LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

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: BP018209.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	17	rVB2	12611	14843	0.70%	0.070%
2	3.617	86	90	96	rBV	25274	49230	2.31%	0.231%
3	3.775	114	117	121	rVB5	6163	8949	0.42%	0.042%
4	3.899	134	138	143	rBV4	4610	8427	0.39%	0.039%
5	4.411	220	225	234	rBV	45800	69624	3.26%	0.326%
6	5.034	326	331	341	rBV2	16196	31849	1.49%	0.149%
7	5.175	351	355	356	rBV4	2826	2148	0.10%	0.010%
8	6.158	518	522	524	rVB5	1847	2390	0.11%	0.011%
9	6.787	625	629	630	rBV3	1936	2234	0.10%	0.010%
10	6.963	653	659	670	rBV2	38584	95238	4.46%	0.446%
11	7.046	670	673	679	rVV5	6130	11292	0.53%	0.053%
12	7.134	682	688	700	rVV	180328	288503	13.51%	1.352%
13	7.305	711	717	728	rBV	190396	351445	16.46%	1.647%
14	7.652	773	776	783	rVB4	3410	6487	0.30%	0.030%
15	7.781	792	798	816	rBV	262308	464695	21.77%	2.178%
16	8.040	838	842	846	rBV7	1410	2632	0.12%	0.012%
17	8.516	917	923	939	rBV	142043	304412	14.26%	1.427%
18	8.640	942	944	949	rVB6	1566	2155	0.10%	0.010%
19	8.993	997	1004	1023	rBV	240037	472759	22.14%	2.216%
20	9.704	1116	1125	1148	rBV	198516	420177	19.68%	1.969%
21	10.234	1209	1215	1242	rBV	211014	539668	25.28%	2.529%
22	10.599	1269	1277	1292	rBV	391173	653355	30.60%	3.062%
23	10.799	1305	1311	1332	rBV	76525	230809	10.81%	1.082%
24	11.416	1405	1416	1419	rBV	5491	16001	0.75%	0.075%
25	11.651	1452	1456	1460	rBV6	1342	2293	0.11%	0.011%
26	12.234	1550	1555	1556	rBV4	3128	4167	0.20%	0.020%
27	12.251	1556	1558	1562	rVB3	2525	3750	0.18%	0.018%
28	12.763	1641	1645	1648	rBV6	1719	2701	0.13%	0.013%
29	13.298	1730	1736	1747	rBV2	62124	110038	5.15%	0.516%
30	13.445	1755	1761	1762	rBV6	1568	2194	0.10%	0.010%
31	13.663	1792	1798	1801	rBV8	1976	3825	0.18%	0.018%
32	13.751	1803	1813	1830	rBV2	97317	401245	18.79%	1.881%
33	13.887	1830	1836	1853	rVB	690585	991438	46.44%	4.647%
34	14.075	1864	1868	1872	rVB7	4955	5577	0.26%	0.026%
35	14.163	1876	1883	1894	rBV	645357	959794	44.96%	4.498%
36	14.463	1928	1934	1944	rBV	567873	807235	37.81%	3.783%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

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	14.798	1983	1991	1997	rBV8	12825	39101	1.83%	0.183%
38	14.981	2018	2022	2029	rVB10	3332	5656	0.26%	0.027%
39	15.328	2078	2081	2085	rBV6	1761	2948	0.14%	0.014%
40	15.469	2098	2105	2118	rBV	858311	1303313	61.05%	6.108%
41	15.628	2125	2132	2144	rBV	270312	471523	22.09%	2.210%
42	15.881	2171	2175	2178	rBV	17335	24154	1.13%	0.113%
43	15.922	2179	2182	2186	rVV	19074	28640	1.34%	0.134%
44	15.963	2186	2189	2194	rVB6	5006	7876	0.37%	0.037%
45	16.092	2207	2211	2212	rBV4	2151	2476	0.12%	0.012%
46	16.187	2222	2227	2232	rBV	26455	38721	1.81%	0.181%
47	16.281	2236	2243	2246	rVV	93314	131636	6.17%	0.617%
48	16.339	2246	2253	2259	rVV2	117230	258916	12.13%	1.213%
49	16.398	2259	2263	2268	rVV3	95519	175410	8.22%	0.822%
50	16.445	2268	2271	2275	rVV	57032	77878	3.65%	0.365%
51	16.498	2275	2280	2282	rVV2	31499	54490	2.55%	0.255%
52	16.545	2286	2288	2292	rVV	33745	43575	2.04%	0.204%
53	16.598	2292	2297	2304	rVV4	82499	162491	7.61%	0.762%
54	16.675	2305	2310	2314	rVV	103159	159676	7.48%	0.748%
55	16.710	2314	2316	2319	rVV3	41853	59496	2.79%	0.279%
56	16.745	2319	2322	2329	rVV2	57169	79426	3.72%	0.372%
57	16.804	2330	2332	2338	rVB7	4973	7441	0.35%	0.035%
58	16.963	2357	2359	2361	rBV3	1884	2167	0.10%	0.010%
59	17.004	2364	2366	2368	rVB3	3885	3100	0.15%	0.015%
60	17.081	2376	2379	2382	rVV5	3285	3432	0.16%	0.016%
61	17.122	2383	2386	2390	rVB6	4045	6663	0.31%	0.031%
62	17.245	2398	2407	2416	rBV	800999	1070422	50.14%	5.017%
63	17.345	2418	2424	2442	rVV	1098085	1610034	75.41%	7.546%
64	17.498	2447	2450	2453	rBV5	2203	3272	0.15%	0.015%
65	17.528	2453	2455	2459	rVB5	5176	6157	0.29%	0.029%
66	17.581	2459	2464	2468	rBV8	5572	9727	0.46%	0.046%
67	17.616	2468	2470	2475	rVB6	3452	5003	0.23%	0.023%
68	17.675	2475	2480	2482	rBV6	4925	8338	0.39%	0.039%
69	17.751	2491	2493	2496	rVB4	3012	3077	0.14%	0.014%
70	17.851	2504	2510	2511	rBV	29478	39954	1.87%	0.187%
71	17.881	2511	2515	2525	rVB	841457	1016486	47.61%	4.764%
72	18.039	2539	2542	2547	rBV7	8042	14646	0.69%	0.069%
73	18.098	2548	2552	2564	rVV2	52731	91327	4.28%	0.428%
74	18.192	2564	2568	2573	rVB3	24747	34310	1.61%	0.161%
75	18.310	2584	2588	2593	rVB	61554	78476	3.68%	0.368%
76	18.680	2647	2651	2656	rVB8	7610	11497	0.54%	0.054%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

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	18.998	2701	2705	2711	rBV6	17899	31361	1.47%	0.147%
78	19.175	2731	2735	2739	rBV3	15953	22354	1.05%	0.105%
79	19.233	2740	2745	2746	rBV3	19313	25165	1.18%	0.118%
80	19.345	2760	2764	2769	rBV2	36002	47524	2.23%	0.223%
81	19.475	2784	2786	2790	rVB4	7503	8574	0.40%	0.040%
82	19.604	2802	2808	2820	rBV	1636242	2072316	97.07%	9.712%
83	21.380	3105	3110	3119	rBV	887211	1214695	56.90%	5.693%
84	23.686	3495	3502	3514	rVB2	1093185	2134942	100.00%	10.006%
85	23.851	3524	3530	3543	rVB	591806	1283459	60.12%	6.015%

Sum of corrected areas: 21336900

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
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Acq On : 16 Nov 2023 12:26
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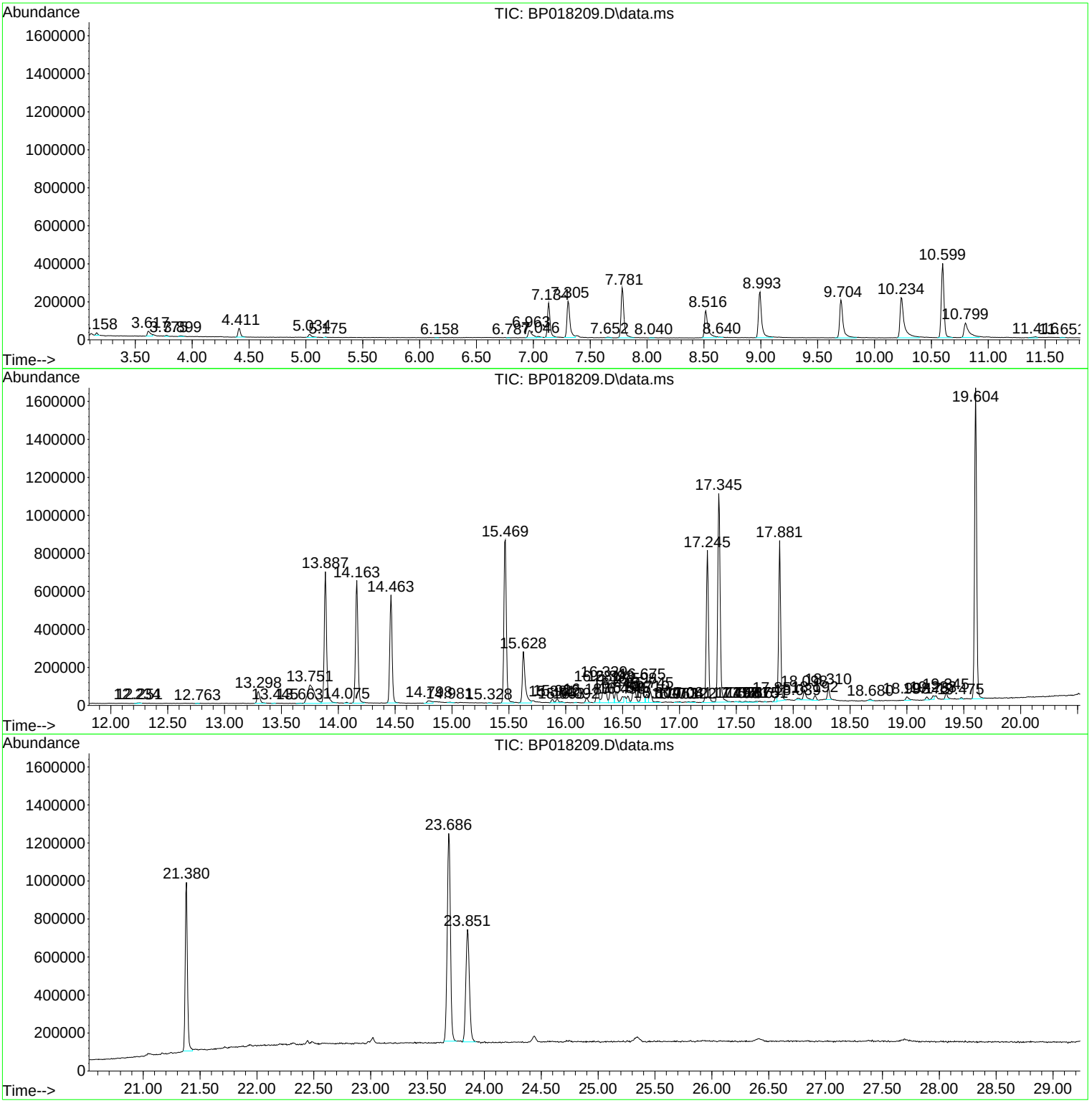
Instrument :
BNA_P
ClientSampleId :
GCDL9

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 : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

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

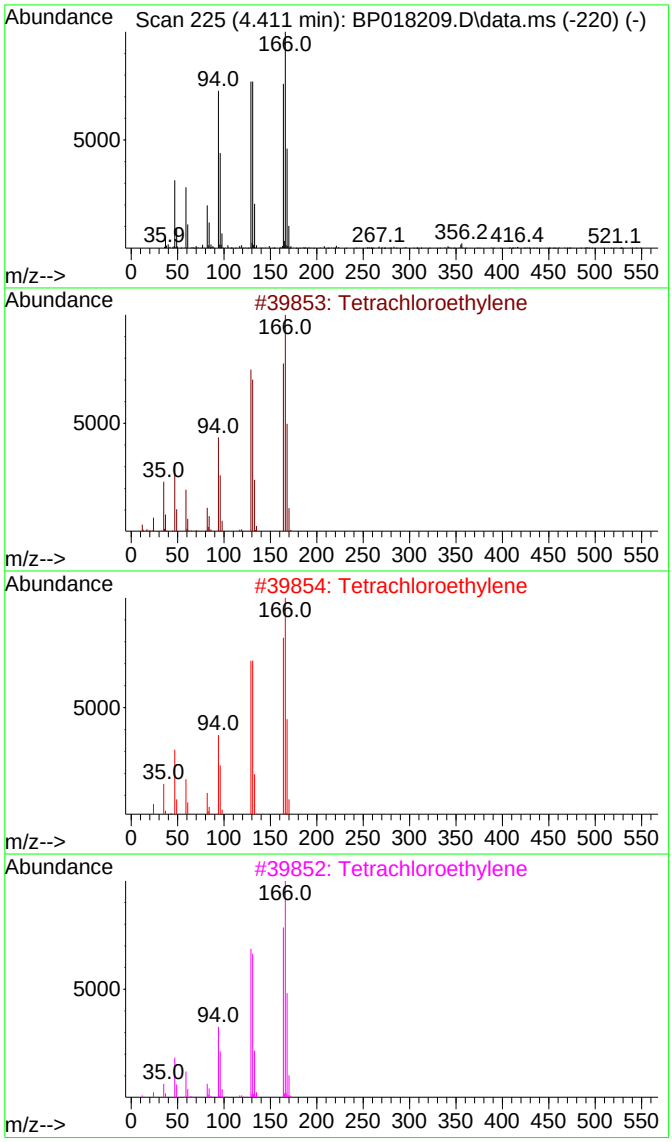
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

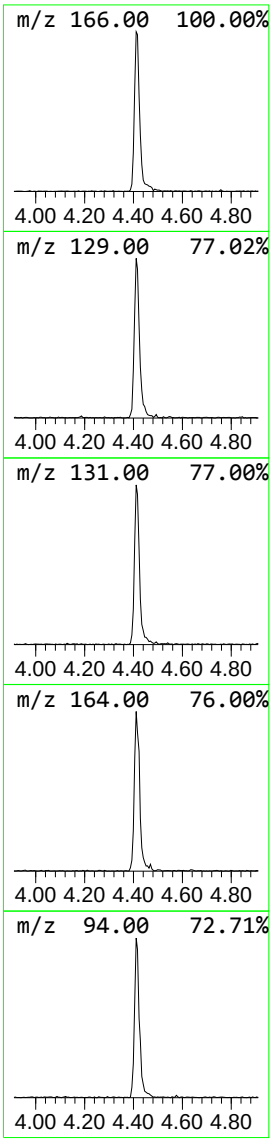
****

Peak Number 1 Tetrachloroethylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.411	3.00 ng/ul	69624	1,4-Dichlorobenzene-d4	7.781

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrachloroethylene	164	C2Cl4	000127-18-4	95
2	Tetrachloroethylene	164	C2Cl4	000127-18-4	95
3	Tetrachloroethylene	164	C2Cl4	000127-18-4	94
4	Tetrachloroethylene	164	C2Cl4	000127-18-4	94
5	Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HCl2FN2	002927-71-1	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

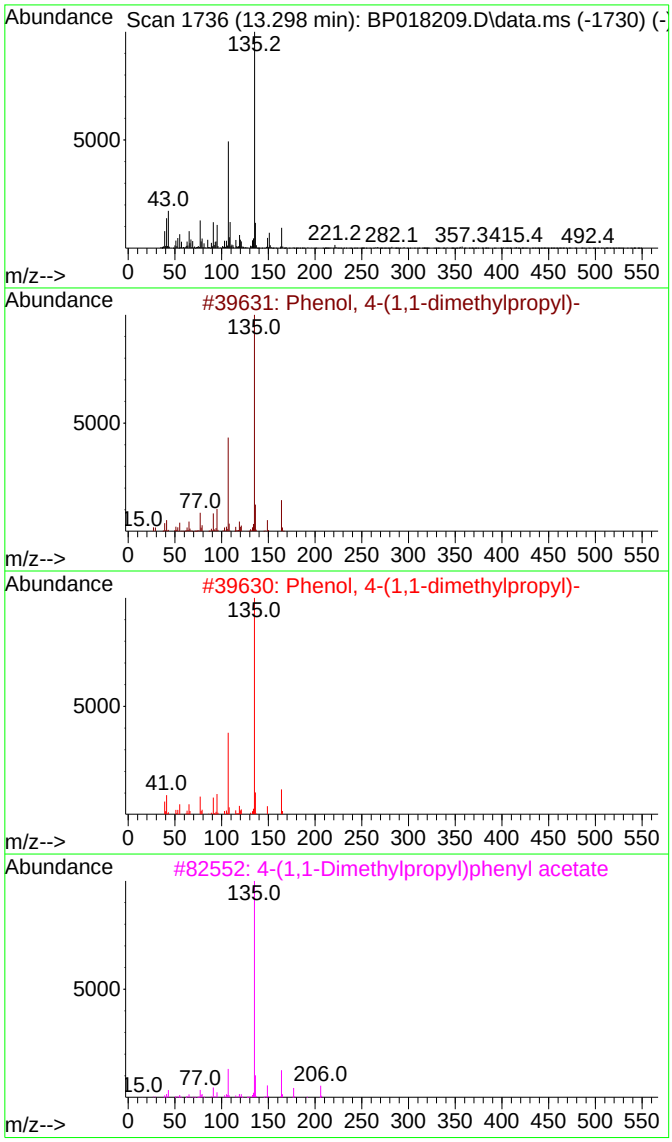
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

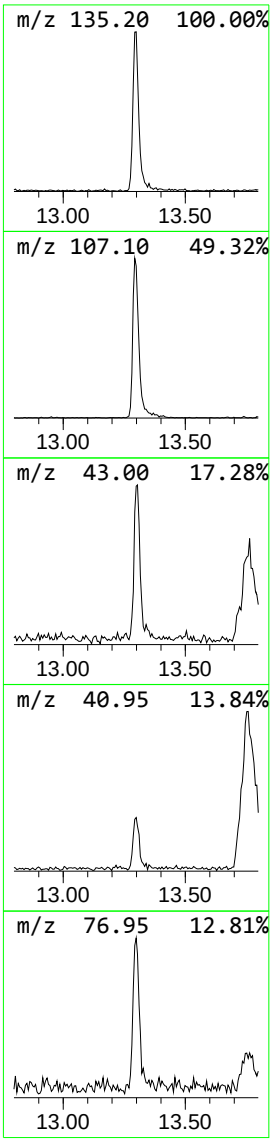
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenol, 4-(1,1-dimethylprop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.298	2.73 ng/ul	110038	Acenaphthene-d10	14.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	91
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	81
3	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	72
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

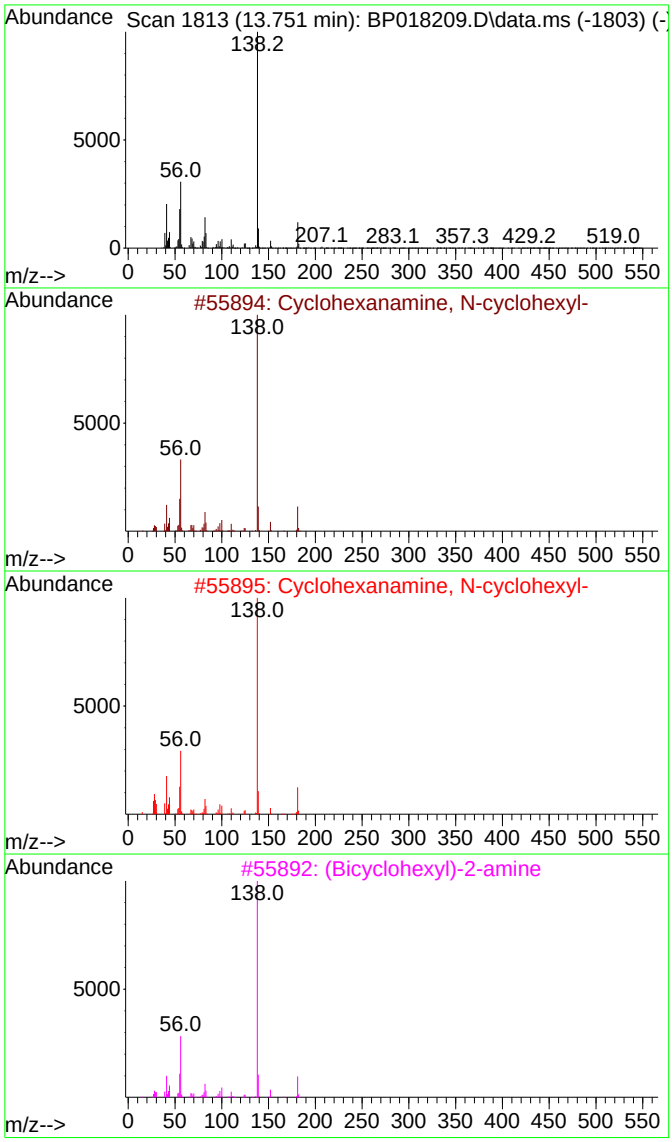
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

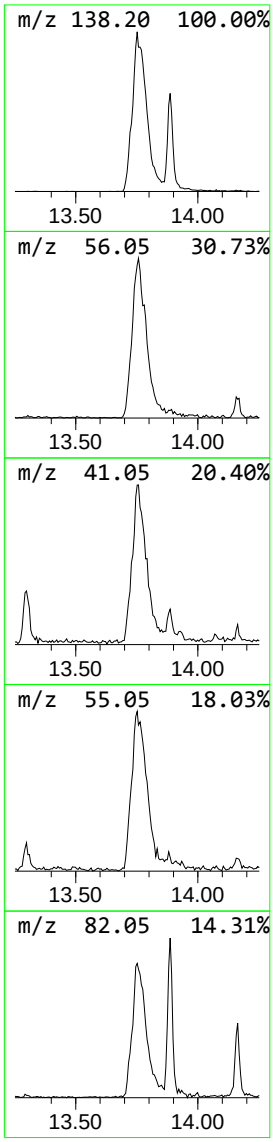
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	9.94 ng/ul	401245	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	94
3	(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	91
4	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	90
5	Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

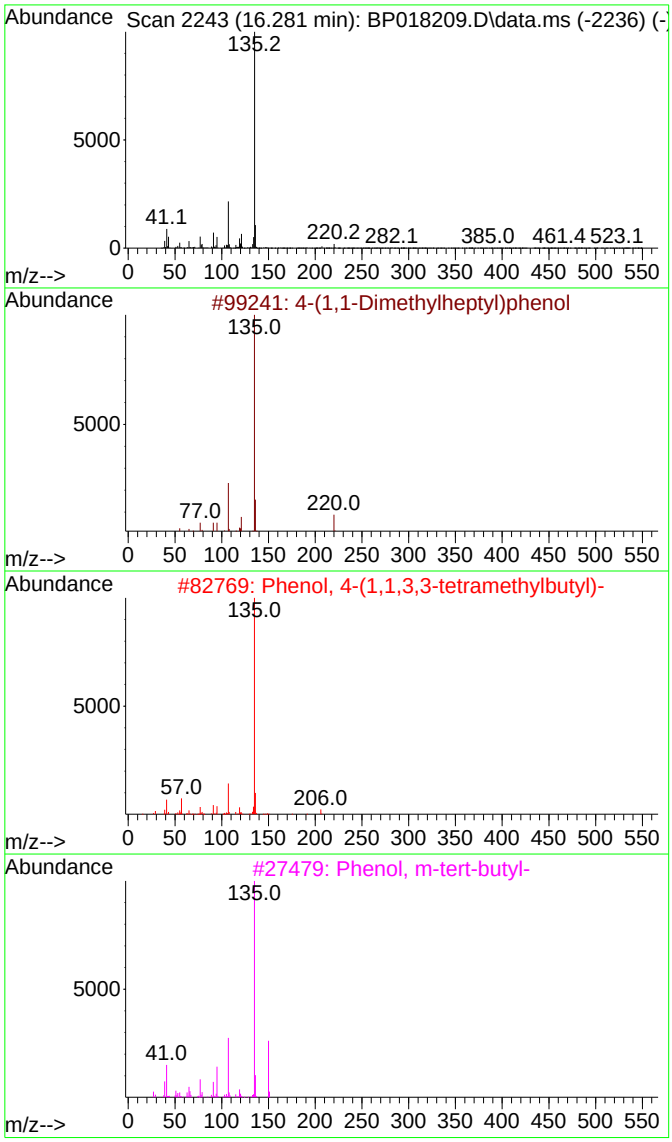
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

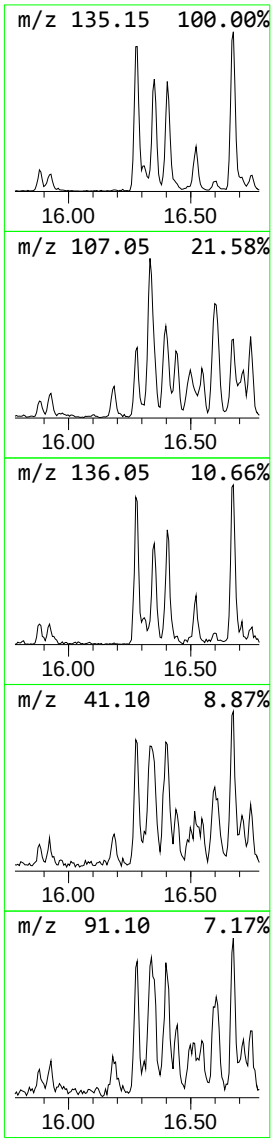
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 4-(1,1-Dimethylheptyl)phenol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.281	2.46 ng/ul	131636	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	90
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	80
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
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 : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

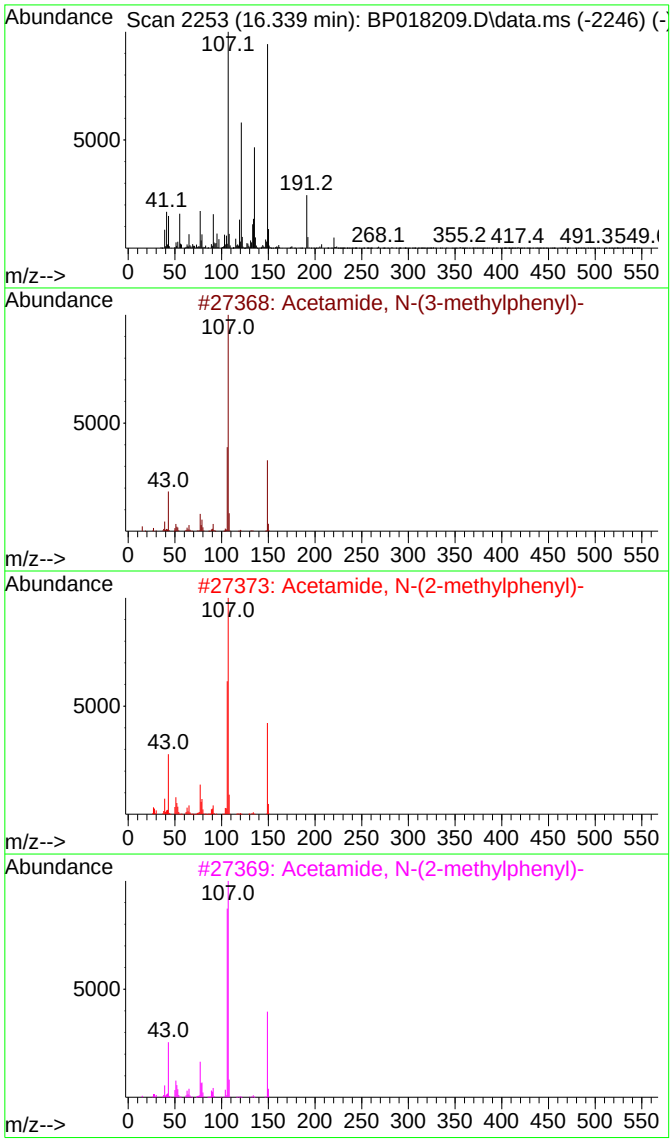
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

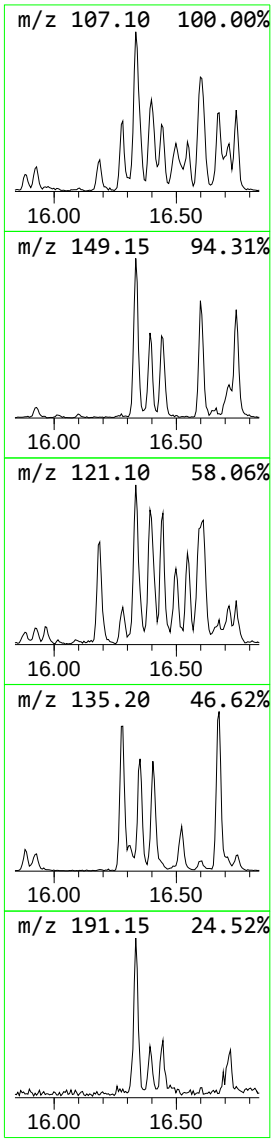
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Acetamide, N-(3-methylphenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.339	4.84 ng/ul	258916	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	53
2	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38
3	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38
4	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	38
5	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

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

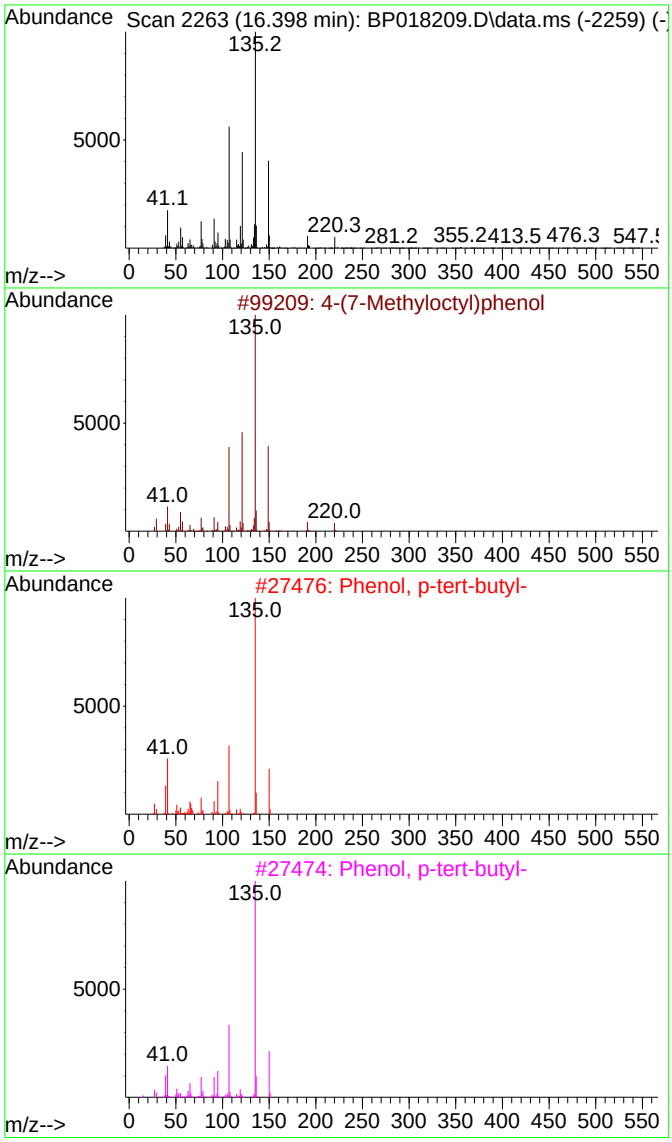
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

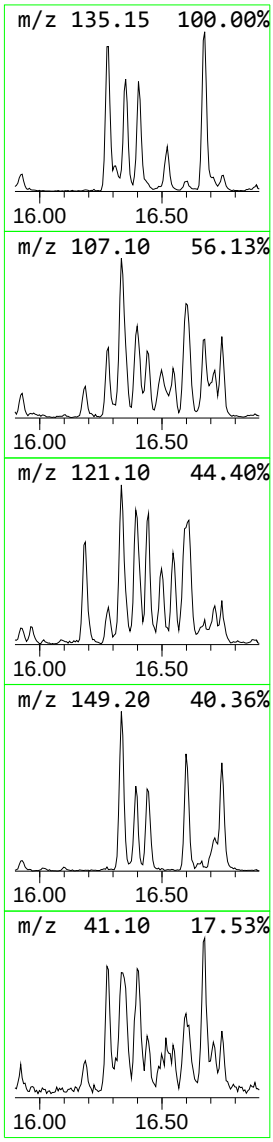
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	3.28 ng/ul	175410	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	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	52
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50





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Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

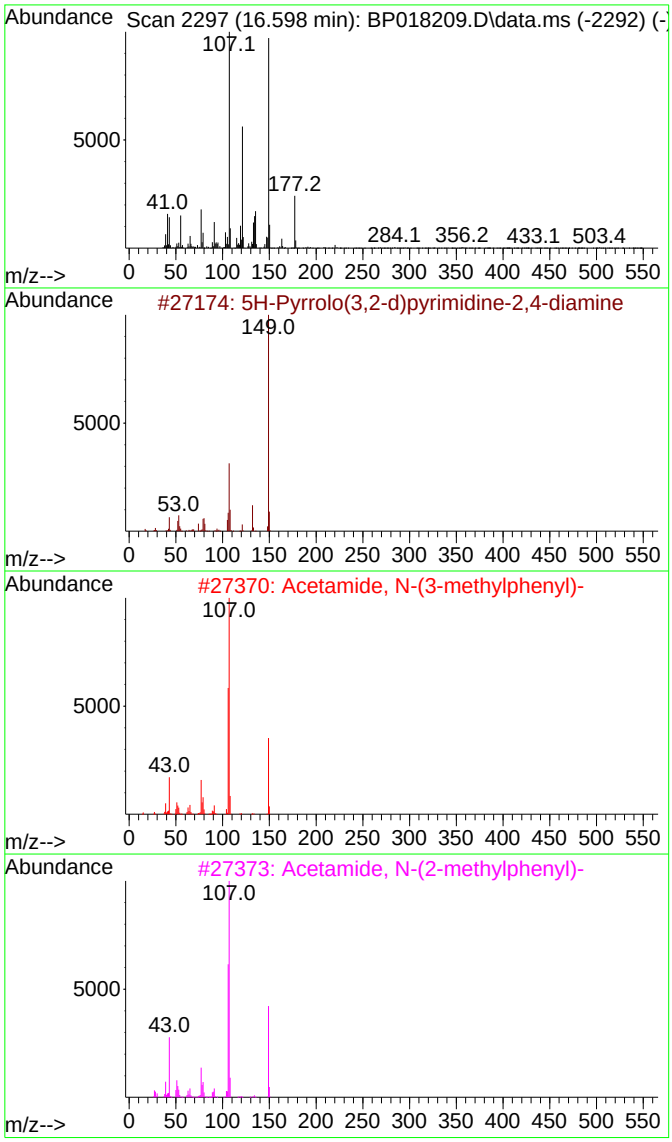
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

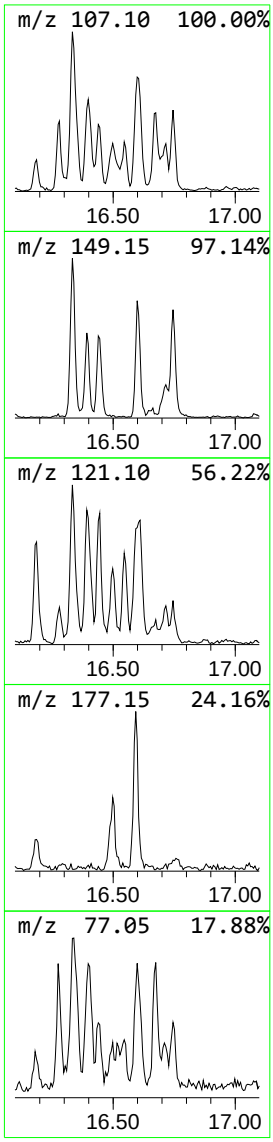
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 5H-Pyrrolo(3,2-d)pyrimidine... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	3.04 ng/ul	162491	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-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	58
4	Benzene, [1-(1-methylethoxy)-3-b...	190	C13H18O	098088-47-2	53
5	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

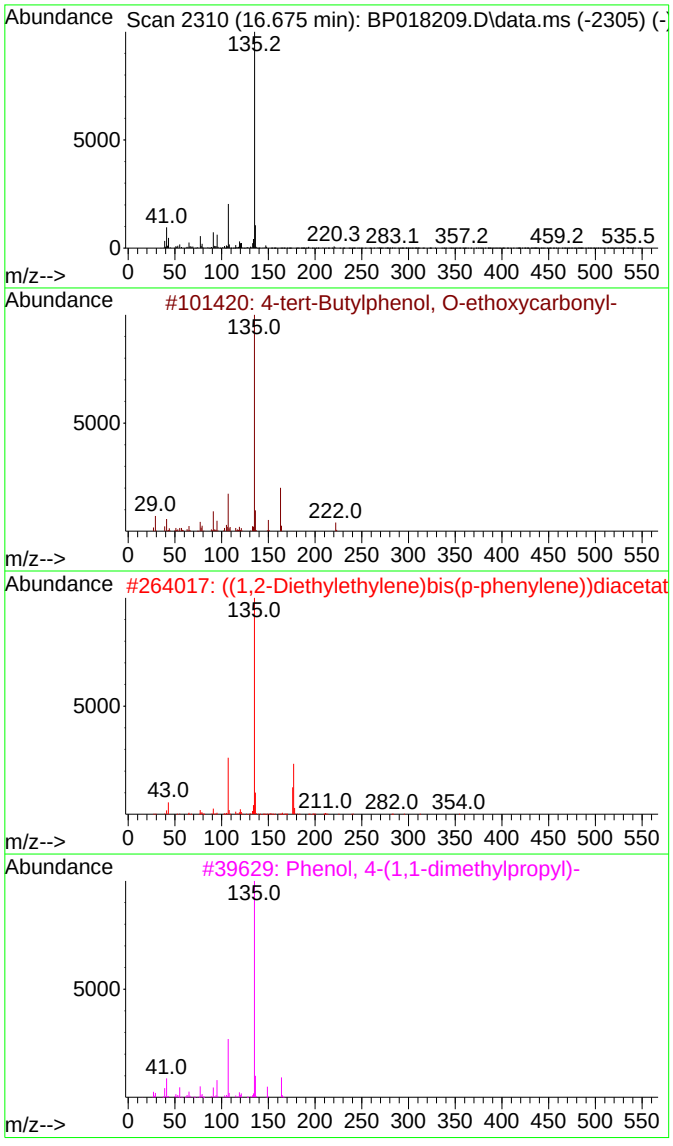
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

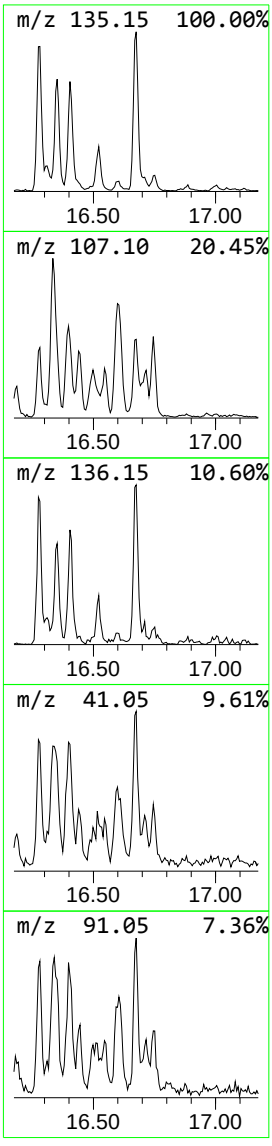
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 4-tert-Butylphenol, O-ethox... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.675	2.98 ng/ul	159676	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-tert-Butylphenol, O-ethoxycarb...	222	C13H18O3	026177-66-2	83
2	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	78
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	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 : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

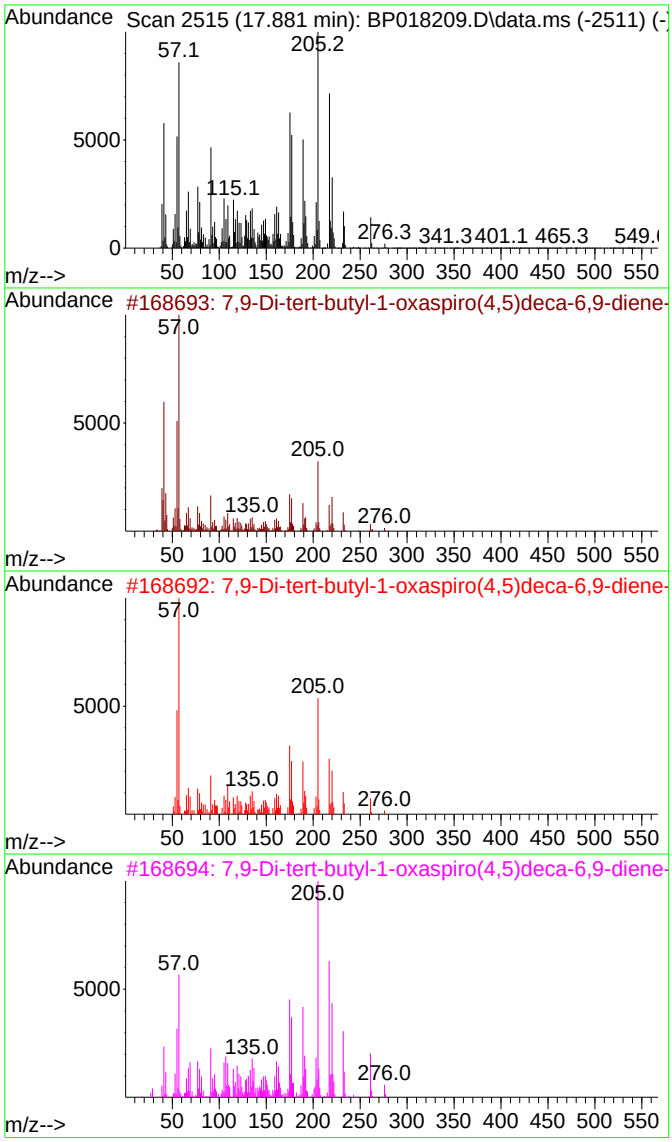
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

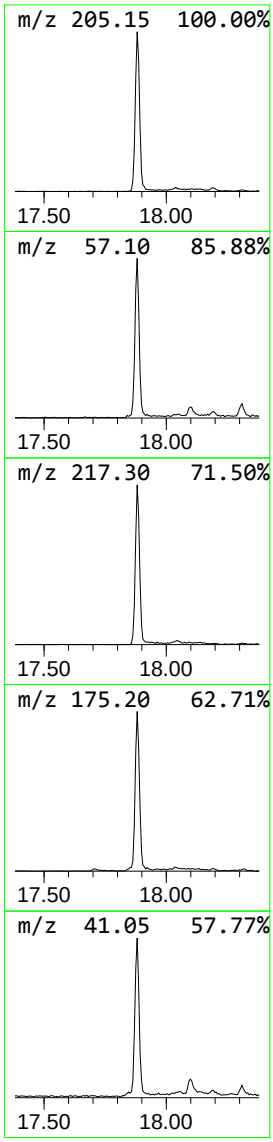
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.881	18.99 ng/ul	1016490	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	93
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	86
4	Benzenamine, N,N-diethyl-4-(2-ni...	220	C12H16N2O2	064462-51-7	52
5	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018209.D
Acq On : 16 Nov 2023 12:26
Operator : MA/JU
Sample : 05338-14
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDL9

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	3.0	ng/ul	69624	1	7.781	464695	20.0
Phenol, 4-(1,1-...	13.298	2.7	ng/ul	110038	3	14.463	807235	20.0
Cyclohexanamine...	13.751	9.9	ng/ul	401245	3	14.463	807235	20.0
4-(1,1-Dimethyl...	16.281	2.5	ng/ul	131636	4	17.245	1070420	20.0
Acetamide, N-(3...	16.339	4.8	ng/ul	258916	4	17.245	1070420	20.0
4-(7-Methylocty...	16.398	3.3	ng/ul	175410	4	17.245	1070420	20.0
5H-Pyrrolo(3,2-...	16.598	3.0	ng/ul	162491	4	17.245	1070420	20.0
4-tert-Butylphe...	16.675	3.0	ng/ul	159676	4	17.245	1070420	20.0
7,9-Di-tert-but...	17.881	19.0	ng/ul	1016490	4	17.245	1070420	20.0