520790	5	21.604	2852510	20.0
Squalene	23.369	34.5	ng/ul	5249190	6	24.939	3046920	20.0