LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCAG9

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-BP030723.M
Title : SVOA CALIBRATION

Signal : TIC: BP013972.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.440	39	43	49	rBV	38454	52077	0.48%	0.030%
2	3.493	49	52	59	rVB3	11202	15522	0.14%	0.009%
3	3.605	65	71	84	rVB2	27734	50617	0.46%	0.029%
4	3.793	98	103	113	rVB	13427	21682	0.20%	0.012%
5	3.875	113	117	136	rBV	466852	813255	7.43%	0.469%
6	4.111	152	157	161	rBV	11541	16380	0.15%	0.009%
7	4.211	169	174	182	rBV2	16367	31827	0.29%	0.018%
8	4.487	215	221	227	rBV	20768	39975	0.37%	0.023%
9	4.546	227	231	235	rVB	12157	16844	0.15%	0.010%
10	5.158	327	335	345	rBV2	105028	170780	1.56%	0.098%
11	5.364	365	370	380	rVB	89962	144810	1.32%	0.083%
12	6.258	516	522	527	rBV	26315	40875	0.37%	0.024%
13	6.463	550	557	562	rBV2	9256	15673	0.14%	0.009%
14	6.705	591	598	601	rBV2	9153	15552	0.14%	0.009%
15	7.240	678	689	706	rBV	701215	1217187	11.13%	0.701%
16	7.410	712	718	729	rBV	536872	825579	7.55%	0.476%
17	7.616	746	753	756	rBV	715674	1220936	11.16%	0.704%
18	7.652	756	759	769	rVB	908109	1333385	12.19%	0.768%
19	8.093	827	834	846	rVB	491414	795326	7.27%	0.458%
20	8.616	912	923	930	rBV2	730482	1849265	16.91%	1.066%
21	8.675	930	933	940	rVB	87191	125183	1.14%	0.072%
22	8.799	947	954	968	rBV	946185	1541726	14.09%	0.888%
23	8.922	968	975	987	rVB	1213433	1933715	17.68%	1.114%
24	9.181	1012	1019	1026	rBV	449499	714214	6.53%	0.412%
25	9.269	1026	1034	1037	rBV	680309	1211857	11.08%	0.698%
26	9.310	1037	1041	1049	rVB	1297213	2068698	18.91%	1.192%
27	9.569	1077	1085	1095	rBV	1756981	2718082	24.85%	1.566%
28	9.675	1095	1103	1108	rVB9	7343	16714	0.15%	0.010%
29	9.993	1148	1157	1159	rBV	646469	1029702	9.41%	0.593%
30	10.028	1159	1163	1168	rVV	937377	1611139	14.73%	0.928%
31	10.081	1168	1172	1180	rVB2	52696	113115	1.03%	0.065%
32	10.534	1239	1249	1252	rBV	980591	1888298	17.26%	1.088%
33	10.722	1275	1281	1287	rVB	22574	34318	0.31%	0.020%
34	10.804	1287	1295	1306	rBV2	53283	104362	0.95%	0.060%
35	10.916	1306	1314	1319	rVV	683076	1196952	10.94%	0.690%
36	10.969	1319	1323	1330	rVB	613417	992673	9.07%	0.572%

LSC Area Percent Report

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Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

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-BP030723.M
Title : SVOA CALIBRATION

37	11.057	1332	1338	1354	rBV	125132	269985	2.47%	0.156%
38	11.616	1425	1433	1442	rBV	53944	91108	0.83%	0.053%
39	12.193	1523	1531	1548	rBV	1674171	2702333	24.70%	1.557%
40	12.434	1564	1572	1580	rBV	1685431	2708239	24.76%	1.561%
41	12.787	1613	1632	1641	rBV	2299356	3591162	32.83%	2.070%
42	12.928	1650	1656	1665	rVB	922630	1387148	12.68%	0.799%
43	13.169	1688	1697	1702	rBV	1292500	1937062	17.71%	1.116%
44	13.216	1702	1705	1715	rVB	50440	81288	0.74%	0.047%
45	13.569	1753	1765	1772	rBV	2420394	3531580	32.29%	2.035%
46	13.816	1798	1807	1819	rBV	1178163	1802105	16.47%	1.039%
47	13.916	1819	1824	1834	rVB	64488	92924	0.85%	0.054%
48	14.134	1855	1861	1869	rBV	1808920	2548685	23.30%	1.469%
49	14.222	1870	1876	1884	rVV	1173936	1696220	15.51%	0.978%
50	14.304	1884	1890	1900	rVV	1422327	2117203	19.36%	1.220%
51	14.428	1904	1911	1913	rVV	1852823	2791298	25.52%	1.609%
52	14.457	1913	1916	1930	rVB	2823926	4010061	36.66%	2.311%
53	14.734	1952	1963	1969	rBV	1119344	1670125	15.27%	0.962%
54	14.792	1969	1973	1983	rVB	272464	410776	3.76%	0.237%
55	14.945	1988	1999	2017	rBV3	1938494	3896972	35.63%	2.246%
56	15.092	2018	2024	2027	rVV	1689901	2425304	22.17%	1.398%
57	15.128	2027	2030	2043	rVB	1165686	1697938	15.52%	0.978%
58	15.375	2063	2072	2082	rBV	371644	551239	5.04%	0.318%
59	15.486	2087	2091	2095	rVV7	11658	17261	0.16%	0.010%
60	15.545	2095	2101	2107	rVV	2810247	3751422	34.30%	2.162%
61	15.728	2122	2132	2137	rBV	2310534	3459390	31.63%	1.994%
62	15.781	2137	2141	2146	rVV	573604	795309	7.27%	0.458%
63	15.839	2146	2151	2162	rVV	931950	1321738	12.08%	0.762%
64	15.986	2165	2176	2189	rVB	3464282	4976890	45.50%	2.868%
65	16.222	2211	2216	2226	rBV	118933	240822	2.20%	0.139%
66	16.357	2235	2239	2245	rVB	36048	50394	0.46%	0.029%
67	16.586	2274	2278	2282	rBV	21053	28115	0.26%	0.016%
68	16.657	2282	2290	2295	rVV2	48524	80409	0.74%	0.046%
69	16.710	2295	2299	2300	rVV	24804	28956	0.26%	0.017%
70	16.745	2300	2305	2308	rVV	579608	845558	7.73%	0.487%
71	16.786	2308	2312	2328	rVB	2539156	3506742	32.06%	2.021%
72	17.133	2364	2371	2387	rBV	2018495	2854099	26.09%	1.645%
73	17.257	2389	2392	2398	rVV8	10761	25300	0.23%	0.015%
74	17.486	2421	2431	2435	rBV	1542466	2327853	21.28%	1.341%
75	17.533	2435	2439	2443	rVV	2622988	3697638	33.80%	2.131%
76	17.586	2443	2448	2456	rVB2	2787948	3860951	35.30%	2.225%

LSC Area Percent Report

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Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCAG9

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
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Title : SVOA CALIBRATION

77	17.816	2483	2487	2492	rVB2	19748	25139	0.23%	0.014%
78	18.086	2527	2533	2538	rBV	159449	204934	1.87%	0.118%
79	18.416	2584	2589	2601	rBV	3495547	4103278	37.51%	2.365%
80	18.692	2633	2636	2643	rVB9	14894	21963	0.20%	0.013%
81	19.380	2749	2753	2760	rBV2	40466	63575	0.58%	0.037%
82	19.457	2761	2766	2769	rVV	36760	53555	0.49%	0.031%
83	19.675	2798	2803	2808	rBV	573324	729202	6.67%	0.420%
84	19.810	2820	2826	2829	rBV	3983223	5869076	53.65%	3.382%
85	19.833	2829	2830	2837	rVB	2137816	2017868	18.45%	1.163%
86	20.022	2857	2862	2869	rBV	216055	291734	2.67%	0.168%
87	20.169	2884	2887	2891	rBV2	31760	33782	0.31%	0.019%
88	20.798	2990	2994	3002	rBV	282925	358228	3.27%	0.206%
89	20.957	3018	3021	3026	rBV5	62428	76322	0.70%	0.044%
90	21.551	3116	3122	3133	rBV2	6070435	10938610	100.00%	6.304%
91	21.839	3167	3171	3180	rBV	273349	396416	3.62%	0.228%
92	22.004	3194	3199	3205	rBV	6488642	8377852	76.59%	4.828%
93	22.151	3219	3224	3236	rVB	595588	970193	8.87%	0.559%
94	22.416	3264	3269	3279	rVB	3010048	4045035	36.98%	2.331%
95	23.327	3418	3424	3435	rVB	3623218	6696458	61.22%	3.859%
96	23.951	3522	3530	3534	rBV2	3063037	6492134	59.35%	3.741%
97	23.998	3534	3538	3549	rVB	1930845	3936290	35.99%	2.268%
98	24.115	3551	3558	3568	rVB2	1646331	3480212	31.82%	2.006%
99	26.827	4007	4019	4031	rVB	3062778	9636641	88.10%	5.553%
100	27.627	4145	4155	4171	rVB	1405027	4839819	44.25%	2.789%

Sum of corrected areas: 173526213

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
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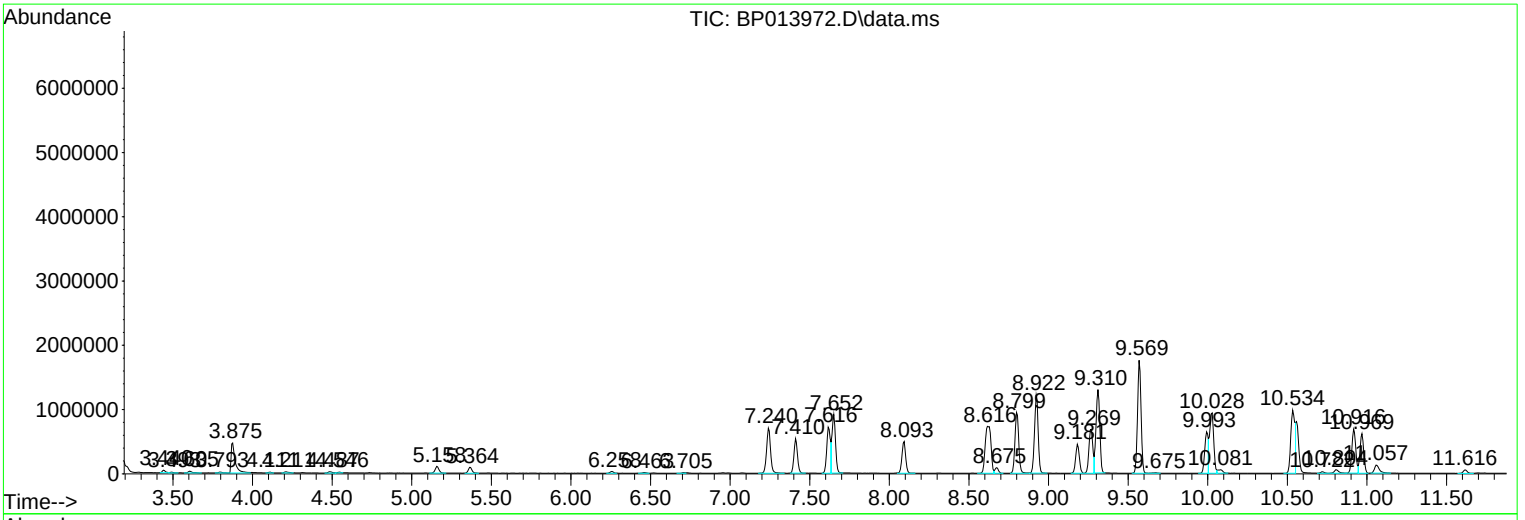
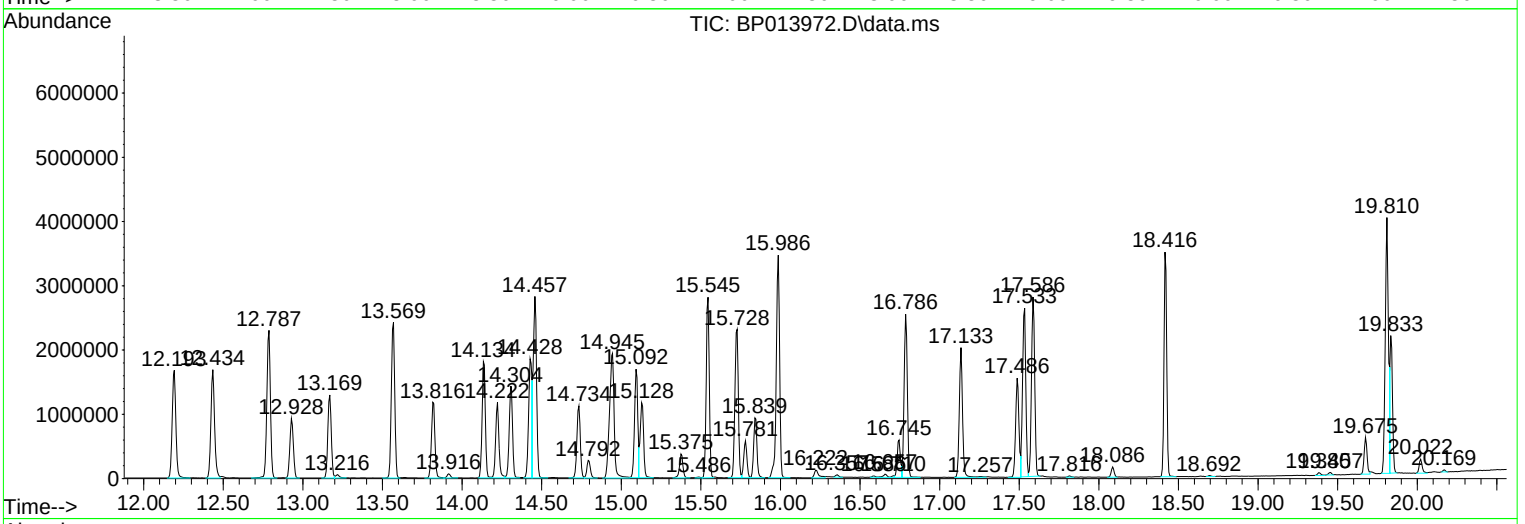
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

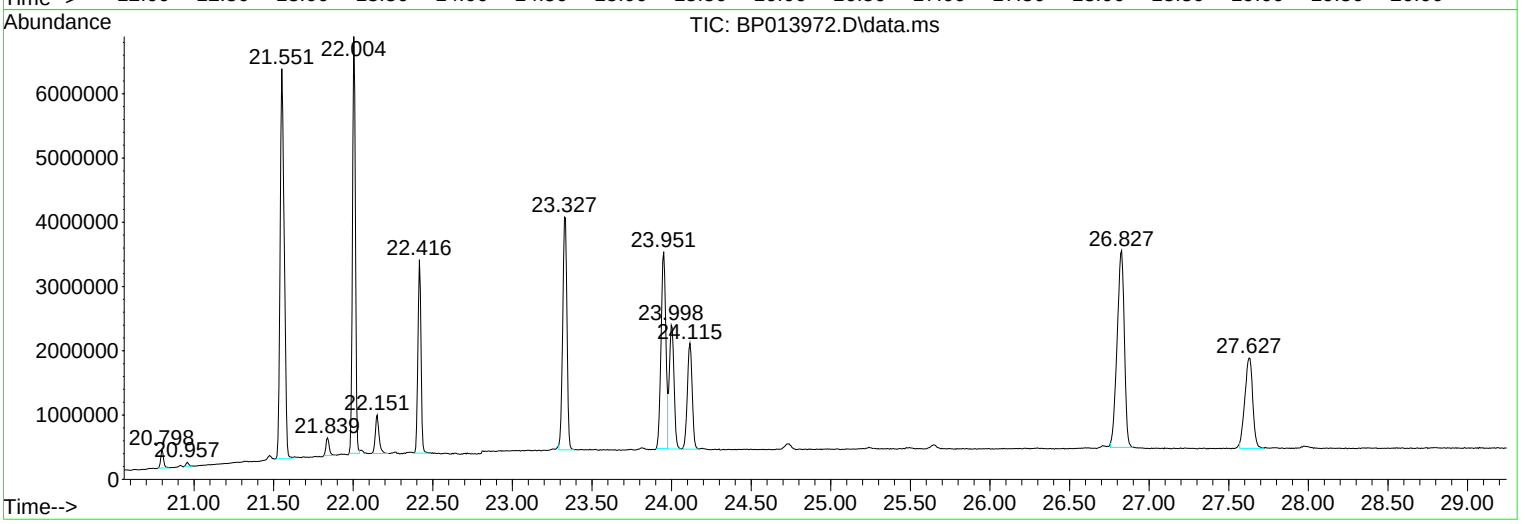
Instrument :
BNA_P
ClientSampleId :
DCAG9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
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Instrument :
BNA_P
ClientSampleId :
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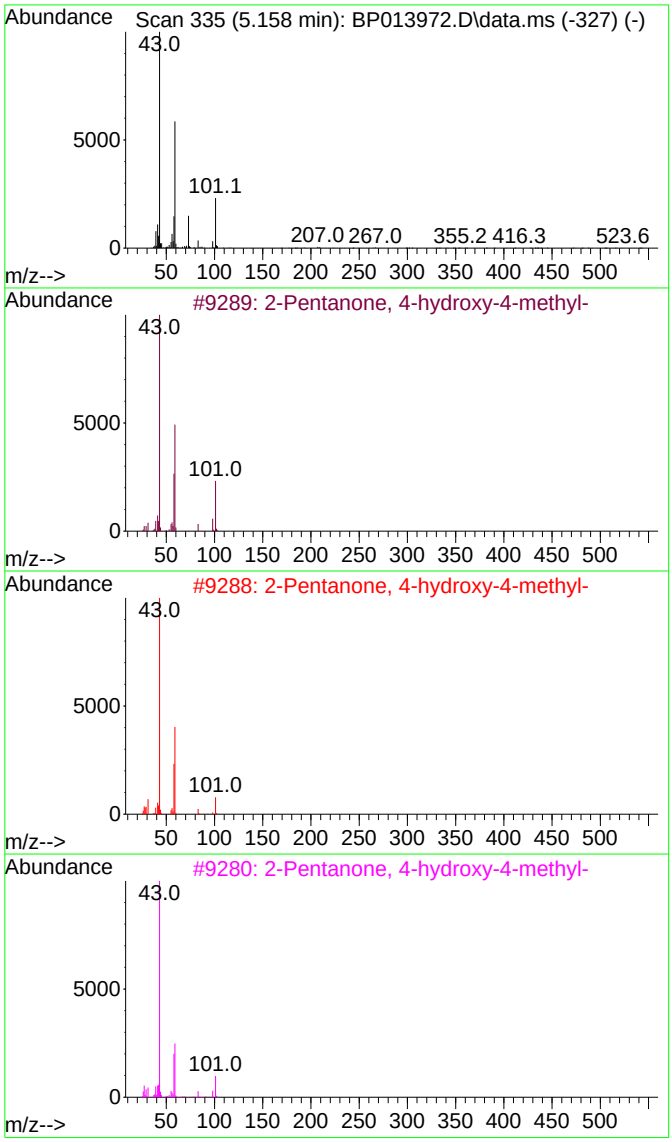
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

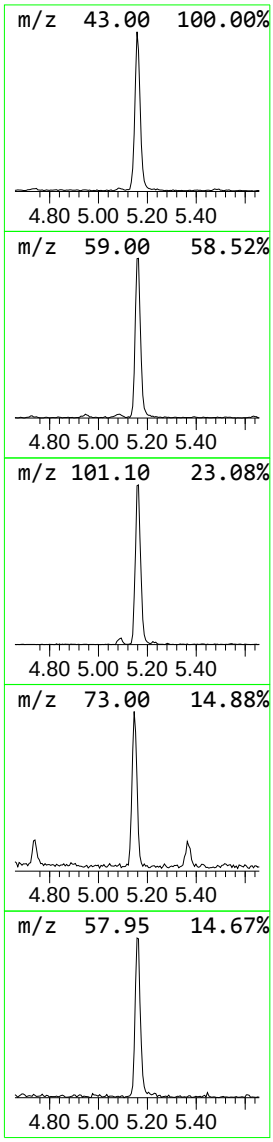
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.158	4.29 ng/ul	170780	1,4-Dichlorobenzene-d4	8.093

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
5	.alpha.-D-Galactopyranoside, met...	249	C10H19NO6	017296-07-0	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

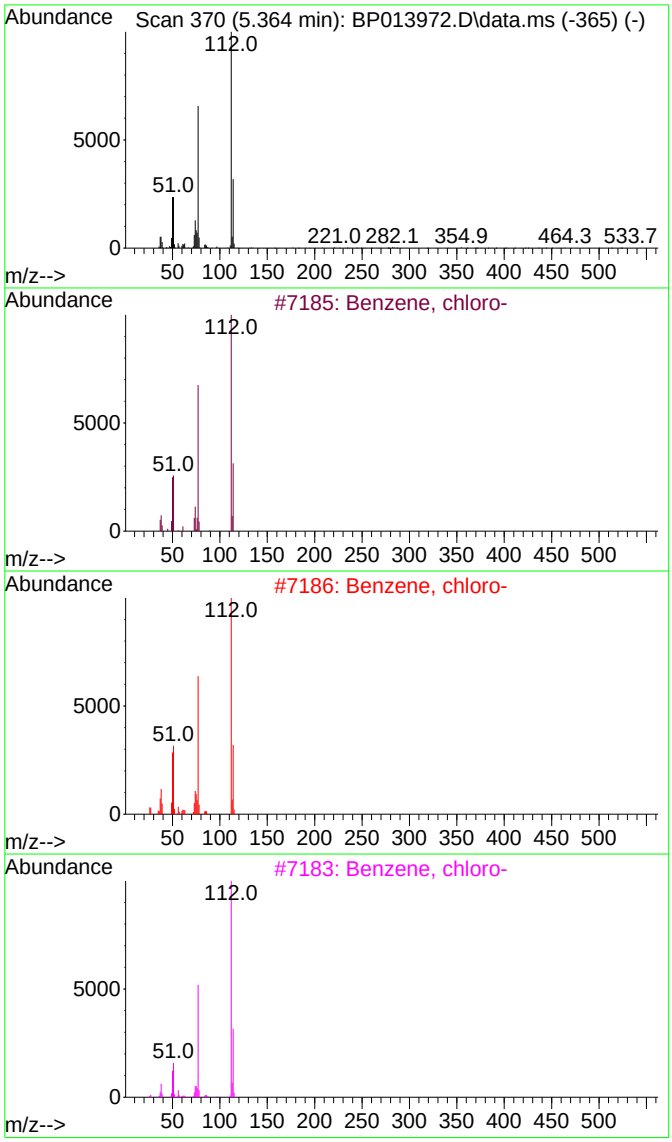
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

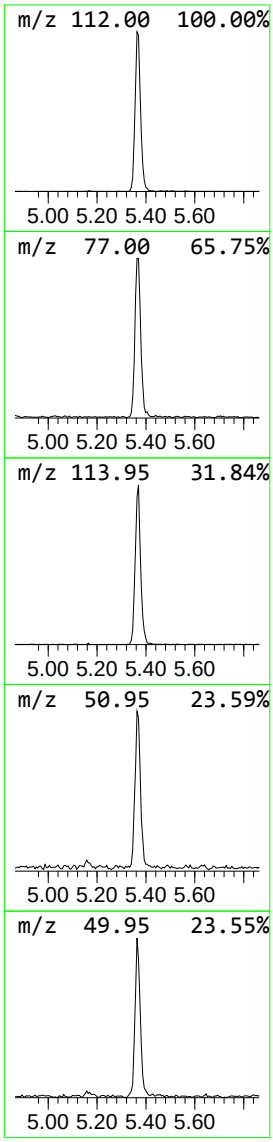
****

Peak Number 2 Benzene, chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.364	3.64 ng/ul	144810	1,4-Dichlorobenzene-d4	8.093

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

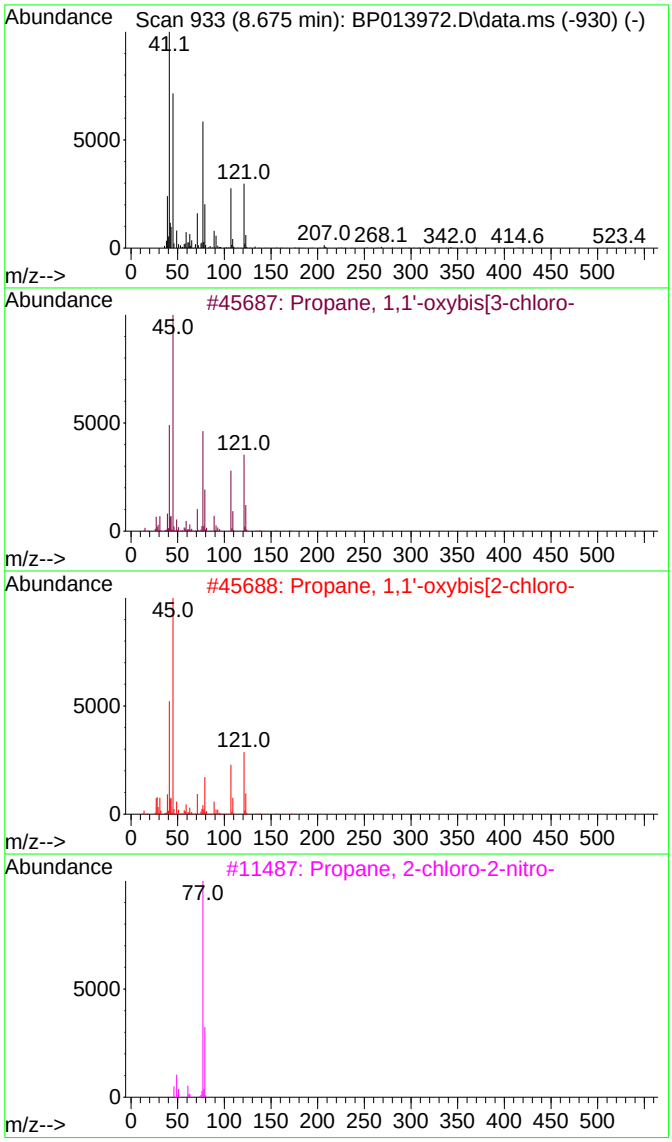
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

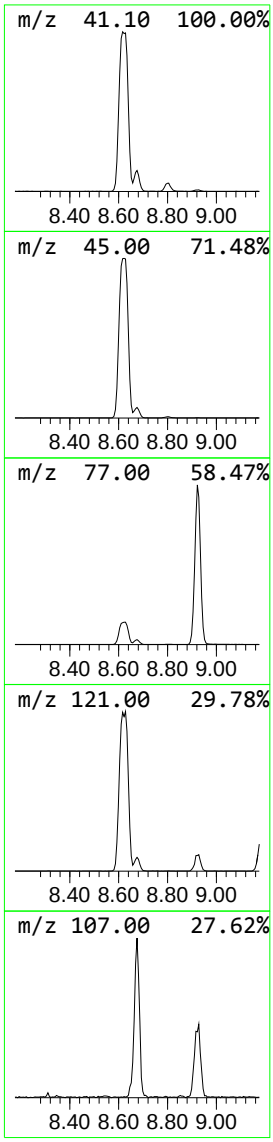
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Propane, 1,1'-oxybis[3-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.675	3.15 ng/ul	125183	1,4-Dichlorobenzene-d4	8.093

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1'-oxybis[3-chloro-	170	C6H12Cl2O	000629-36-7	83
2	Propane, 1,1'-oxybis[2-chloro-	170	C6H12Cl2O	054460-96-7	45
3	Propane, 2-chloro-2-nitro-	123	C3H6ClNO2	000594-71-8	43
4	Bis(2-chloro-1-methylethyl) ether	170	C6H12Cl2O	000108-60-1	32
5	Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
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Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
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Instrument :
BNA_P
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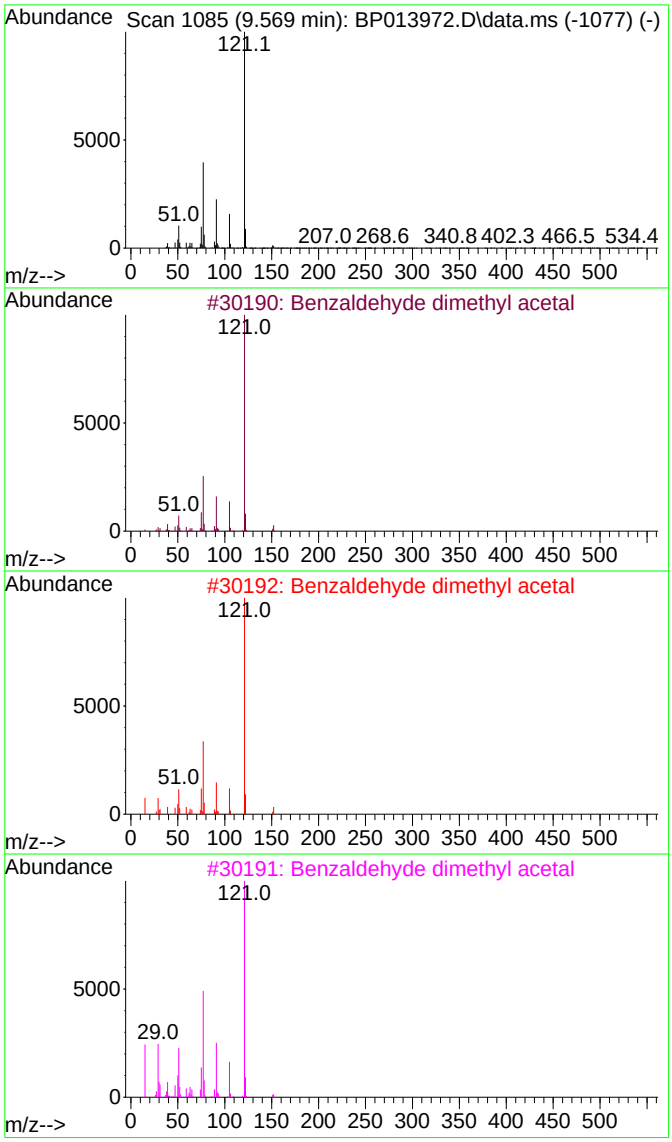
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

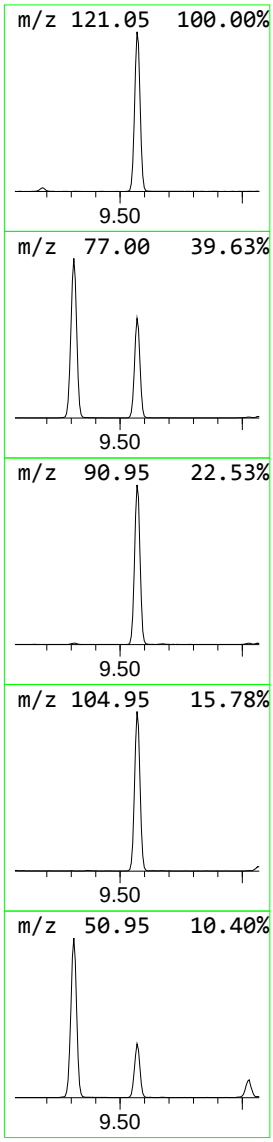
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzaldehyde dimethyl acetal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.569	45.42 ng/ul	2718080	Naphthalene-d8	10.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	91
2	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	90
3	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	83
4	Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	78
5	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

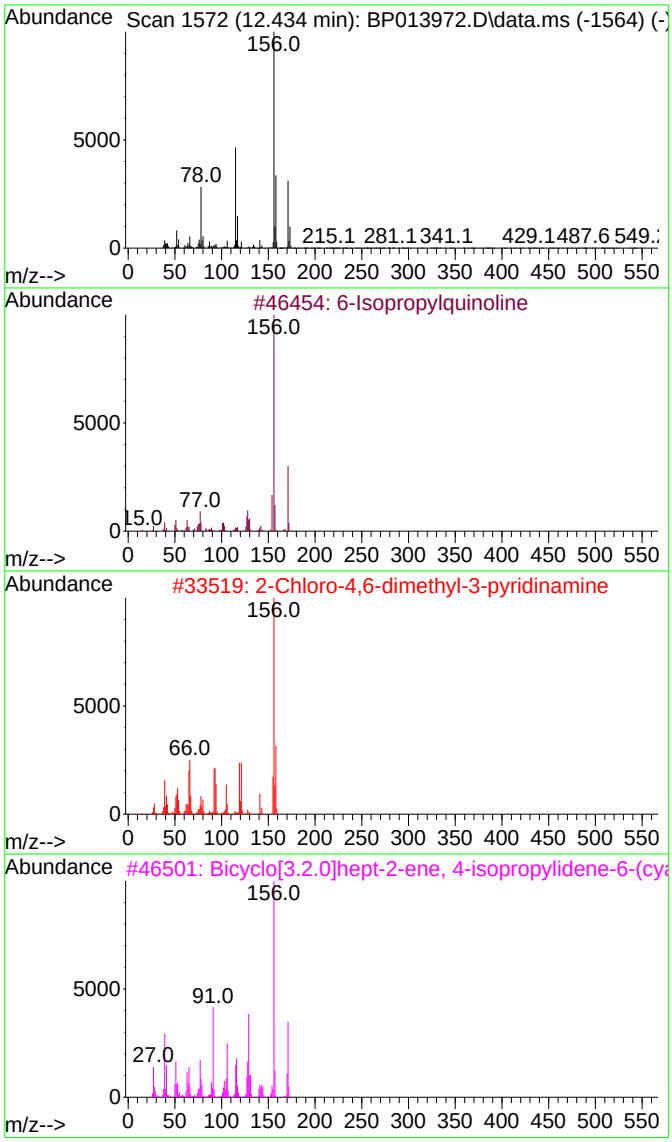
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

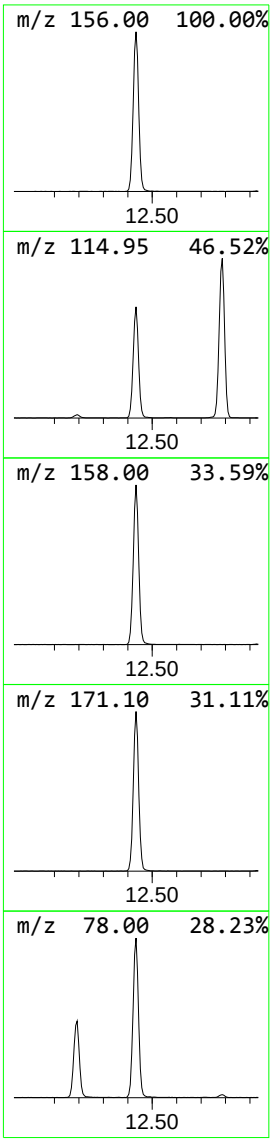
****

Peak Number 5 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.434	45.25 ng/ul	2708240	Naphthalene-d8	10.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Isopropylquinoline	171	C12H13N	000135-79-5	43
2	2-Chloro-4,6-dimethyl-3-pyridina...	156	C7H9ClN2	140413-40-7	32
3	Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
4	Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	32
5	Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28





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Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

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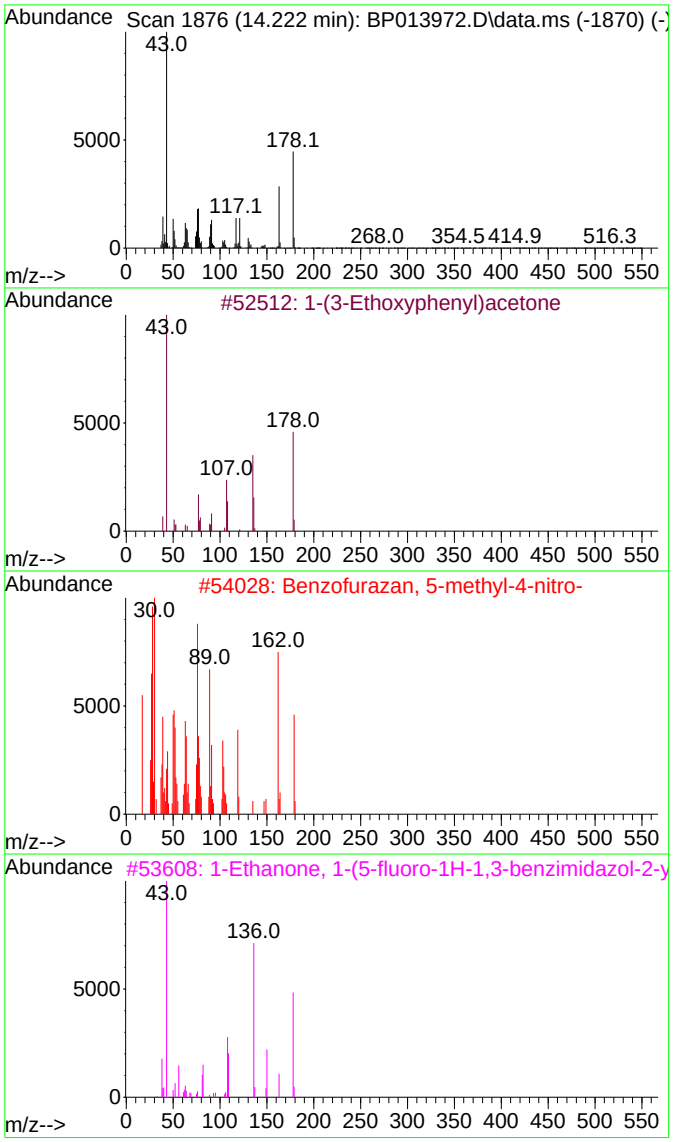
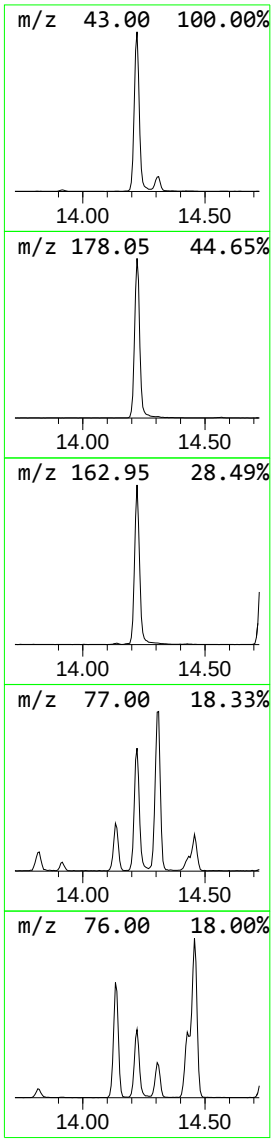
Peak Number 6 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.222	20.31 ng/ul	1696220	Acenaphthene-d10	14.734

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(3-Ethoxyphenyl)acetone	178	C11H14O2	023037-47-0	25
2	Benzofurazan, 5-methyl-4-nitro-	179	C7H5N3O3	016322-21-7	11
3	1-Ethanone, 1-(5-fluoro-1H-1,3-b...	178	C9H7FN2O	1000362-63-4	9
4	3-Carene, 4-acetyl-	178	C12H18O	1000156-14-1	9
5	Benzeneethanol, .alpha.-methyl-3...	178	C12H18O	054518-11-5	9





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Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

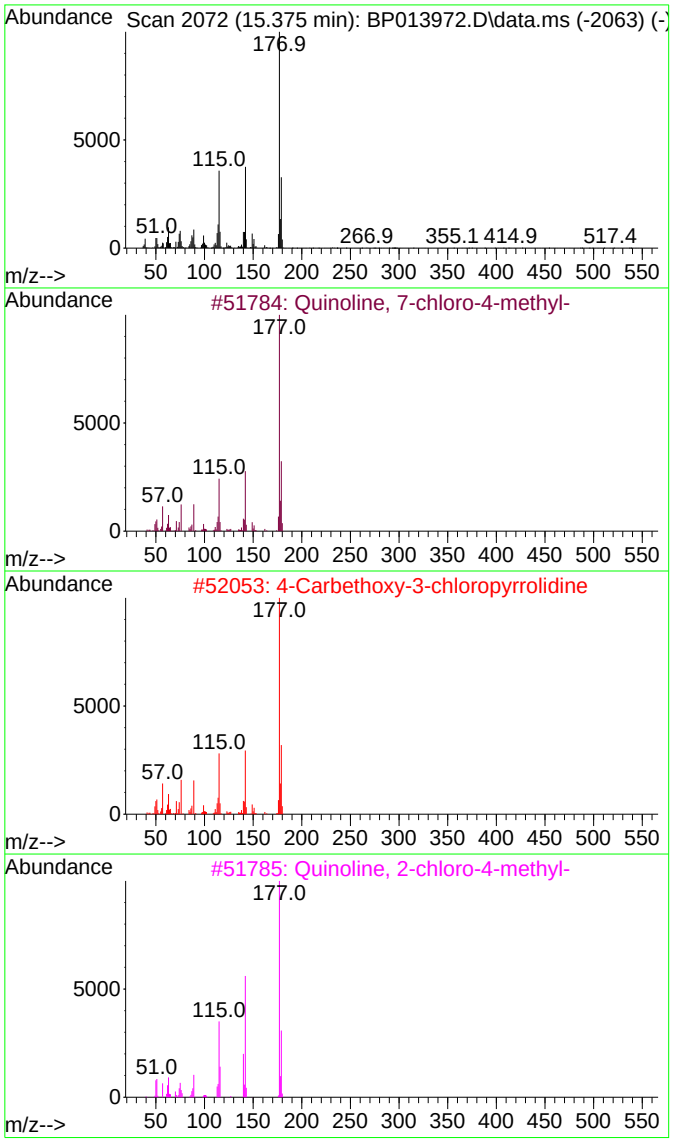
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

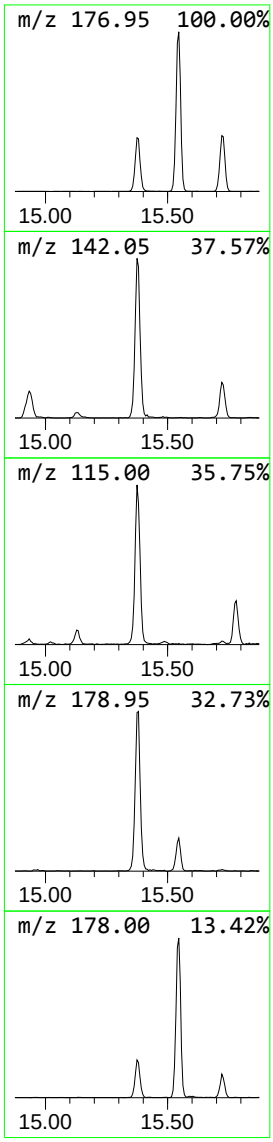
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Quinoline, 7-chloro-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.375	6.60 ng/ul	551239	Acenaphthene-d10	14.734

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinoline, 7-chloro-4-methyl-	177	C10H8ClN	040941-53-5	94
2	4-Carbethoxy-3-chloropyrrolidine	177	C7H12ClNO2	1010255-97-9	94
3	Quinoline, 2-chloro-4-methyl-	177	C10H8ClN	000634-47-9	90
4	6-Chloro-2-methylquinoline	177	C10H8ClN	000092-46-6	90
5	Quinoline, 7-chloro-2-methyl-	177	C10H8ClN	004965-33-7	86





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Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

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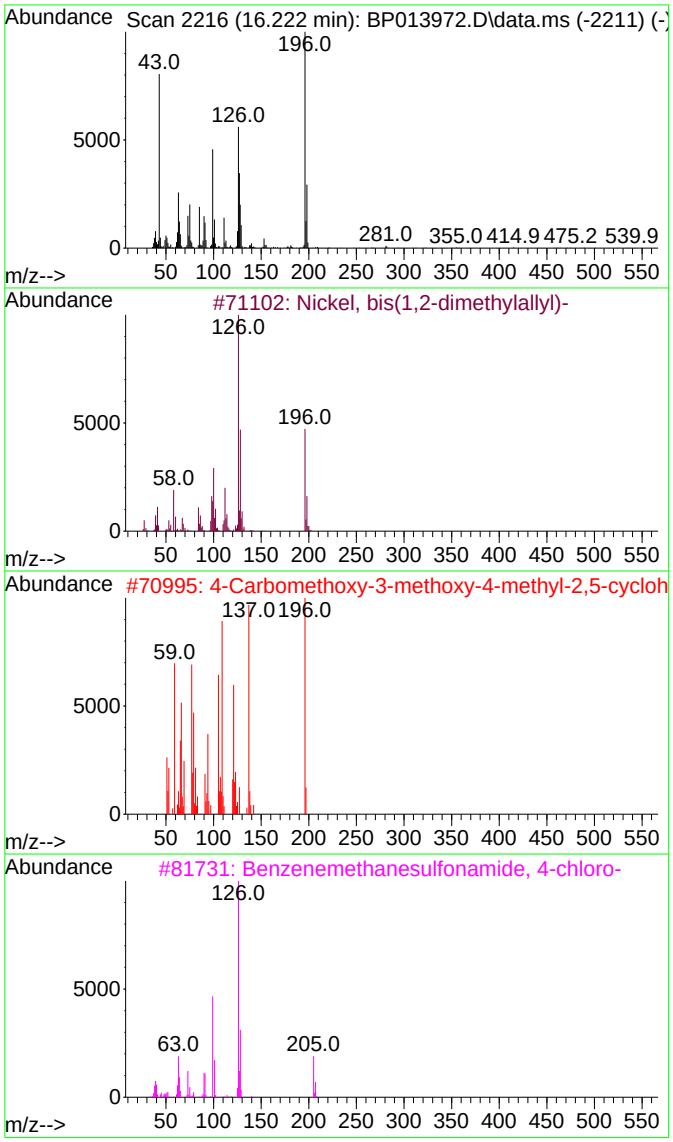
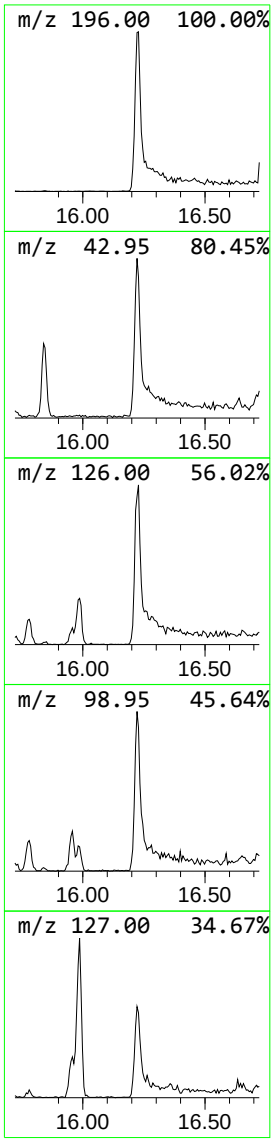
Peak Number 8 Nickel, bis(1,2-dimethylall... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.222	2.07 ng/ul	240822	Phenanthrene-d10	17.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nickel, bis(1,2-dimethylallyl)-	196	C10H18Ni	1000150-48-3	50
2	4-Carbomethoxy-3-methoxy-4-methy...	196	C10H12O4	120696-19-7	35
3	Benzenemethanesulfonamide, 4-chl...	205	C7H8ClNO2S	071799-35-4	30
4	2-Amino-3-chloro-5-trifluorometh...	196	C6H4ClF3N2	079456-26-1	27
5	2-Chloro-4-(3-thienyl)pyrimidine	196	C8H5ClN2S	063558-68-9	27





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Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

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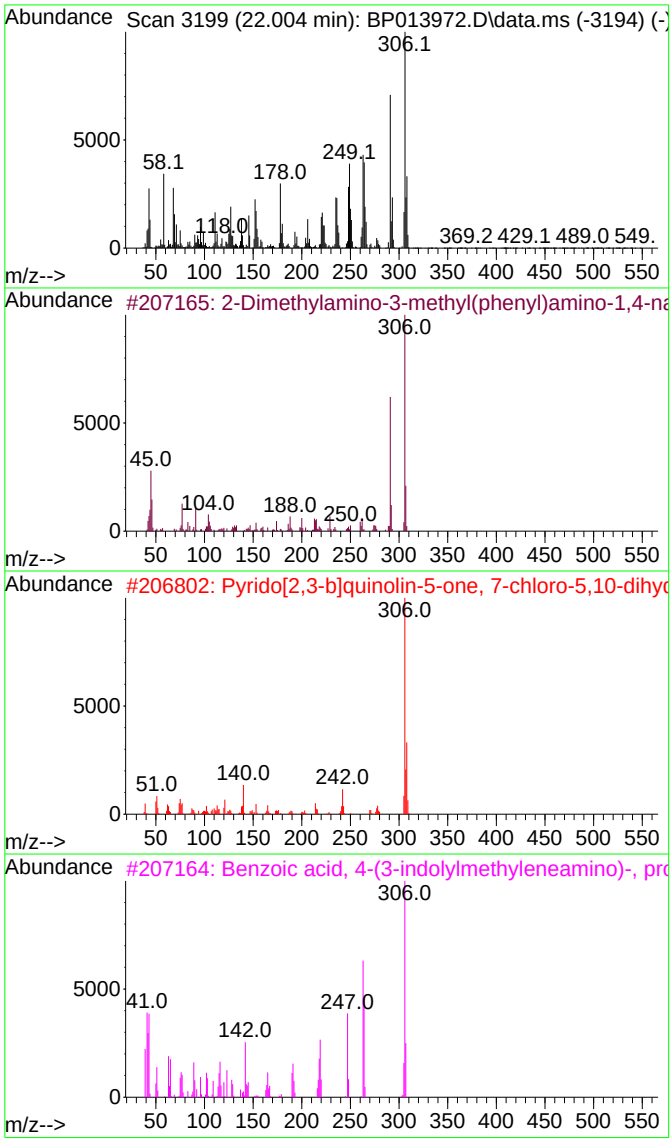
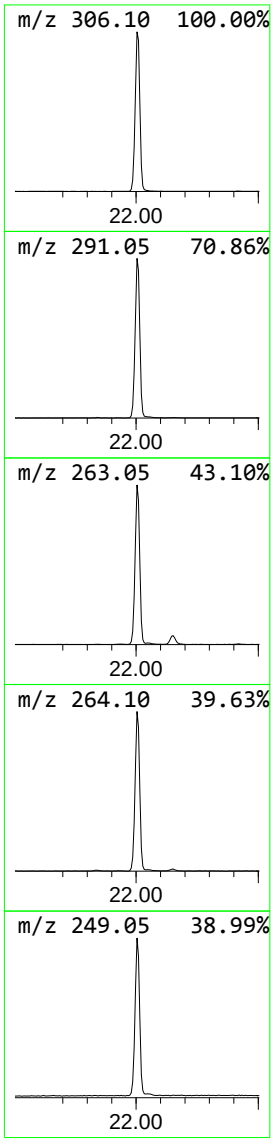
Peak Number 9 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.004	15.32 ng/ul	8377850	Chrysene-d12	21.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dimethylamino-3-methyl(phenyl)...	306	C19H18N2O2	134614-10-1	38
2	Pyrido[2,3-b]quinolin-5-one, 7-c...	306	C18H11ClN2O	069750-07-8	38
3	Benzoic acid, 4-(3-indolylmethyl...	306	C19H18N2O2	304669-02-1	35
4	Acetamide, 2-[(4,5-dihydro-6-met...	306	C13H14N4O3S	1000317-32-8	25
5	4-Amino-2-chloro-5-iodobenzonitr...	306	C9H8ClIN2	1000499-50-6	22





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
Data File : BP013972.D
Acq On : 08 Mar 2023 11:53
Operator : CG/JU
Sample : 01746-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCAG9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.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
2-Pentanone, 4-...	5.158	4.3	ng/ul	170780	1	8.093	795326	20.0
Benzene, chloro-	5.364	3.6	ng/ul	144810	1	8.093	795326	20.0
Propane, 1,1'-o...	8.675	3.1	ng/ul	125183	1	8.093	795326	20.0
Benzaldehyde di...	9.569	45.4	ng/ul	2718080	2	10.916	1196950	20.0
unknown-01	12.434	45.3	ng/ul	2708240	2	10.916	1196950	20.0
unknown-02	14.222	20.3	ng/ul	1696220	3	14.734	1670130	20.0
Quinoline, 7-ch...	15.375	6.6	ng/ul	551239	3	14.734	1670130	20.0
Nickel, bis(1,2...	16.222	2.1	ng/ul	240822	4	17.486	2327850	20.0
unknown-03	22.004	15.3	ng/ul	8377850	5	21.569	10938600	20.0