LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

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: BP018258.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	9	12	18	rVB	13588	16975	0.89%	0.095%
2	3.369	46	48	51	rBV4	2243	2588	0.14%	0.015%
3	3.605	85	88	102	rBV	50705	103744	5.46%	0.583%
4	3.769	114	116	122	rVB7	5284	8167	0.43%	0.046%
5	3.899	134	138	142	rBV6	2818	4053	0.21%	0.023%
6	4.487	233	238	241	rBV7	1942	3528	0.19%	0.020%
7	6.963	652	659	679	rBV	38520	107697	5.67%	0.605%
8	7.128	681	687	701	rBV	207890	345510	18.17%	1.941%
9	7.299	709	716	739	rBV	214276	401604	21.13%	2.256%
10	7.510	750	752	756	rVB5	2316	2148	0.11%	0.012%
11	7.781	791	798	813	rBV	161340	300292	15.80%	1.687%
12	7.881	813	815	817	rVV3	2609	2639	0.14%	0.015%
13	7.922	821	822	826	rVB4	1832	1942	0.10%	0.011%
14	8.510	916	922	944	rBV	146625	314123	16.52%	1.764%
15	8.987	996	1003	1015	rBV	251970	474596	24.97%	2.666%
16	9.146	1028	1030	1035	rVB5	4179	5216	0.27%	0.029%
17	9.699	1117	1124	1145	rBV	213716	439610	23.12%	2.469%
18	10.228	1208	1214	1243	rBV	232044	560284	29.47%	3.147%
19	10.593	1268	1276	1290	rVV	231632	392170	20.63%	2.203%
20	10.781	1301	1308	1330	rBV	175655	465553	24.49%	2.615%
21	10.987	1341	1343	1350	rVB8	4000	5944	0.31%	0.033%
22	11.040	1350	1352	1358	rVB7	1883	3211	0.17%	0.018%
23	11.375	1405	1409	1410	rBV4	2264	2417	0.13%	0.014%
24	11.410	1410	1415	1419	rVB7	3785	7405	0.39%	0.042%
25	11.528	1433	1435	1439	rVB5	2121	2225	0.12%	0.012%
26	11.628	1450	1452	1458	rVB7	1461	1939	0.10%	0.011%
27	11.763	1473	1475	1478	rVB4	2334	2119	0.11%	0.012%
28	12.251	1556	1558	1563	rVB6	3129	2744	0.14%	0.015%
29	12.445	1587	1591	1595	rVV7	1696	2508	0.13%	0.014%
30	12.757	1641	1644	1647	rBV4	1836	2130	0.11%	0.012%
31	13.298	1729	1736	1742	rBV	39040	66226	3.48%	0.372%
32	13.669	1793	1799	1804	rBV	91822	147054	7.74%	0.826%
33	13.739	1804	1811	1830	rVV2	56076	243294	12.80%	1.367%
34	13.881	1830	1835	1851	rVB	667972	979742	51.54%	5.503%
35	14.016	1857	1858	1863	rBV5	2038	2181	0.11%	0.012%
36	14.157	1874	1882	1895	rBV	715184	1052349	55.36%	5.911%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

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	14.457	1926	1933	1946	rVB	323670	471533	24.80%	2.649%
38	14.816	1989	1994	1995	rBV5	7405	9701	0.51%	0.054%
39	14.881	2001	2005	2010	rVB8	2833	4008	0.21%	0.023%
40	15.104	2040	2043	2047	rVB6	2209	2597	0.14%	0.015%
41	15.463	2097	2104	2115	rBV	920759	1366398	71.88%	7.675%
42	15.628	2126	2132	2154	rBV	233642	461839	24.29%	2.594%
43	15.875	2170	2174	2178	rVV	16462	21544	1.13%	0.121%
44	15.922	2178	2182	2187	rVV3	15921	23538	1.24%	0.132%
45	15.963	2187	2189	2193	rVV4	3809	4215	0.22%	0.024%
46	16.181	2220	2226	2235	rBV	32223	42401	2.23%	0.238%
47	16.275	2237	2242	2246	rVV	114055	154586	8.13%	0.868%
48	16.333	2246	2252	2258	rVV2	127763	258883	13.62%	1.454%
49	16.398	2258	2263	2267	rVV3	105470	189302	9.96%	1.063%
50	16.439	2267	2270	2274	rVV	69131	90642	4.77%	0.509%
51	16.492	2274	2279	2281	rVV	37051	56340	2.96%	0.316%
52	16.516	2281	2283	2285	rVV	35840	41794	2.20%	0.235%
53	16.539	2285	2287	2291	rVV	37435	50317	2.65%	0.283%
54	16.598	2291	2297	2304	rVV3	91135	181064	9.52%	1.017%
55	16.669	2304	2309	2313	rVV	110159	158696	8.35%	0.891%
56	16.710	2314	2316	2318	rVV3	28716	34448	1.81%	0.193%
57	16.739	2318	2321	2327	rVV	49344	70674	3.72%	0.397%
58	16.810	2330	2333	2337	rVB6	2037	3477	0.18%	0.020%
59	16.875	2340	2344	2349	rBV5	3667	7429	0.39%	0.042%
60	16.992	2361	2364	2368	rVB5	4418	7168	0.38%	0.040%
61	17.039	2371	2372	2376	rVV4	2901	2699	0.14%	0.015%
62	17.075	2376	2378	2382	rVV4	3321	4939	0.26%	0.028%
63	17.192	2395	2398	2401	rBV5	2358	3456	0.18%	0.019%
64	17.245	2401	2407	2415	rBV	458798	647152	34.04%	3.635%
65	17.339	2417	2423	2437	rBV	1050736	1593207	83.81%	8.949%
66	17.875	2509	2514	2519	rVB7	10305	14298	0.75%	0.080%
67	18.092	2547	2551	2558	rBV3	25830	42154	2.22%	0.237%
68	18.186	2565	2567	2571	rVB4	2371	1949	0.10%	0.011%
69	18.304	2583	2587	2594	rBV3	23667	36593	1.92%	0.206%
70	18.527	2623	2625	2629	rBV5	1735	2549	0.13%	0.014%
71	18.563	2629	2631	2635	rVB4	2141	2582	0.14%	0.015%
72	18.733	2658	2660	2663	rVB4	2411	2371	0.12%	0.013%
73	18.769	2663	2666	2668	rBV5	2975	3366	0.18%	0.019%
74	18.904	2687	2689	2692	rVB3	2471	2173	0.11%	0.012%
75	19.016	2705	2708	2712	rVB5	7663	9550	0.50%	0.054%
76	19.174	2732	2735	2739	rBV2	10050	12218	0.64%	0.069%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

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	19.227	2742	2744	2745	rVV2	5153	4270	0.22%	0.024%
78	19.251	2746	2748	2757	rVV4	14616	23588	1.24%	0.132%
79	19.345	2760	2764	2776	rVB9	19334	33220	1.75%	0.187%
80	19.598	2802	2807	2820	rBV	1390526	1898774	99.88%	10.665%
81	21.380	3105	3110	3117	rBV	485672	677561	35.64%	3.806%
82	23.686	3495	3502	3514	rBV2	949116	1901018	100.00%	10.678%
83	23.857	3525	3531	3542	rVB	324215	694891	36.55%	3.903%

Sum of corrected areas: 17803099

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
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Operator : MA/JU
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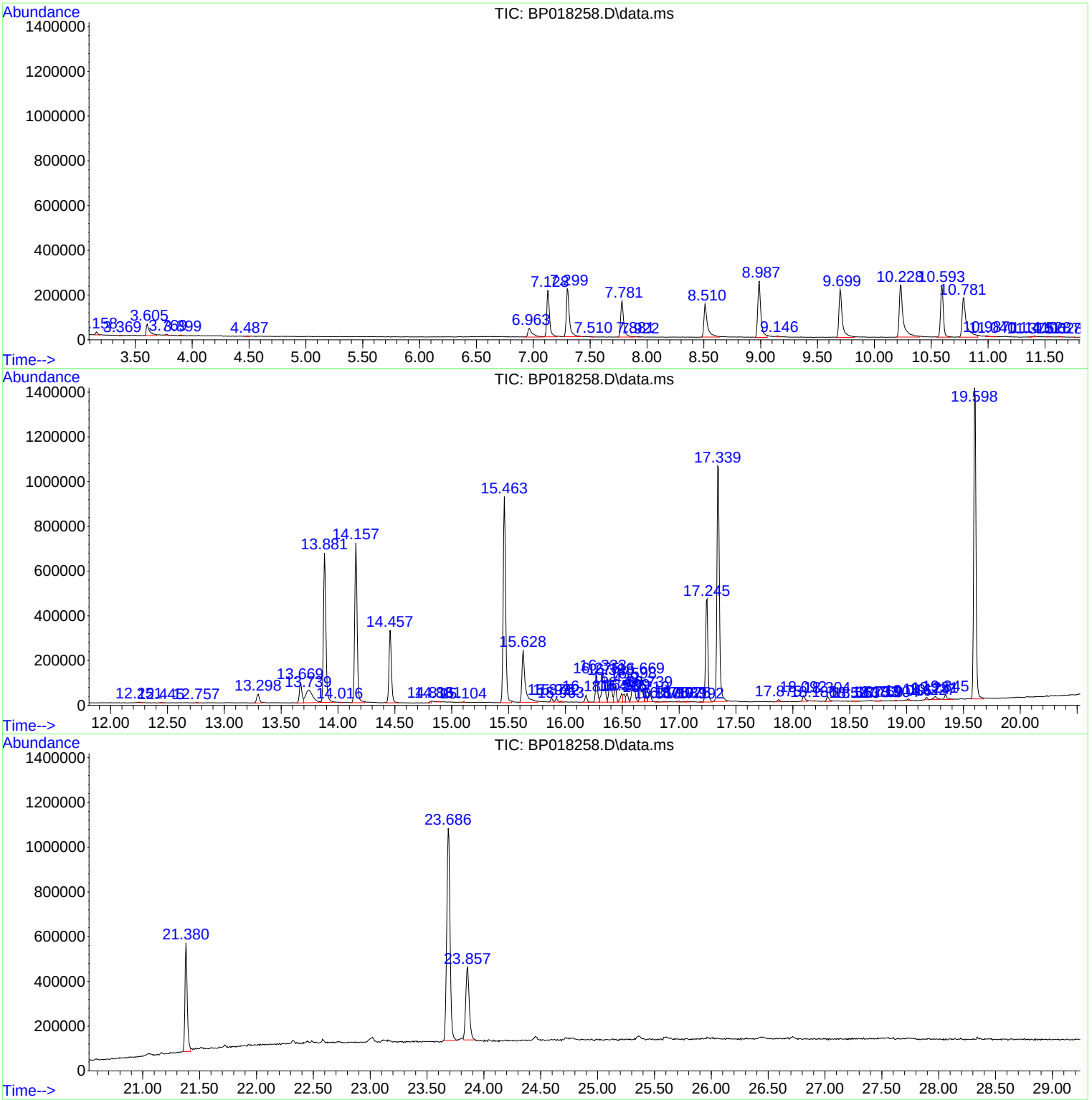
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

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\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

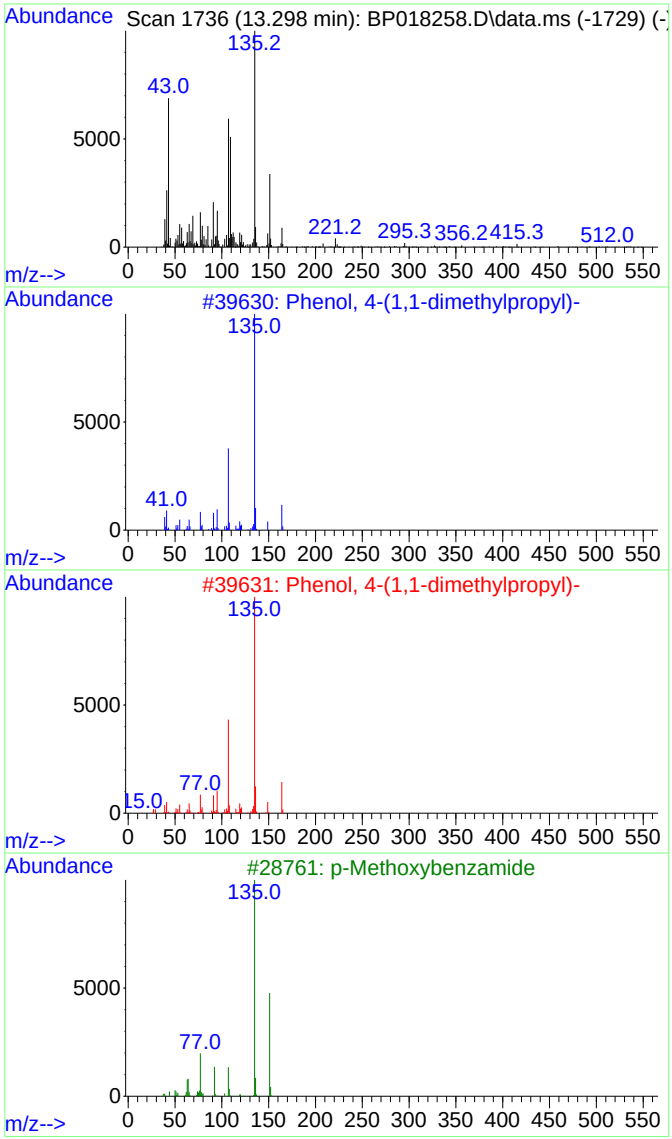
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

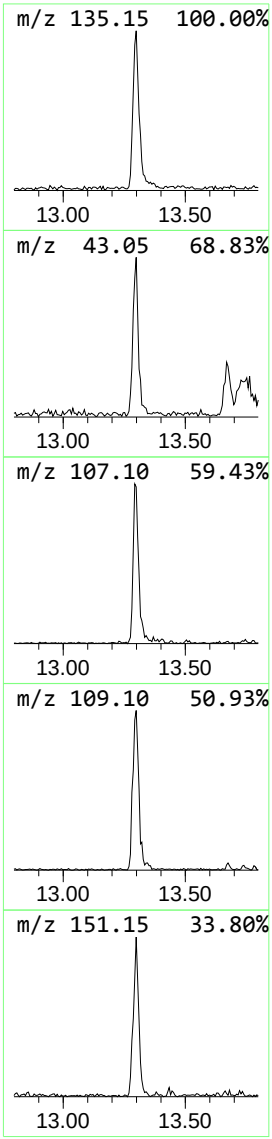
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Phenol, 4-(1,1-dimethylprop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.298	2.81 ng/ul	66226	Acenaphthene-d10	14.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	60
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	60
3	p-Methoxybenzamide	151	C8H9NO2	003424-93-9	49
4	Thymol	150	C10H14O	000089-83-8	35
5	3-(Adamantan-1-yl)propanoic acid	208	C13H20O2	016269-16-2	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

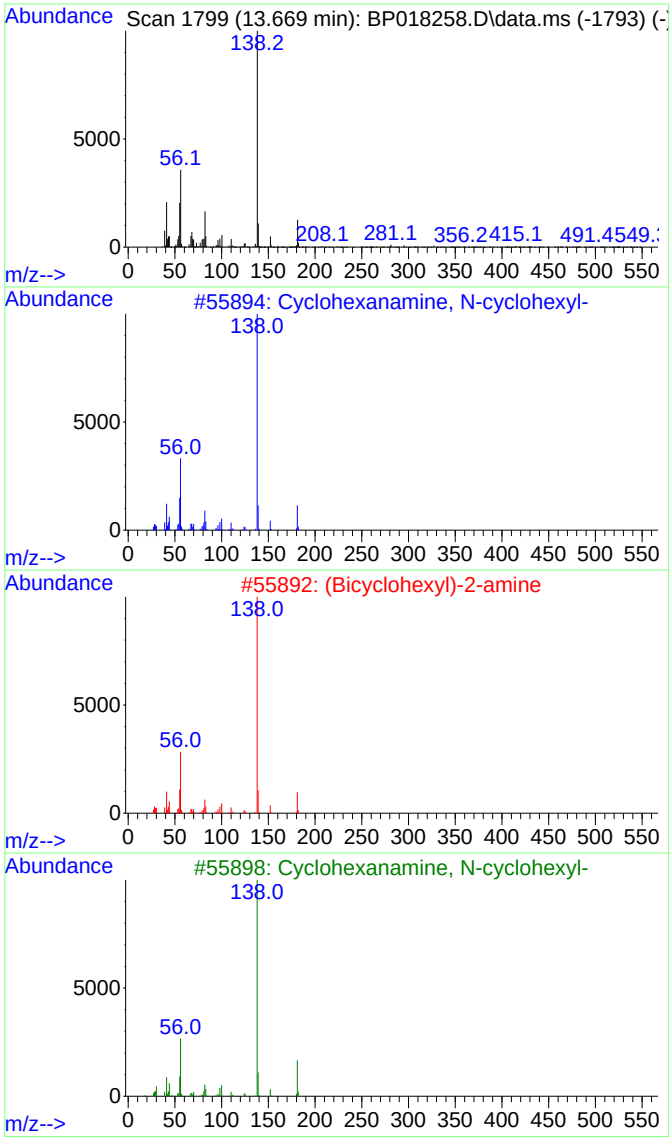
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

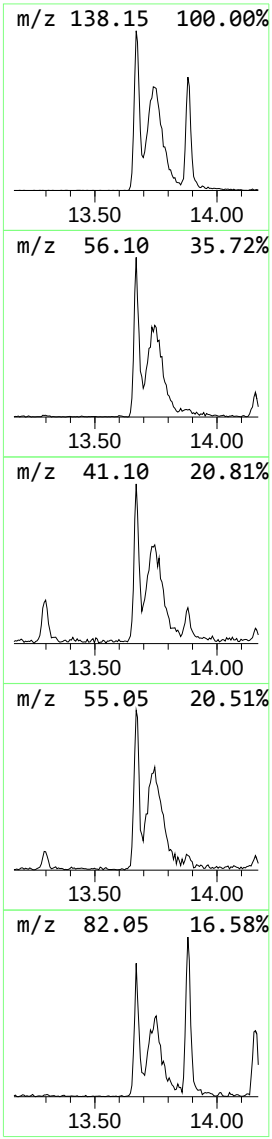
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclohexanamine, N-cyclohexyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.669	6.24 ng/ul	147054	Acenaphthene-d10	14.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	97
2	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	95
3	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	91
4	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
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Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

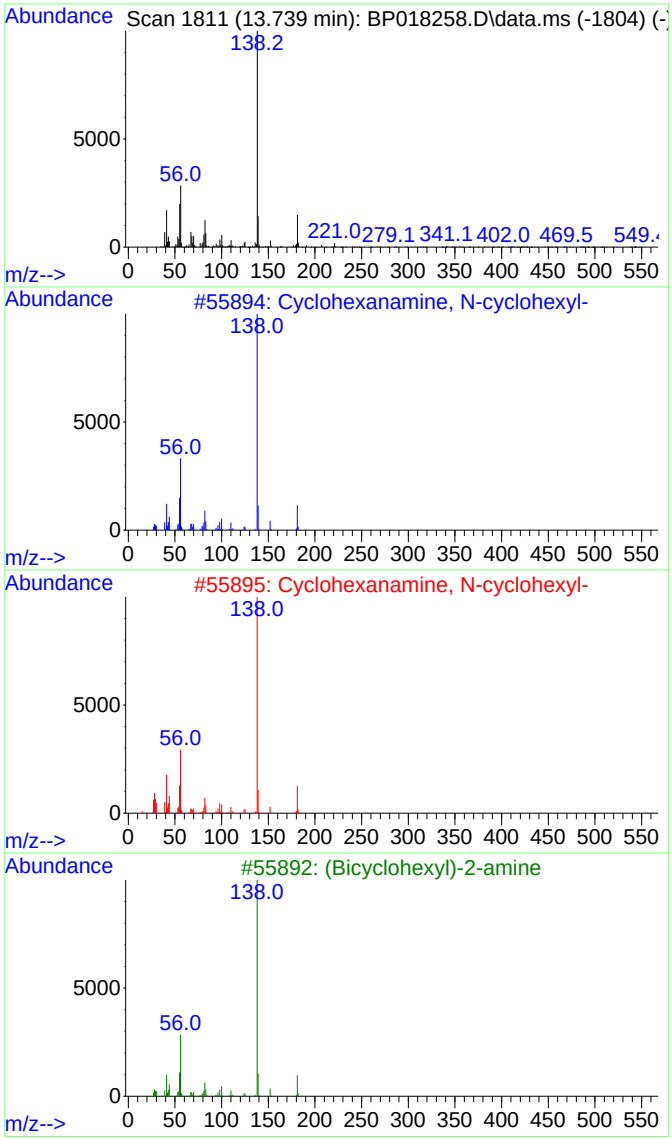
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

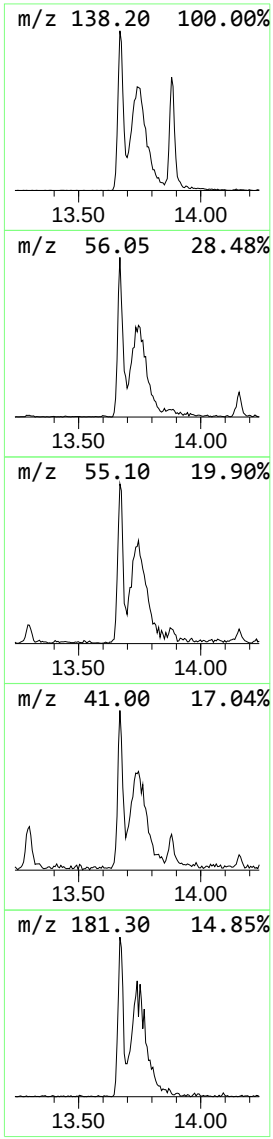
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (Bicyclohexyl)-2-amine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.740	10.32 ng/ul	243294	Acenaphthene-d10	14.457

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
2	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	93
3	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	91
4	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	91
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
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Acq On : 18 Nov 2023 13:15
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Instrument :
BNA_P
ClientSampleId :
GCDQ7

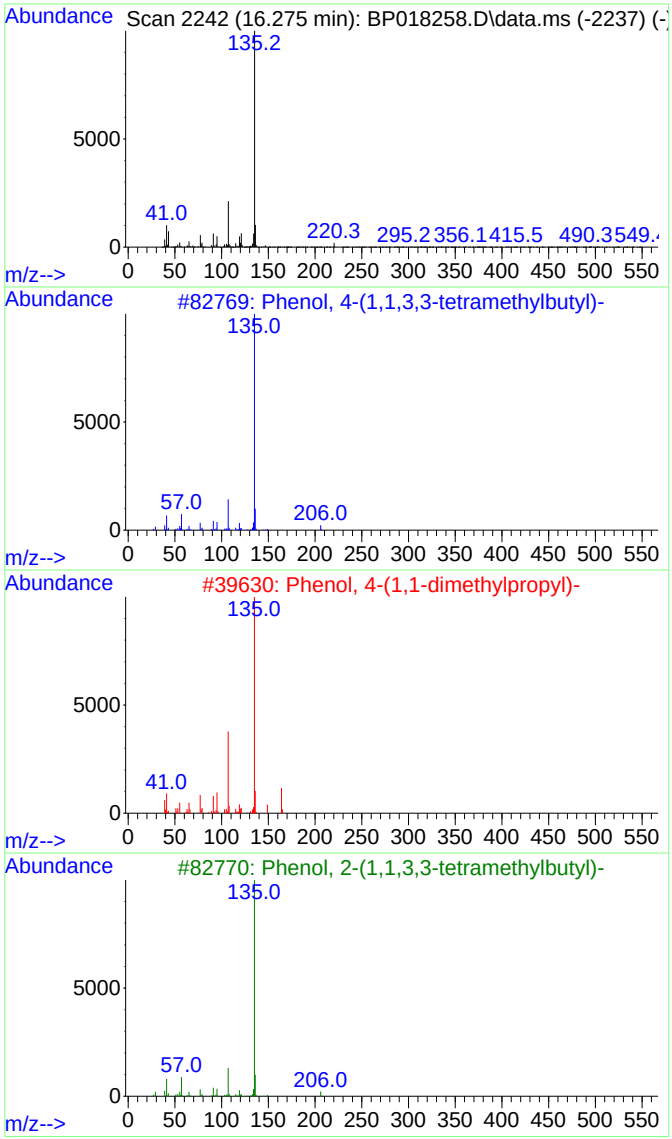
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

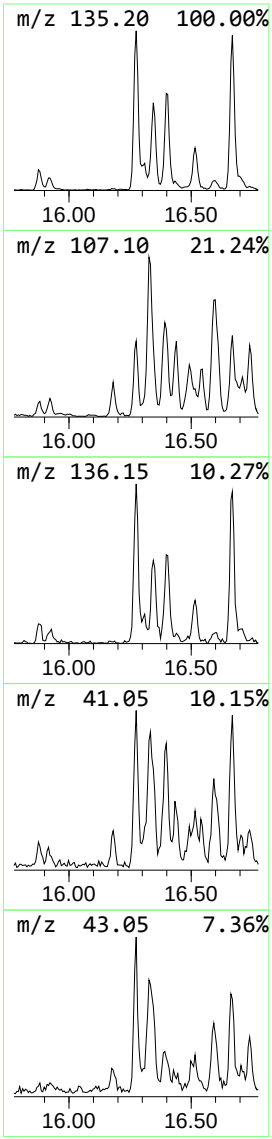
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	4.78 ng/ul	154586	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	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
4	Methanol, (cyclohexyl)(2,3-dimet...	218	C15H22O	349498-47-1	72
5	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
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Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

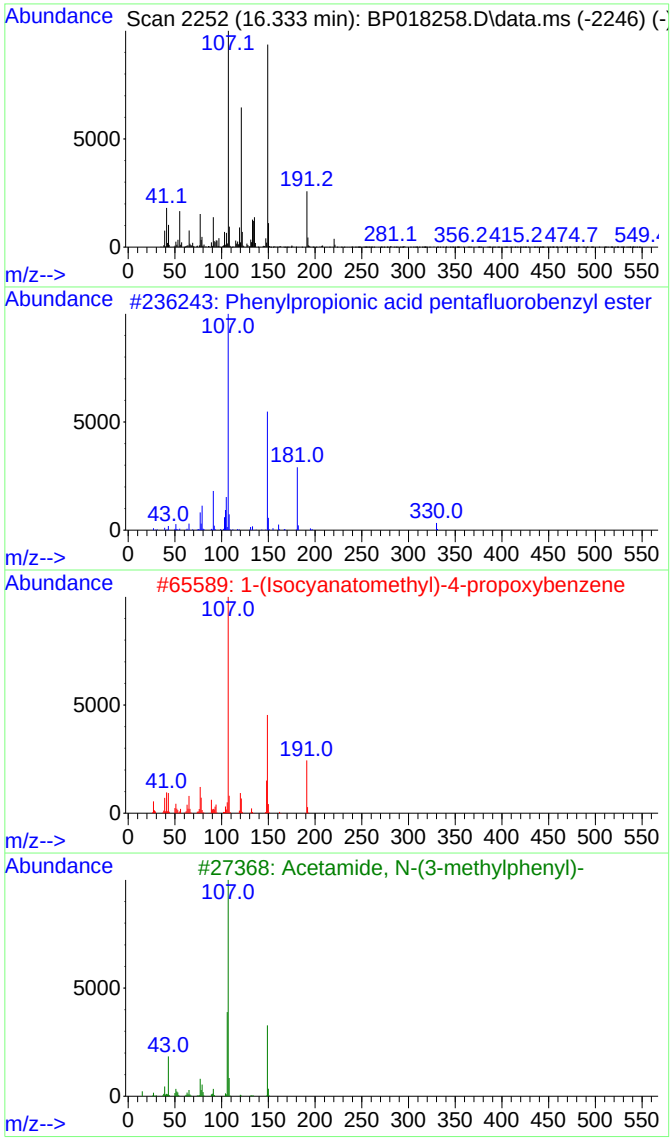
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

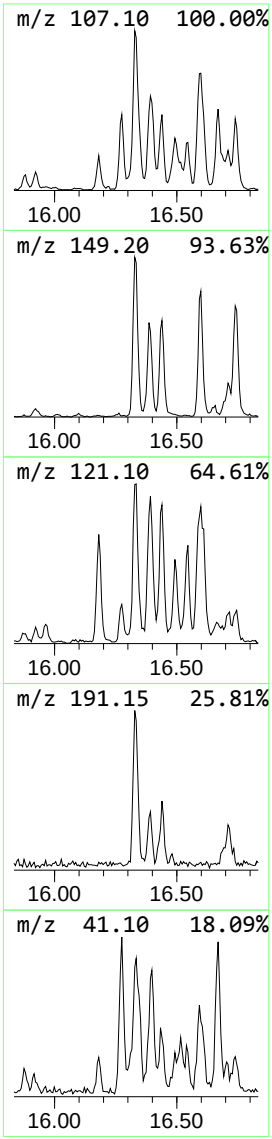
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenylpropionic acid pentaf... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.334	8.00 ng/ul	258883	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenylpropionic acid pentafluoro...	330	C16H11F5O2	062434-95-1	53
2	1-(Isocyanatomethyl)-4-propoxybe...	191	C11H13NO2	936095-07-7	46
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
4	Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	43
5	4-Isopropoxy-4-phenylbutyronitrile	203	C13H17NO	1010370-35-3	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

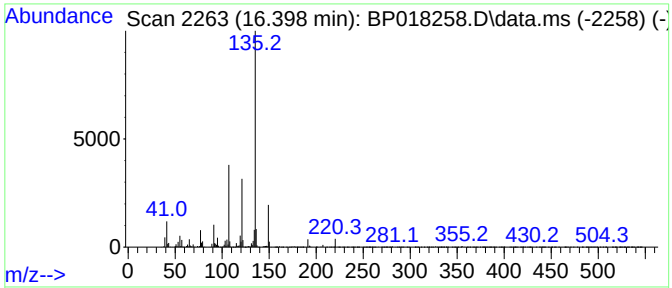
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

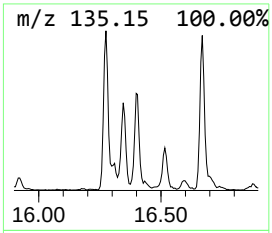
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

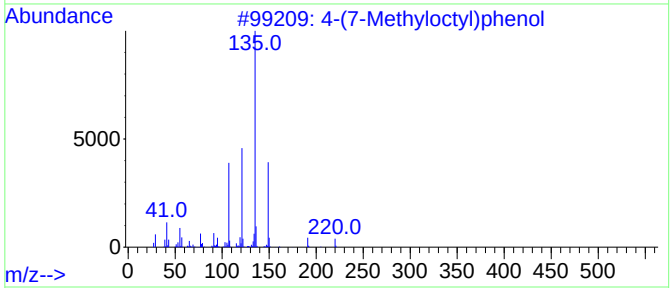
****
Peak Number 6 4-(7-Methyloctyl)phenol Concentration Rank 4

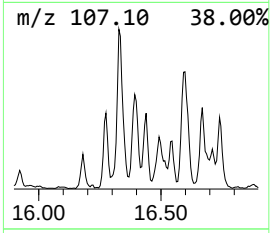
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	5.85 ng/ul	189302	Phenanthrene-d10	17.245

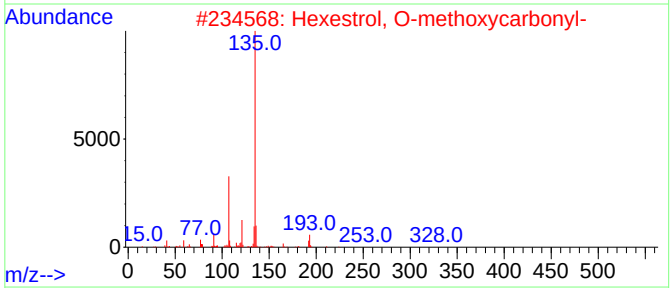
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	81
2	Hexestrol, O-methoxycarbonyl-	328	C20H24O4	1000487-66-3	64
3	meso-Hexestrol, O,O'-bis(n-propy...	442	C26H34O6	1000487-65-0	59
4	Pentafluoropropionic acid, 1-ada...	312	C14H17F5O2	1000283-03-8	53
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	53

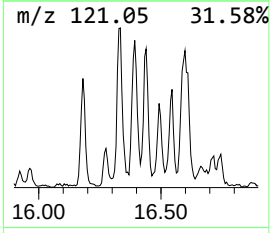


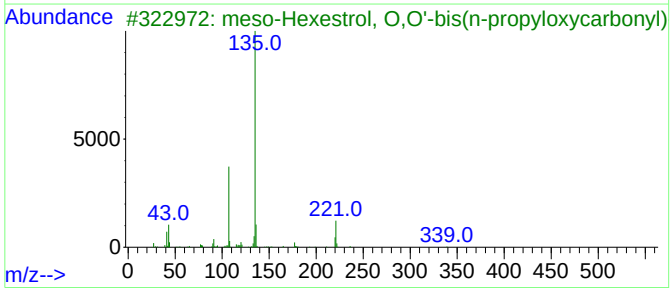


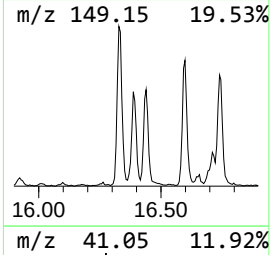


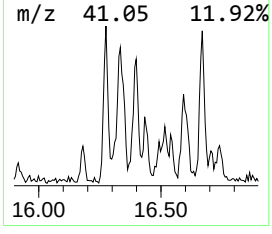












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

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

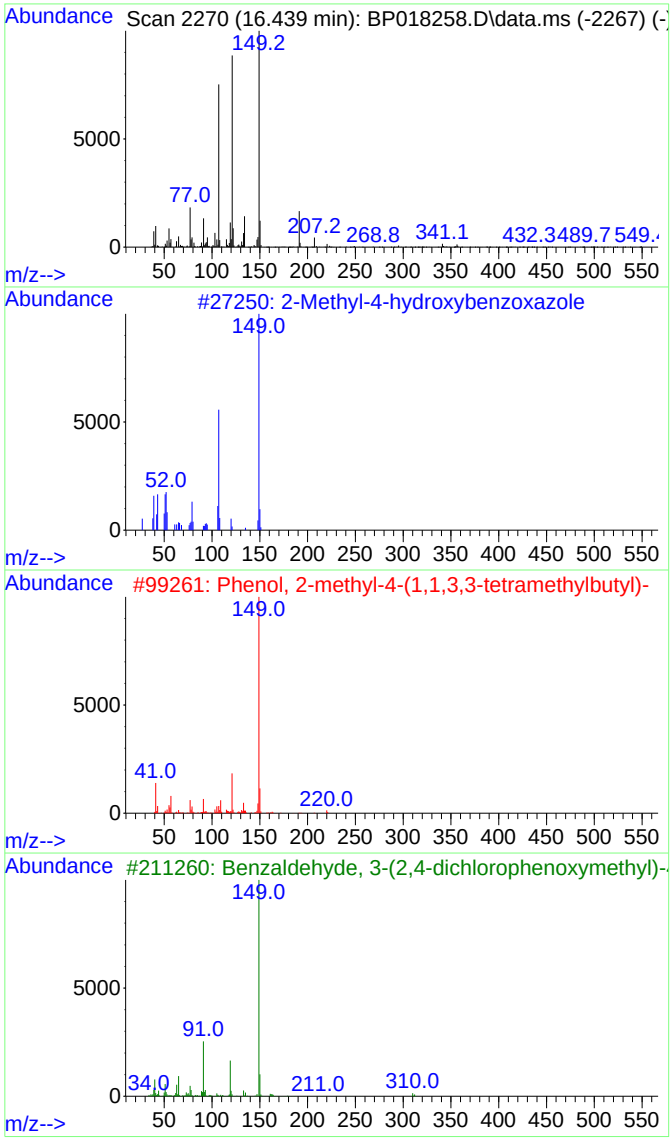
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

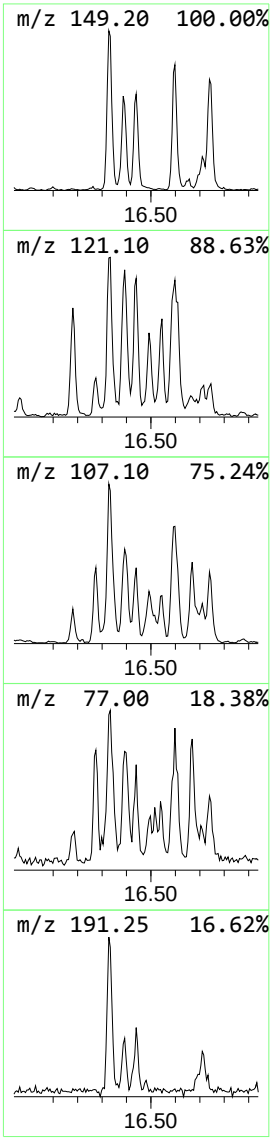
****

Peak Number 7 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.439	2.80 ng/ul	90642	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	49
2	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	25
3	Benzaldehyde, 3-(2,4-dichlorophe...	310	C15H12Cl2O3	1000273-78-9	22
4	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	22
5	Fenobucarb	207	C12H17NO2	003766-81-2	22





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Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

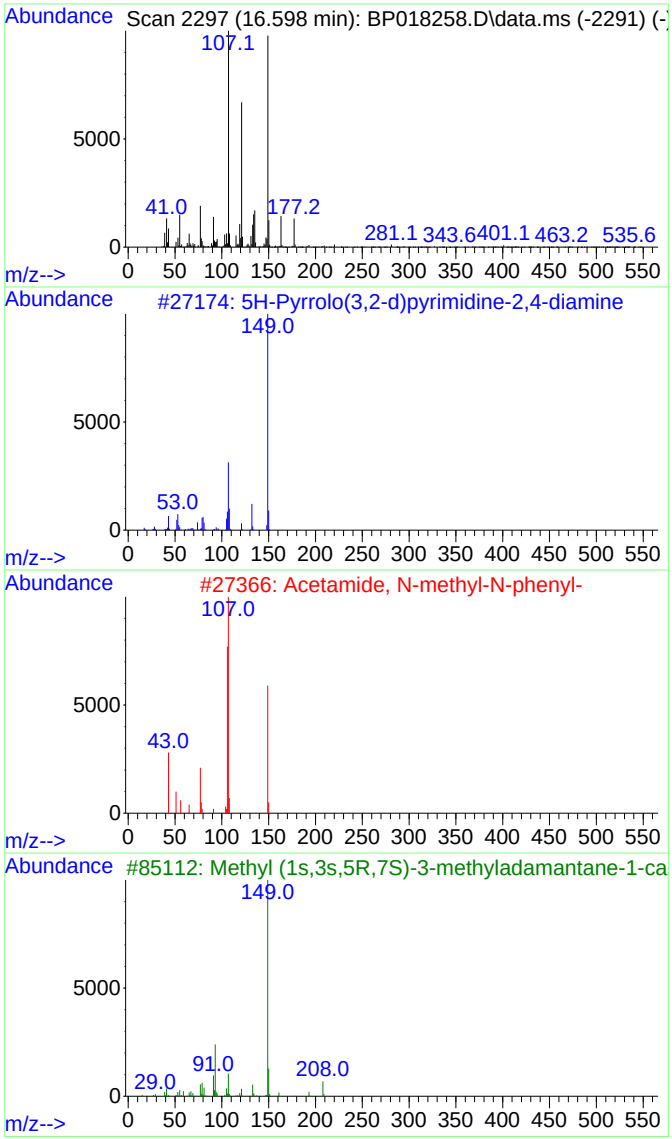
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

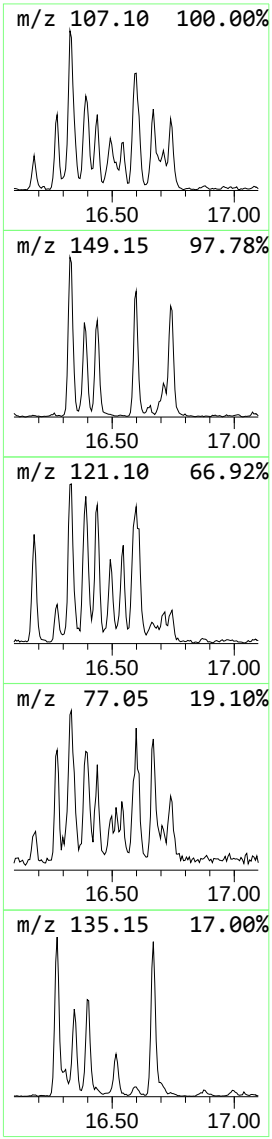
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 5H-Pyrrolo(3,2-d)pyrimidine... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	5.60 ng/ul	181064	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	64
2	Acetamide, N-methyl-N-phenyl-	149	C9H11NO	000579-10-2	46
3	Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	38
4	Benzene, 1-butyl-4-methoxy-	164	C11H16O	018272-84-9	35
5	1-Isopropenyl-3-propenylcyclopen...	150	C11H18	1000192-81-2	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

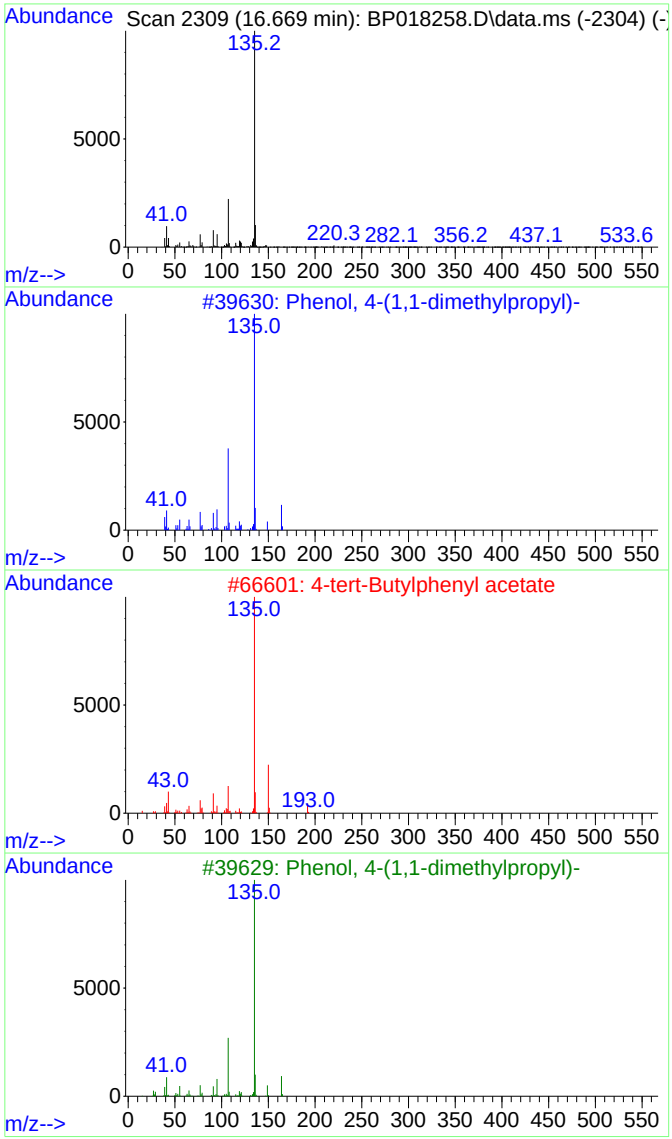
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

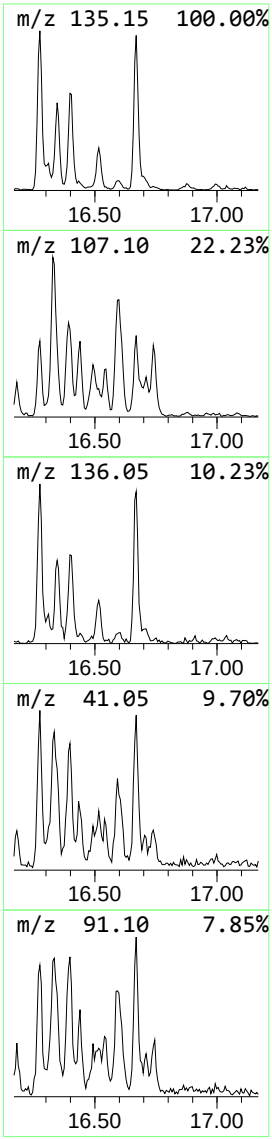
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 4-tert-Butylphenyl acetate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	4.90 ng/ul	158696	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	83
2	4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	78
5	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

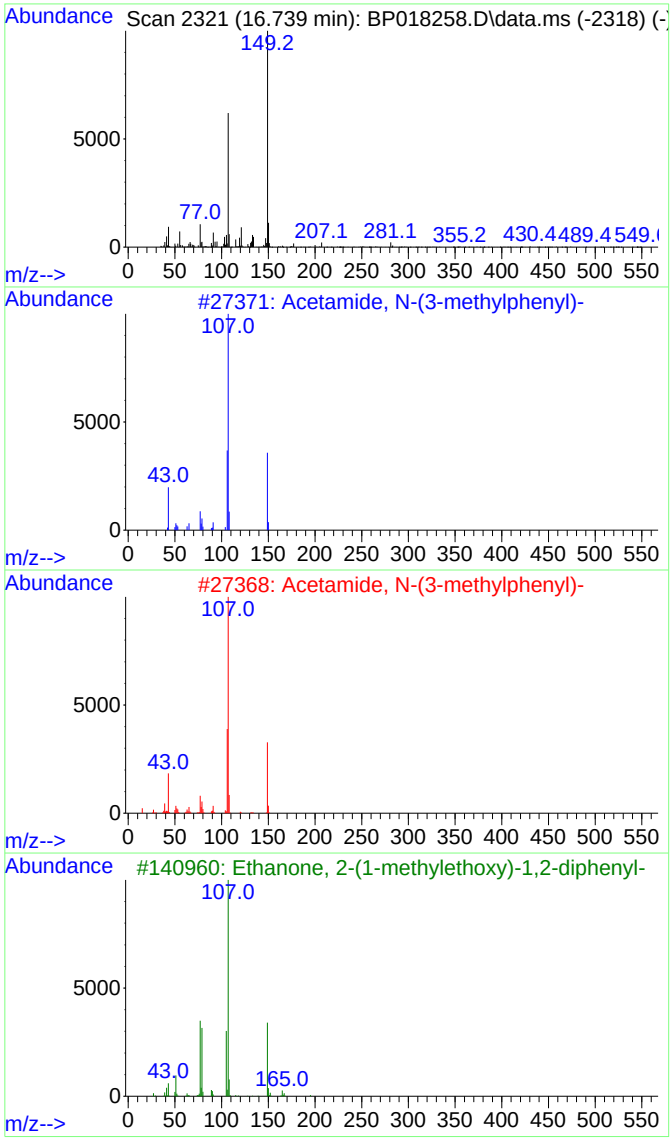
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

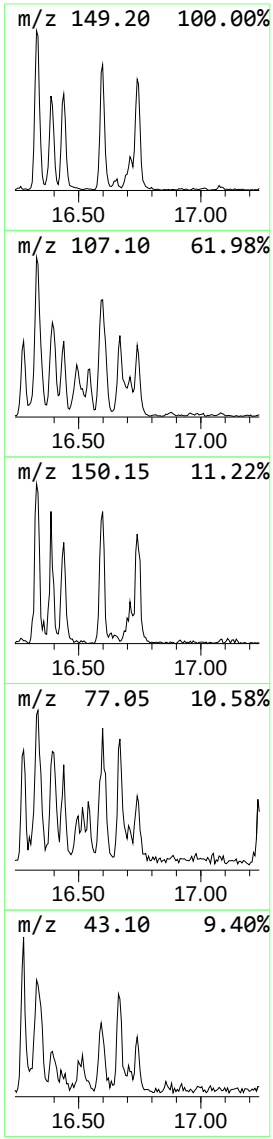
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Acetamide, N-(3-methylphenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.739	2.18 ng/ul	70674	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	72
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
3	Ethanone, 2-(1-methylethoxy)-1,2...	254	C17H18O2	006652-28-4	64
4	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	64
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018258.D
Acq On : 18 Nov 2023 13:15
Operator : MA/JU
Sample : 05370-13
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDQ7

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
Phenol, 4-(1,1-...	13.298	2.8	ng/ul	66226	3	14.457	471533	20.0
Cyclohexanamine...	13.669	6.2	ng/ul	147054	3	14.457	471533	20.0
(Bicyclohexyl)-...	13.740	10.3	ng/ul	243294	3	14.457	471533	20.0
Phenol, 4-(1,1,...	16.275	4.8	ng/ul	154586	4	17.245	647152	20.0
Phenylpropionic...	16.334	8.0	ng/ul	258883	4	17.245	647152	20.0
4-(7-Methylocty...	16.398	5.8	ng/ul	189302	4	17.245	647152	20.0
unknown-01	16.439	2.8	ng/ul	90642	4	17.245	647152	20.0
5H-Pyrrolo(3,2-...	16.598	5.6	ng/ul	181064	4	17.245	647152	20.0
4-tert-Butylphe...	16.669	4.9	ng/ul	158696	4	17.245	647152	20.0
Acetamide, N-(3...	16.739	2.2	ng/ul	70674	4	17.245	647152	20.0