Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

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-BP111023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP018195.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	3	4	8	rVB	36462	32182	1.46%	0.145%
2	3.158	10	12	19	rVB	16158	21711	0.99%	0.098%
3	3.605	84	88	101	rBV	49269	100224	4.56%	0.453%
4	3.775	113	117	123	rVB3	8869	13041	0.59%	0.059%
5	3.899	135	138	146	rVB4	5860	9782	0.44%	0.044%
6	4.411	220	225	233	rBV2	39544	63302	2.88%	0.286%
7	6.352	552	555	559	rVB5	1680	2265	0.10%	0.010%
8	6.963	652	659	673	rBV	43023	108829	4.95%	0.492%
9	7.134	682	688	702	rBV	229148	366674	16.67%	1.658%
10	7.305	710	717	736	rBV	224000	418787	19.04%	1.893%
11	7.781	790	798	810	rBV	286939	499222	22.70%	2.257%
12	7.969	828	830	835	rVB6	1548	2464	0.11%	0.011%
13	8.516	915	923	937	rBV	146965	310274	14.11%	1.403%
14	8.658	946	947	954	rVB7	2064	3512	0.16%	0.016%
15	8.993	997	1004	1017	rBV	262436	504953	22.96%	2.283%
16	9.116	1022	1025	1027	rVV4	3084	4084	0.19%	0.018%
17	9.163	1027	1033	1036	rVB6	5161	9720	0.44%	0.044%
18	9.705	1117	1125	1144	rBV	224081	447178	20.33%	2.022%
19	10.234	1208	1215	1235	rBV	241128	572577	26.03%	2.588%
20	10.599	1269	1277	1285	rBV	377952	629063	28.60%	2.844%
21	10.652	1285	1286	1290	rVB4	3247	2944	0.13%	0.013%
22	10.693	1290	1293	1298	rVB6	2075	3808	0.17%	0.017%
23	10.787	1302	1309	1323	rBV	147000	378896	17.23%	1.713%
24	10.993	1337	1344	1355	rVB7	13348	37802	1.72%	0.171%
25	11.093	1357	1361	1362	rBV4	2406	2832	0.13%	0.013%
26	11.146	1362	1370	1378	rBV9	7400	22580	1.03%	0.102%
27	11.410	1405	1415	1423	rVV5	18652	58067	2.64%	0.263%
28	11.481	1424	1427	1432	rVV7	6120	12386	0.56%	0.056%
29	11.540	1432	1437	1442	rVB7	6236	13677	0.62%	0.062%
30	11.657	1453	1457	1462	rVB6	5313	8448	0.38%	0.038%
31	12.234	1550	1555	1556	rBV5	3066	4147	0.19%	0.019%
32	13.298	1731	1736	1743	rBV2	22865	40374	1.84%	0.183%
33	13.693	1793	1803	1810	rBV	386118	911458	41.44%	4.120%
34	13.769	1810	1816	1819	rBV4	65476	137708	6.26%	0.623%
35	13.887	1830	1836	1863	rVB	814294	1623282	73.80%	7.338%
36	14.163	1876	1883	1898	rBV	680982	1029937	46.83%	4.656%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : O5338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

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-BP111023.MA.M
Title : SVOA CALIBRATION

37	14.463	1928	1934	1949	rVV	538725	788332	35.84%	3.564%
38	14.810	1985	1993	1996	rBV9	10617	25219	1.15%	0.114%
39	15.198	2055	2059	2061	rBV5	1952	2225	0.10%	0.010%
40	15.245	2064	2067	2073	rVB8	1730	2469	0.11%	0.011%
41	15.334	2078	2082	2087	rVB7	2971	4764	0.22%	0.022%
42	15.463	2098	2104	2114	rBV	890126	1329595	60.45%	6.011%
43	15.628	2126	2132	2144	rBV	244330	432926	19.68%	1.957%
44	15.769	2154	2156	2160	rVB5	3578	3721	0.17%	0.017%
45	15.881	2171	2175	2179	rBV	17332	22318	1.01%	0.101%
46	15.922	2179	2182	2187	rVV3	7705	10864	0.49%	0.049%
47	15.963	2187	2189	2196	rVB6	5427	7400	0.34%	0.033%
48	16.098	2207	2212	2213	rBV5	1918	2523	0.11%	0.011%
49	16.181	2221	2226	2232	rBV	47816	67585	3.07%	0.306%
50	16.275	2236	2242	2246	rVV	157805	206347	9.38%	0.933%
51	16.334	2247	2252	2258	rVV2	178888	364544	16.57%	1.648%
52	16.398	2258	2263	2267	rVV3	154830	263663	11.99%	1.192%
53	16.439	2267	2270	2275	rVV	92234	127825	5.81%	0.578%
54	16.498	2275	2280	2282	rVV	42997	75413	3.43%	0.341%
55	16.545	2286	2288	2292	rVV	44308	50188	2.28%	0.227%
56	16.598	2292	2297	2304	rVV4	96877	195838	8.90%	0.885%
57	16.669	2304	2309	2313	rVV	143308	202920	9.23%	0.917%
58	16.710	2314	2316	2319	rVV	41458	58533	2.66%	0.265%
59	16.745	2319	2322	2331	rVB	66207	90711	4.12%	0.410%
60	16.822	2331	2335	2337	rVB5	2193	2768	0.13%	0.013%
61	16.881	2340	2345	2347	rBV6	5827	8833	0.40%	0.040%
62	16.957	2355	2358	2359	rBV3	2486	2500	0.11%	0.011%
63	16.992	2361	2364	2366	rVV3	4965	7018	0.32%	0.032%
64	17.045	2370	2373	2376	rBV5	3898	5318	0.24%	0.024%
65	17.081	2376	2379	2382	rVB4	3343	3880	0.18%	0.018%
66	17.245	2398	2407	2415	rBV	709982	1006203	45.75%	4.549%
67	17.345	2417	2424	2434	rBV	1103263	1571222	71.44%	7.103%
68	17.581	2461	2464	2467	rVB5	4754	5342	0.24%	0.024%
69	17.669	2476	2479	2483	rVB4	5718	7160	0.33%	0.032%
70	17.751	2491	2493	2497	rVB5	3996	4313	0.20%	0.019%
71	17.792	2497	2500	2503	rBV5	3879	5064	0.23%	0.023%
72	17.875	2511	2514	2518	rBV5	13895	17982	0.82%	0.081%
73	18.092	2547	2551	2564	rBV2	47093	89762	4.08%	0.406%
74	18.186	2564	2567	2571	rVB3	5691	8696	0.40%	0.039%
75	18.316	2584	2589	2593	rVB5	6202	10410	0.47%	0.047%
76	19.175	2732	2735	2739	rBV3	13341	14574	0.66%	0.066%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

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-BP111023.MA.M
Title : SVOA CALIBRATION

77	19.228	2741	2744	2746	rBV3	19633	22532	1.02%	0.102%
78	19.339	2759	2763	2769	rBV5	39627	63121	2.87%	0.285%
79	19.598	2802	2807	2817	rBV	1395930	1916304	87.13%	8.663%
80	21.374	3104	3109	3119	rBV2	877998	1162198	52.84%	5.254%
81	23.680	3493	3501	3510	rBV	1136537	2199486	100.00%	9.943%
82	23.845	3522	3529	3540	rVB	585052	1241745	56.46%	5.614%

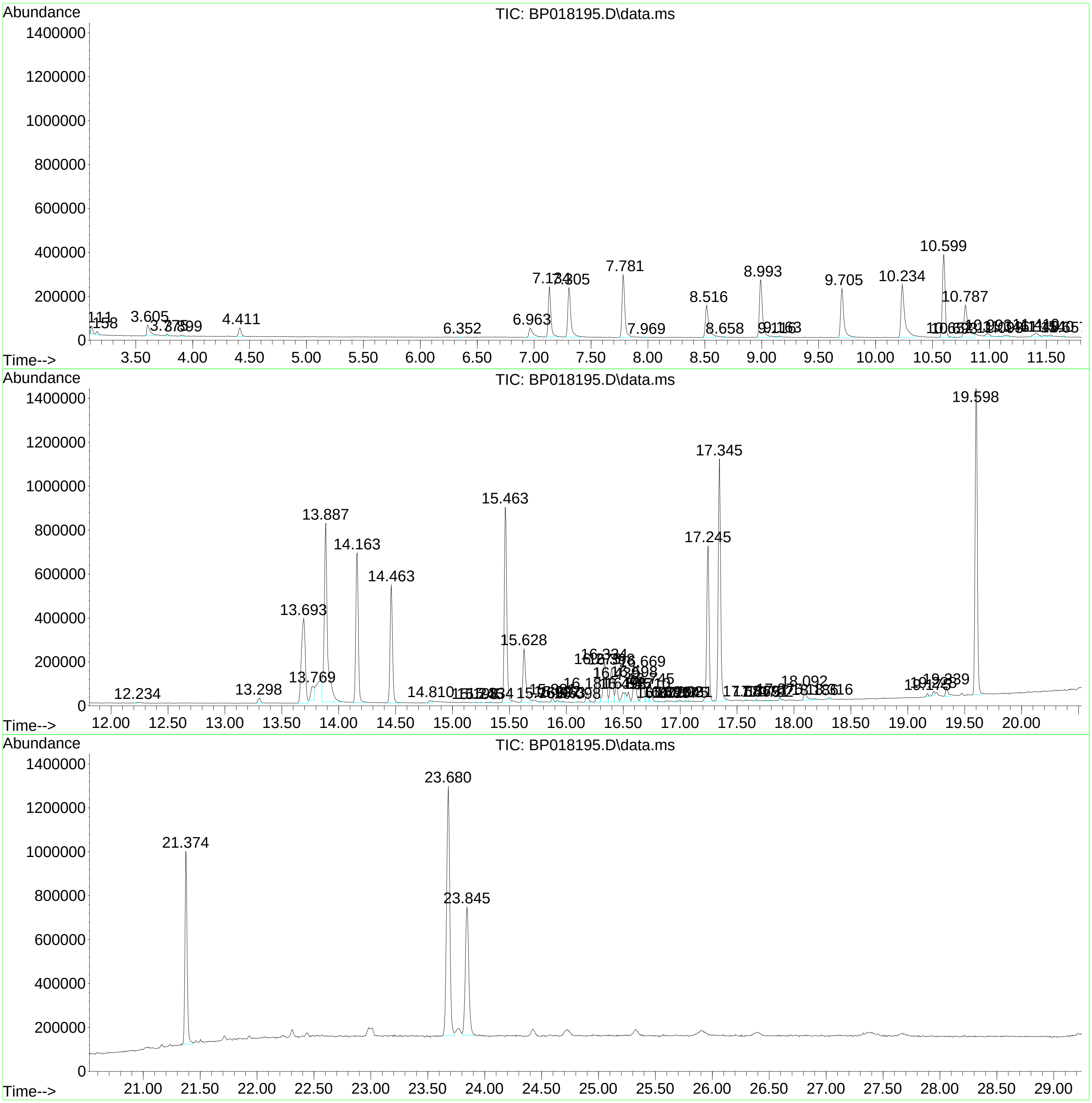
Sum of corrected areas: 22120544

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.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\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

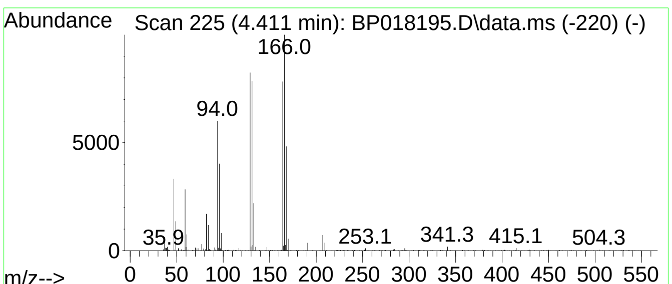
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

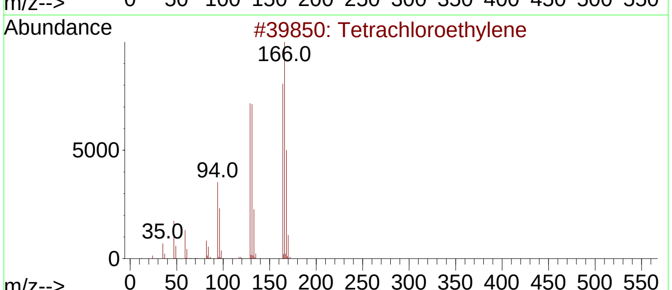
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

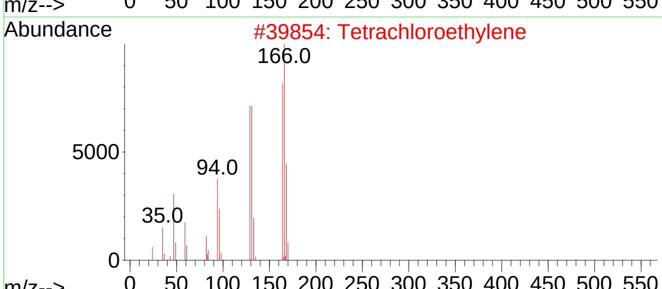
****
Peak Number 1 Tetrachloroethylene Concentration Rank 9

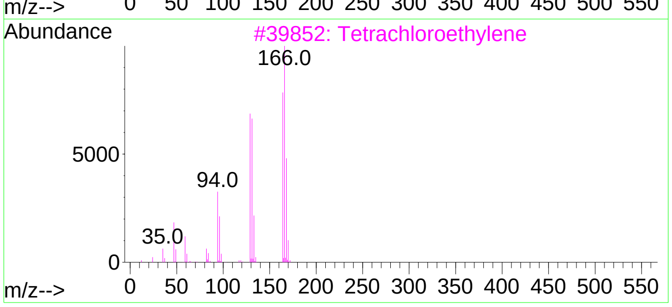
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.411	2.54 ng/ul	63302	1,4-Dichlorobenzene-d4	7.781

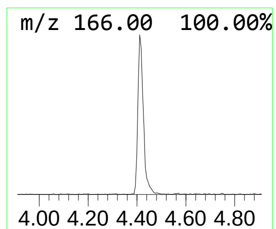
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrachloroethylene	164	C2Cl4	000127-18-4	94
2	Tetrachloroethylene	164	C2Cl4	000127-18-4	94
3	Tetrachloroethylene	164	C2Cl4	000127-18-4	94
4	Tetrachloroethylene	164	C2Cl4	000127-18-4	93
5	2-Chloroquinoxaline	164	C8H5ClN2	001448-87-9	12

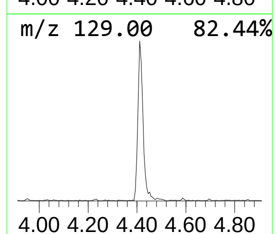


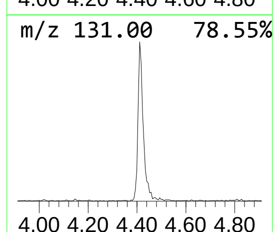


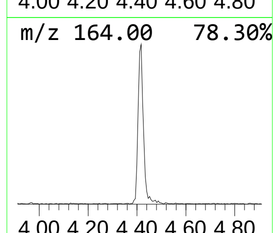


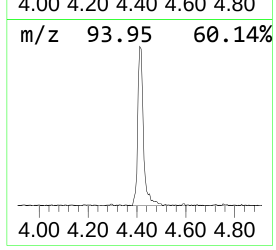












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

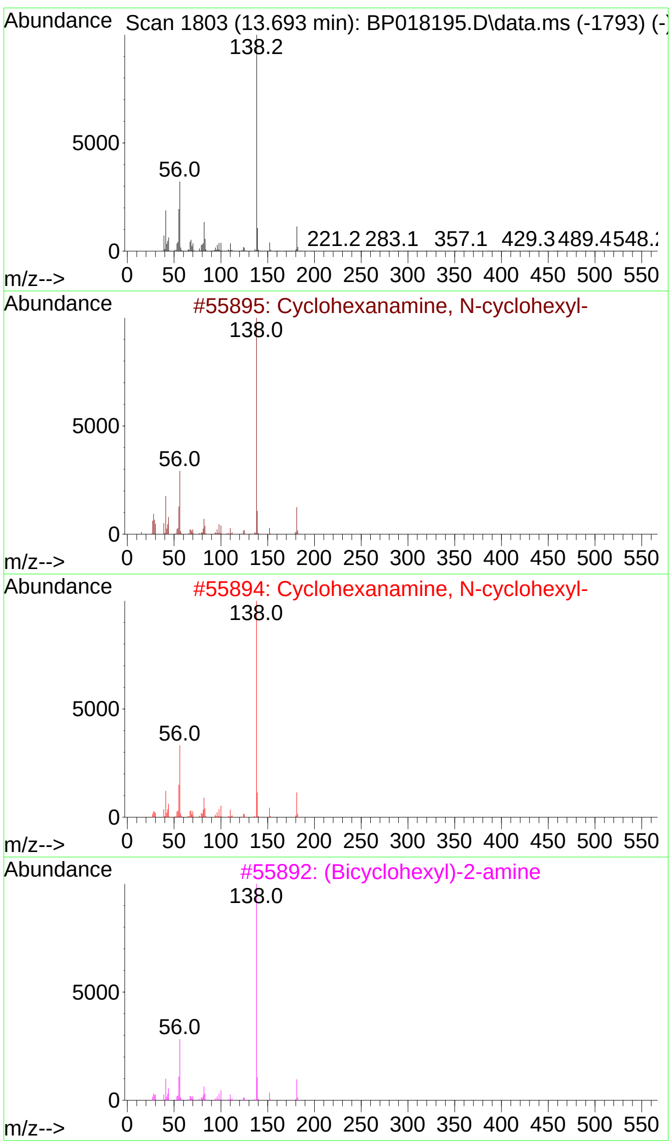
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

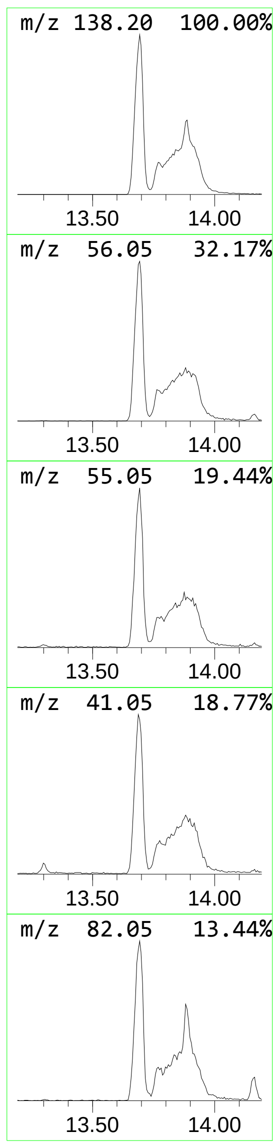
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclohexanamine, N-cyclohexyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.693	23.12 ng/ul	911458	Acenaphthene-d10	14.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	97
2	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	96
3	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	94
4	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	91
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

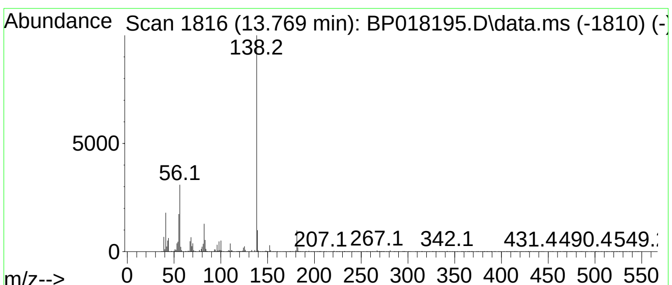
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

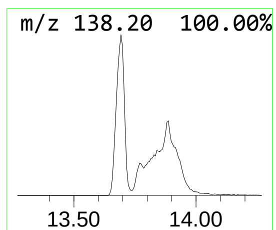
****

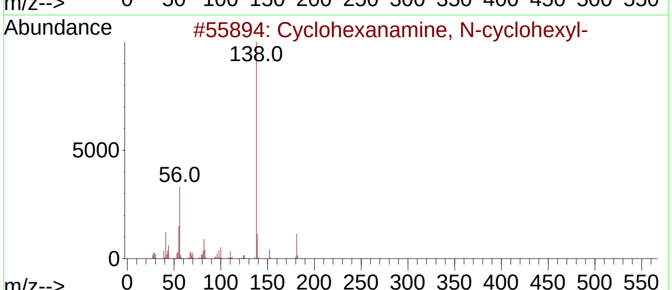
Peak Number 3 (Bicyclohexyl)-2-amine Concentration Rank 7

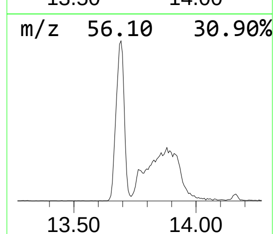
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.769	3.49 ng/ul	137708	Acenaphthene-d10	14.463

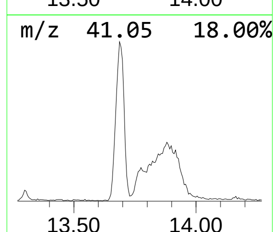
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	96
2	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
3	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
4	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	93
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90

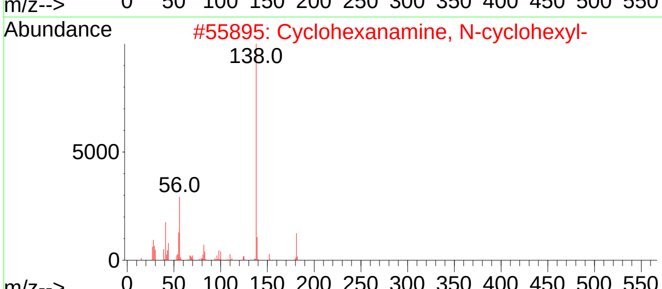


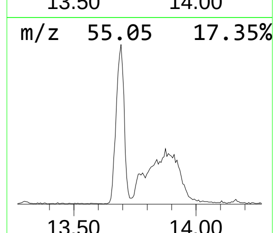


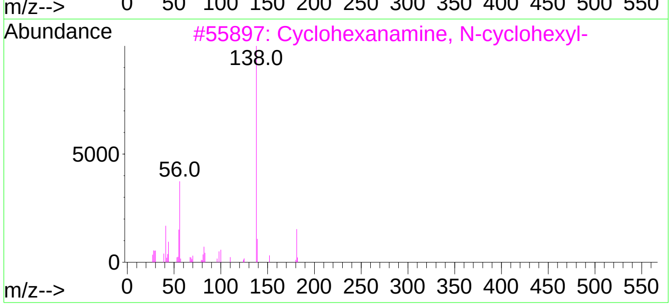


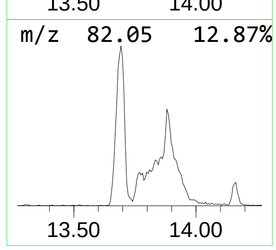












Library Search Compound Report

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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On    : 16 Nov 2023  03:51
Operator  : MA/JU
Sample    : O5338-05
Misc      :
ALS Vial  : 22   Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title  : SVOA CALIBRATION

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

Instrument :
BNA_P
ClientSampleId :
GCDL0

```
*********************************************************************
Peak Number  4  Phenol, 4-(1,1,3,3-tetramet...  Concentration Rank  4

  R.T.     EstConc          Area       Relative to ISTD         R.T.
---------------------------------------------------------------------
 16.275    4.10 ng/ul      206347    Phenanthrene-d10           17.245

 Hit# of  5    Tentative ID              MW  MolForm       CAS#       Qual
---------------------------------------------------------------------
 1 Phenol, 4-(1,1,3,3-tetramethylbu... 206 C14H22O       000140-66-9 86
 2 Phenol, 4-(1,1-dimethylpropyl)-     164 C11H16O       000080-46-6 78
 3 ((1,2-Diethylethylene)bis(p-phen... 354 C22H26O4      004547-76-6 72
 4 2-tert-Butylphenol, acetate         192 C12H16O2      003245-25-8 72
 5 Phenol, 4-(1,1,3,3-tetramethylbu... 206 C14H22O       000140-66-9 72
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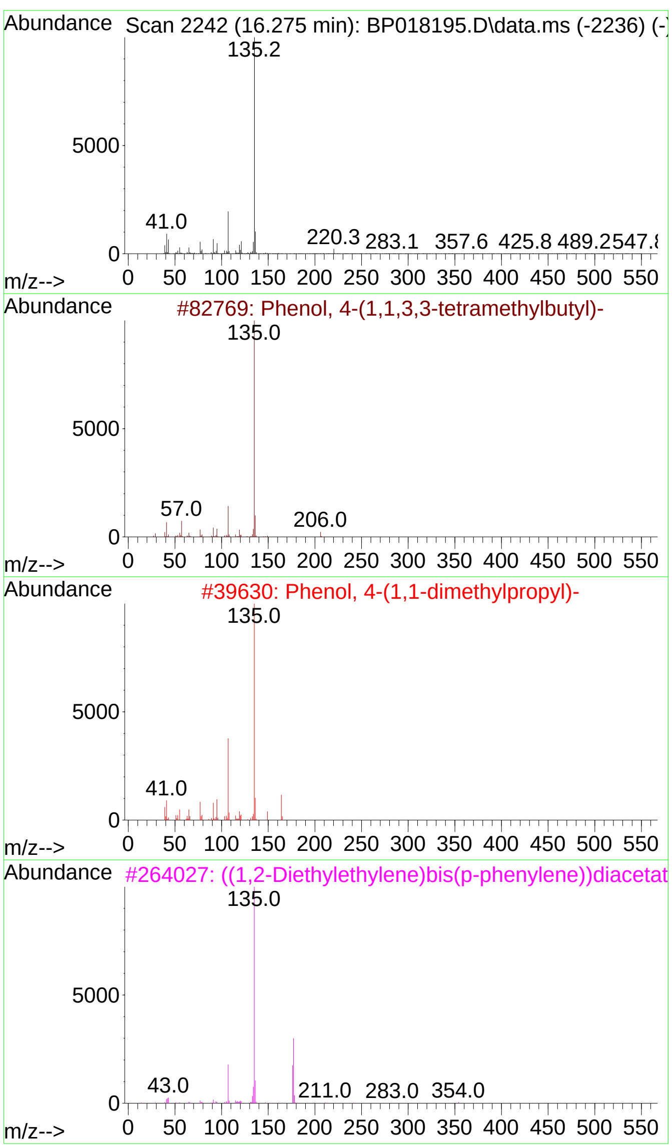

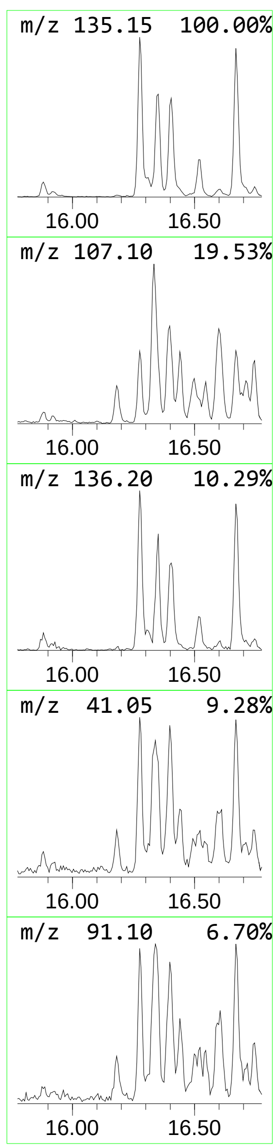

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Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

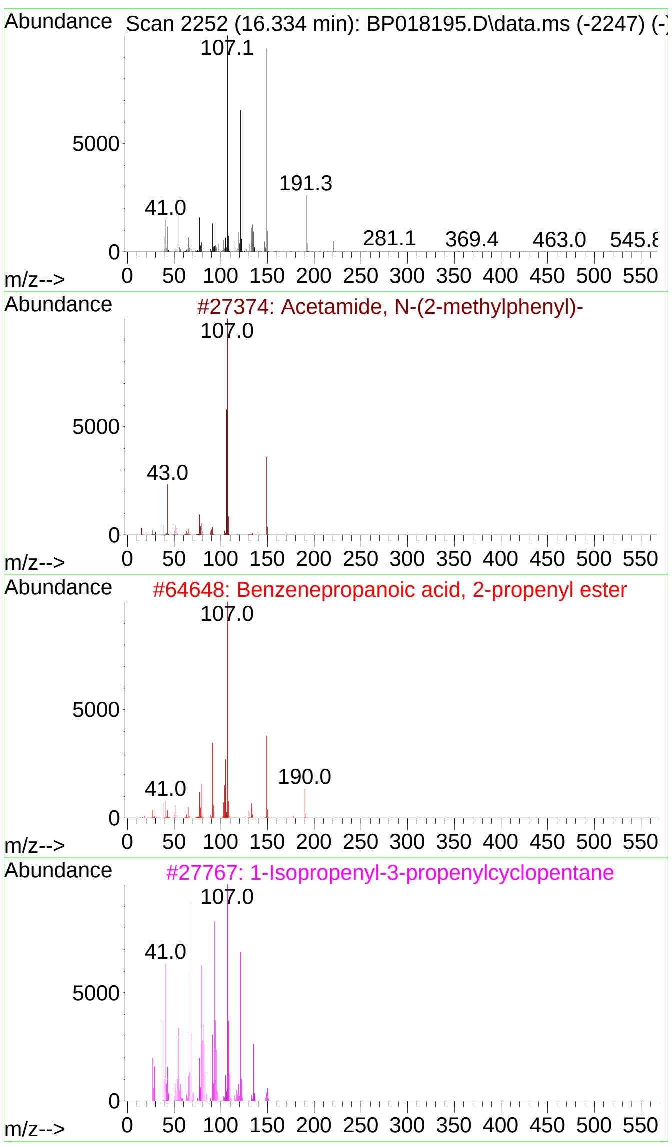
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

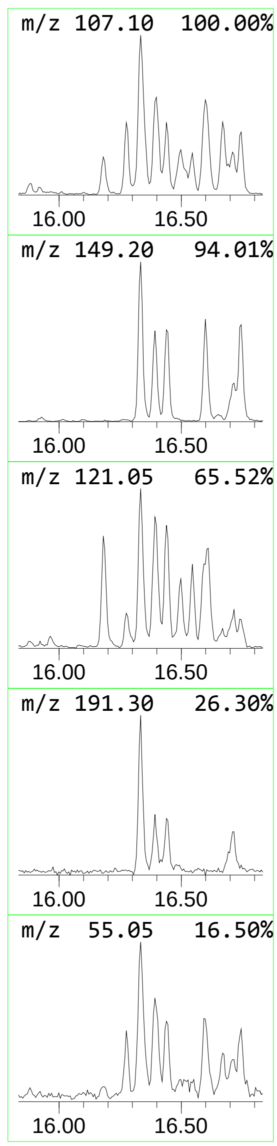
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Acetamide, N-(2-methylphenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.334	7.25 ng/ul	364544	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	53
2	Benzenepropanoic acid, 2-propeny...	190	C12H14O2	015814-45-6	47
3	1-Isopropenyl-3-propenylcyclopen...	150	C11H18	1000192-81-2	46
4	Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	43
5	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43





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Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

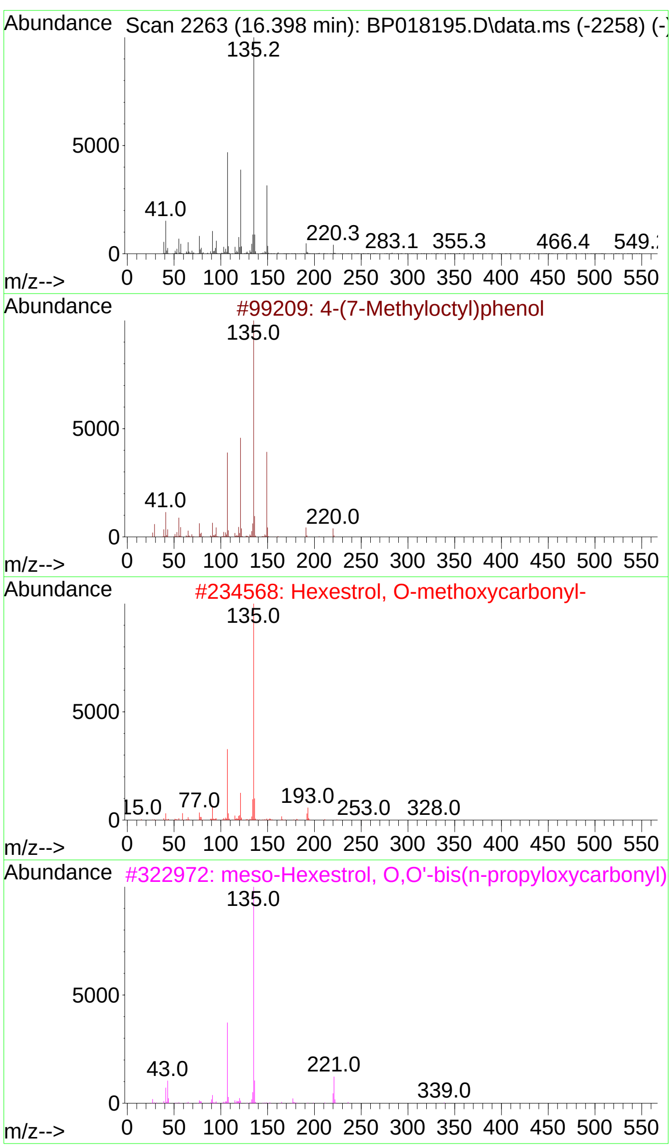
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

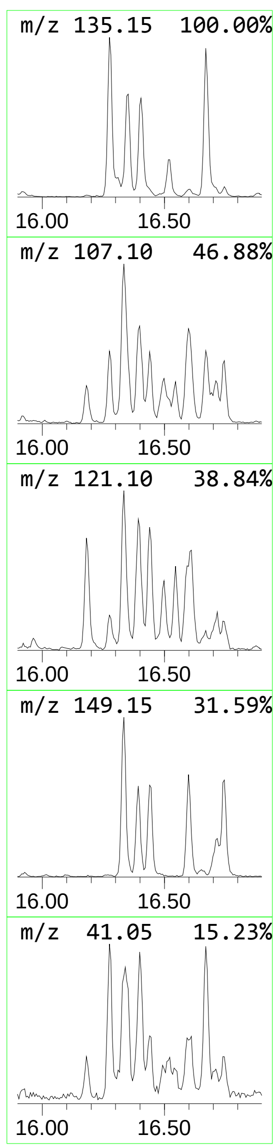
****

Peak Number 6 4-(7-Methyloctyl)phenol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	5.24 ng/ul	263663	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	98
2	Hexestrol, O-methoxycarbonyl-	328	C20H24O4	1000487-66-3	64
3	meso-Hexestrol, O,O'-bis(n-propy...	442	C26H34O6	1000487-65-0	59
4	Hexestrol, O-(n-propyloxycarbonyl)-	356	C22H28O4	1000487-65-1	59
5	4,6-Bis(4-ethoxybenzylthio)-5-ni...	457	C22H23N3O4S2	325957-76-4	53





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Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : O5338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

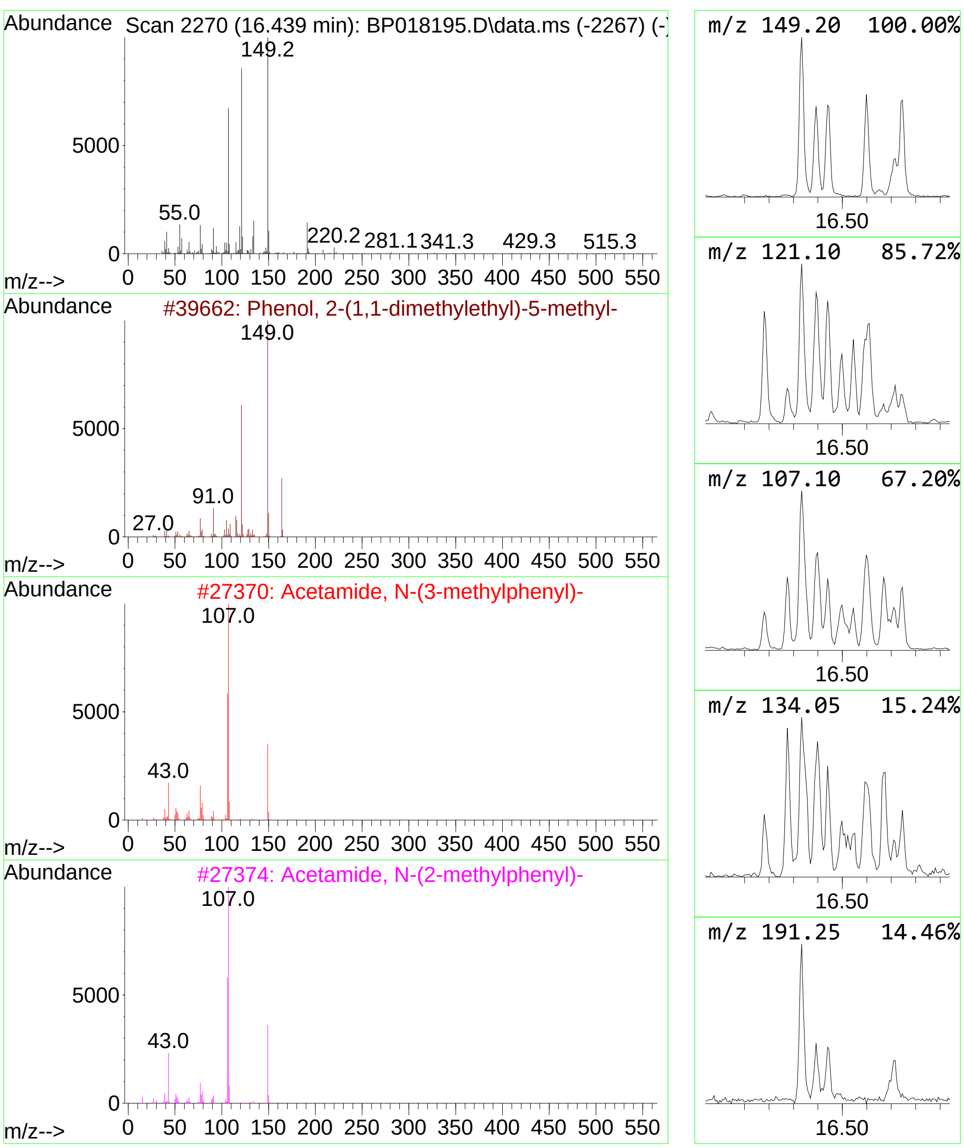
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

****
Peak Number 7 Phenol, 2-(1,1-dimethylethy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.439	2.54 ng/ul	127825	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	53
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
3	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38
4	3',4'-Formoxylidide	149	C9H11NO	006639-60-7	38
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	35



Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

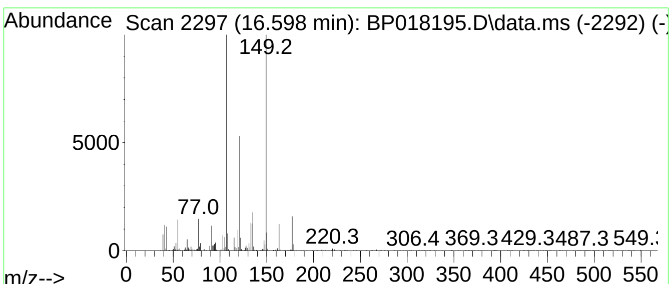
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

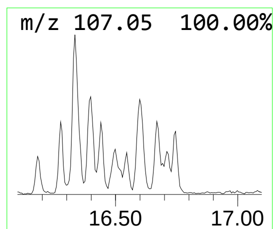
****

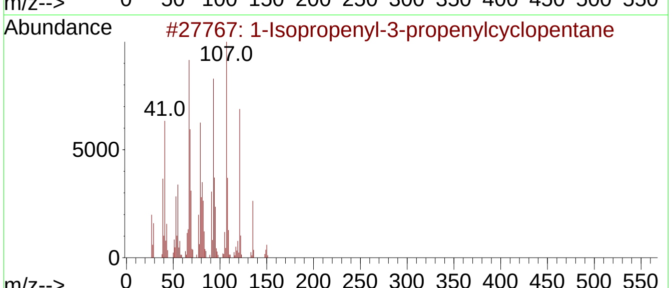
Peak Number 8 unknown-01 Concentration Rank 6

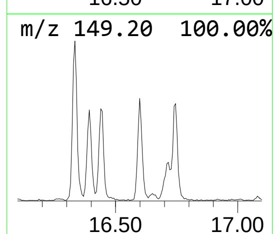
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	3.89 ng/ul	195838	Phenanthrene-d10	17.245

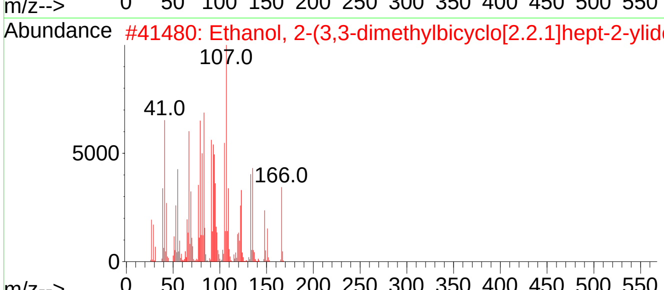
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Isopropenyl-3-propenylcyclopen...	150	C11H18	1000192-81-2	42
2	Ethanol, 2-(3,3-dimethylbicyclo[...	166	C11H18O	002226-05-3	38
3	Benzenepropanoic acid, 2-propeny...	190	C12H14O2	015814-45-6	35
4	2-(4-Methoxy-phenyl)-2-methyl-pr...	208	C12H16O3	006274-50-6	35
5	1-Methyl-3-ethyladamantane	178	C13H22	001687-34-9	35

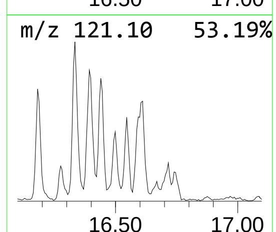


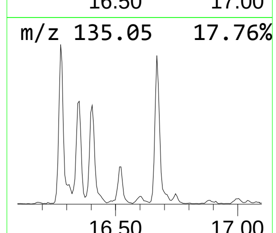


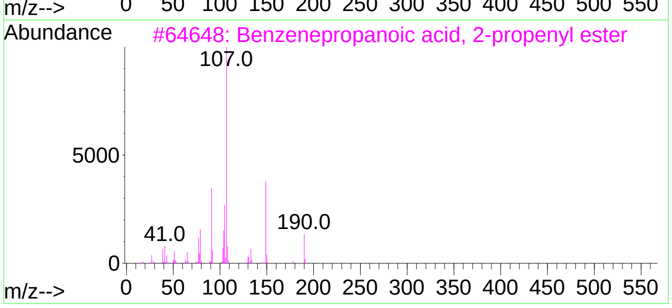


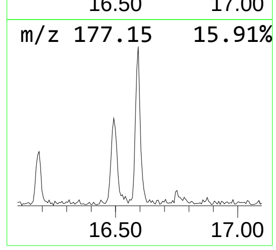












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Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : 05338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

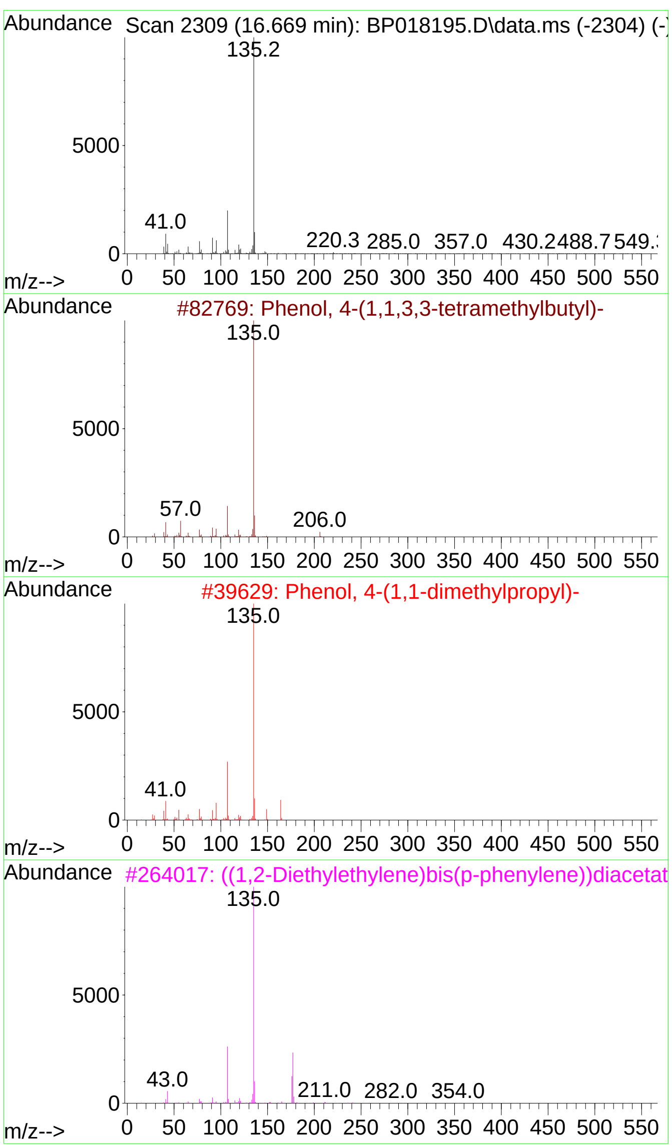
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

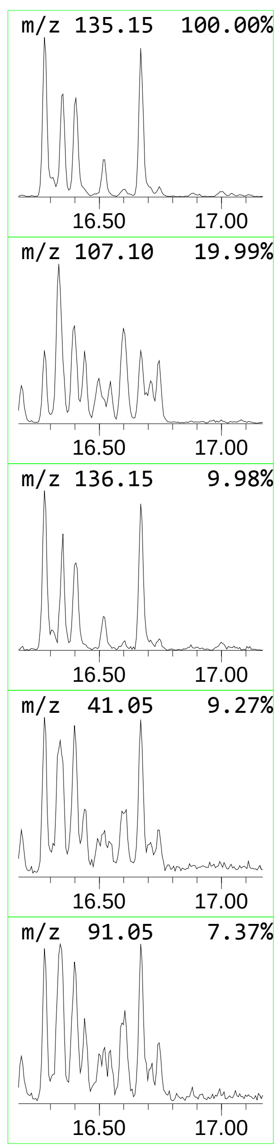
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, 4-(1,1-dimethylprop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	4.03 ng/ul	202920	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	91
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	72
4	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	72
5	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018195.D
Acq On : 16 Nov 2023 03:51
Operator : MA/JU
Sample : O5338-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.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
Tetrachloroethy...	4.411	2.5	ng/ul	63302	1	7.781	499222	20.0
Cyclohexanamine...	13.693	23.1	ng/ul	911458	3	14.463	788332	20.0
(Bicyclohexyl)-...	13.769	3.5	ng/ul	137708	3	14.463	788332	20.0
Phenol, 4-(1,1,...	16.275	4.1	ng/ul	206347	4	17.245	1006200	20.0
Acetamide, N-(2...	16.334	7.3	ng/ul	364544	4	17.245	1006200	20.0
4-(7-Methylocty...	16.398	5.2	ng/ul	263663	4	17.245	1006200	20.0
Phenol, 2-(1,1-...	16.439	2.5	ng/ul	127825	4	17.245	1006200	20.0
unknown-01	16.598	3.9	ng/ul	195838	4	17.245	1006200	20.0
Phenol, 4-(1,1-...	16.669	4.0	ng/ul	202920	4	17.245	1006200	20.0