LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

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-BP040722.M
Title : SVOA CALIBRATION

Signal : TIC: BP009829.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.123	3	6	21	rVB	696638	955775	35.48%	2.691%
2	3.381	47	50	55	rBV2	11098	14197	0.53%	0.040%
3	3.805	119	122	131	rBV	70670	120234	4.46%	0.338%
4	3.870	131	133	137	rVV5	12326	16521	0.61%	0.047%
5	3.905	137	139	146	rVV6	9509	13696	0.51%	0.039%
6	4.140	173	179	188	rVB2	23231	39816	1.48%	0.112%
7	4.358	212	216	223	rVB5	10394	17718	0.66%	0.050%
8	4.781	282	288	298	rVB	55441	88589	3.29%	0.249%
9	4.969	315	320	325	rBV4	6477	10584	0.39%	0.030%
10	5.058	329	335	341	rBV2	7225	14293	0.53%	0.040%
11	5.269	364	371	382	rBV	572900	898766	33.36%	2.530%
12	6.628	596	602	608	rVB7	8926	15058	0.56%	0.042%
13	6.964	654	659	663	rVB8	3421	8252	0.31%	0.023%
14	7.116	676	685	694	rBV	214509	362206	13.45%	1.020%
15	7.287	707	714	723	rBV	171614	278124	10.32%	0.783%
16	7.493	742	749	762	rBV	210482	368968	13.70%	1.039%
17	7.605	762	768	769	rVV5	4825	9421	0.35%	0.027%
18	7.963	821	829	838	rBV	509844	830595	30.83%	2.338%
19	8.040	838	842	847	rVB6	5959	11102	0.41%	0.031%
20	8.258	874	879	884	rBV	47135	77093	2.86%	0.217%
21	8.316	884	889	897	rVB2	27647	46260	1.72%	0.130%
22	8.393	897	902	907	rBV8	5543	11251	0.42%	0.032%
23	8.575	927	933	940	rVV2	65211	110493	4.10%	0.311%
24	8.663	940	948	954	rVV	288854	471192	17.49%	1.327%
25	8.716	954	957	963	rVV3	13920	26824	1.00%	0.076%
26	8.781	963	968	976	rVB	27407	44730	1.66%	0.126%
27	9.116	1018	1025	1032	rBV	225184	374206	13.89%	1.053%
28	9.457	1077	1083	1090	rBV	76743	127175	4.72%	0.358%
29	9.781	1130	1138	1143	rBV	77694	130643	4.85%	0.368%
30	9.846	1143	1149	1161	rVB	168996	293127	10.88%	0.825%
31	10.193	1201	1208	1214	rBV	6075	15320	0.57%	0.043%
32	10.381	1233	1240	1253	rBV	274022	471922	17.52%	1.329%
33	10.646	1277	1285	1291	rVV	15114	29289	1.09%	0.082%
34	10.769	1298	1306	1313	rBV	744637	1272730	47.25%	3.583%
35	10.899	1321	1328	1340	rBV	280118	489478	18.17%	1.378%
36	10.987	1340	1343	1356	rVB	10858	24936	0.93%	0.070%

LSC Area Percent Report

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Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

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-BP040722.M
Title : SVOA CALIBRATION

37	11.122	1361	1366	1371	rVB9	3999	7521	0.28%	0.021%
38	11.222	1377	1383	1388	rVB9	4224	8312	0.31%	0.023%
39	11.316	1394	1399	1405	rVB8	7971	15108	0.56%	0.043%
40	12.199	1542	1549	1556	rBV9	6820	13507	0.50%	0.038%
41	12.357	1571	1576	1582	rVV7	11806	20922	0.78%	0.059%
42	12.998	1681	1685	1695	rVB7	5324	13403	0.50%	0.038%
43	13.093	1695	1701	1706	rBV9	6035	10913	0.41%	0.031%
44	13.246	1723	1727	1733	rVB4	23391	35505	1.32%	0.100%
45	13.328	1733	1741	1746	rVB3	14835	33490	1.24%	0.094%
46	13.440	1756	1760	1763	rVB2	7657	10710	0.40%	0.030%
47	13.957	1842	1848	1850	rBV2	27417	40200	1.49%	0.113%
48	13.998	1850	1855	1862	rVB	647163	899045	33.37%	2.531%
49	14.287	1892	1904	1916	rBV	601448	916060	34.01%	2.579%
50	14.593	1948	1956	1970	rBV	1238452	1838365	68.24%	5.175%
51	14.769	1978	1986	1998	rBV	222340	388917	14.44%	1.095%
52	15.004	2021	2026	2030	rVB8	7176	11644	0.43%	0.033%
53	15.216	2056	2062	2065	rBV8	7893	14583	0.54%	0.041%
54	15.334	2078	2082	2086	rVB7	6099	9188	0.34%	0.026%
55	15.404	2089	2094	2100	rBV8	7550	13421	0.50%	0.038%
56	15.498	2105	2110	2114	rBV7	5032	8068	0.30%	0.023%
57	15.581	2116	2124	2137	rBV	882201	1319489	48.98%	3.715%
58	15.692	2137	2143	2155	rVB	206981	297544	11.05%	0.838%
59	15.822	2161	2165	2171	rBV4	9159	13796	0.51%	0.039%
60	16.375	2254	2259	2262	rVB7	5928	10127	0.38%	0.029%
61	16.845	2336	2339	2344	rBV7	5654	9048	0.34%	0.025%
62	16.992	2359	2364	2375	rVB6	15338	29383	1.09%	0.083%
63	17.110	2379	2384	2385	rBV5	5542	7776	0.29%	0.022%
64	17.345	2416	2424	2434	rBV	1559971	2331535	86.55%	6.564%
65	17.439	2434	2440	2452	rVV	1091155	1623921	60.28%	4.572%
66	17.528	2452	2455	2459	rVV5	15050	23305	0.87%	0.066%
67	17.563	2459	2461	2467	rVB7	6682	8413	0.31%	0.024%
68	17.857	2506	2511	2517	rVV4	17838	30846	1.15%	0.087%
69	17.963	2524	2529	2533	rBV3	21245	29283	1.09%	0.082%
70	18.175	2562	2565	2569	rBV5	8253	8992	0.33%	0.025%
71	18.222	2569	2573	2582	rBV3	57537	93535	3.47%	0.263%
72	18.298	2584	2586	2589	rVV2	7748	7654	0.28%	0.022%
73	18.351	2593	2595	2602	rVV5	6469	10543	0.39%	0.030%
74	18.769	2661	2666	2673	rBV3	35133	53101	1.97%	0.149%
75	18.981	2698	2702	2704	rBV2	20246	25306	0.94%	0.071%
76	19.016	2704	2708	2713	rVB	53632	78054	2.90%	0.220%

LSC Area Percent Report

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Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Title : SVOA CALIBRATION

77	19.245	2744	2747	2754	rVB7	18377	37001	1.37%	0.104%
78	19.316	2756	2759	2762	rBV2	38669	52435	1.95%	0.148%
79	19.386	2767	2771	2773	rVV	59372	77742	2.89%	0.219%
80	19.416	2773	2776	2787	rVB	505166	652256	24.21%	1.836%
81	19.586	2801	2805	2810	rBV2	25629	42093	1.56%	0.119%
82	19.628	2810	2812	2814	rVV3	15643	17188	0.64%	0.048%
83	19.669	2814	2819	2823	rVV	1581152	2114717	78.50%	5.954%
84	19.704	2823	2825	2831	rVB	226001	247321	9.18%	0.696%
85	19.804	2837	2842	2847	rBV	649435	810353	30.08%	2.281%
86	19.851	2847	2850	2854	rVV5	15537	16743	0.62%	0.047%
87	19.904	2854	2859	2865	rVV	562359	686357	25.48%	1.932%
88	20.180	2902	2906	2911	rVB3	34238	47783	1.77%	0.135%
89	20.298	2921	2926	2931	rBV	912713	1051210	39.02%	2.959%
90	20.351	2931	2935	2942	rVB	387633	455136	16.90%	1.281%
91	20.586	2972	2975	2980	rVB3	31858	37005	1.37%	0.104%
92	20.727	2994	2999	3006	rVB	1833807	2073388	76.97%	5.837%
93	21.039	3049	3052	3056	rBV	144416	165186	6.13%	0.465%
94	21.133	3063	3068	3077	rBV2	436162	680660	25.27%	1.916%
95	21.422	3112	3117	3122	rBV	2083751	2693783	100.00%	7.584%
96	21.527	3131	3135	3142	rVB3	71030	100934	3.75%	0.284%
97	21.945	3202	3206	3212	rBV	301777	389180	14.45%	1.096%
98	22.433	3286	3289	3296	rVB	120424	182644	6.78%	0.514%
99	23.727	3502	3509	3517	rBV2	962205	1903488	70.66%	5.359%
100	23.886	3529	3536	3548	rVB2	1218819	2574831	95.58%	7.249%

Sum of corrected areas: 35520551

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
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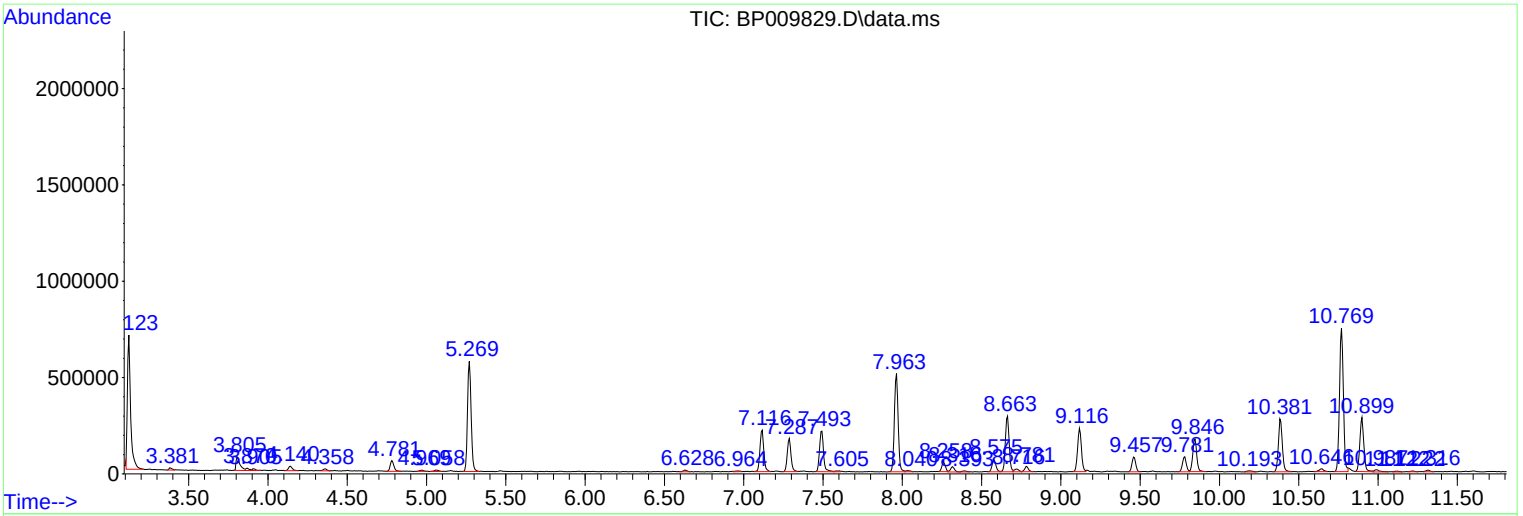
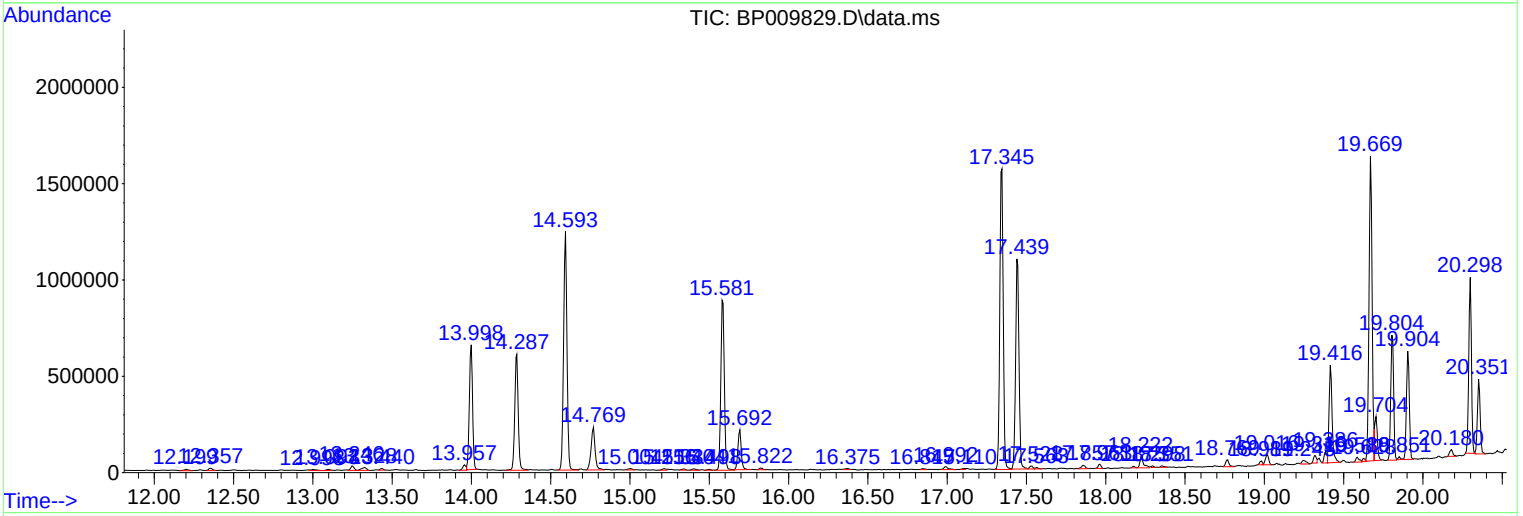
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

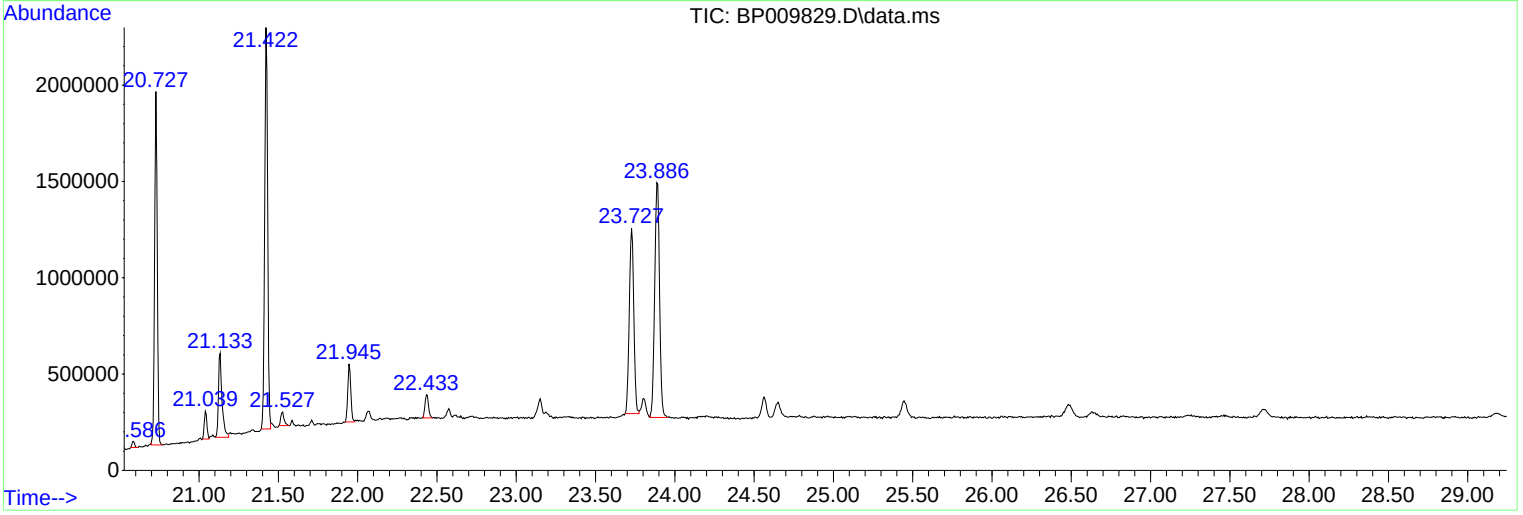
Instrument :
BNA_P
ClientSampleId :
ETNN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
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Operator : CG/JU
Sample : N2293-07
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Instrument :
BNA_P
ClientSampleId :
ETNN7

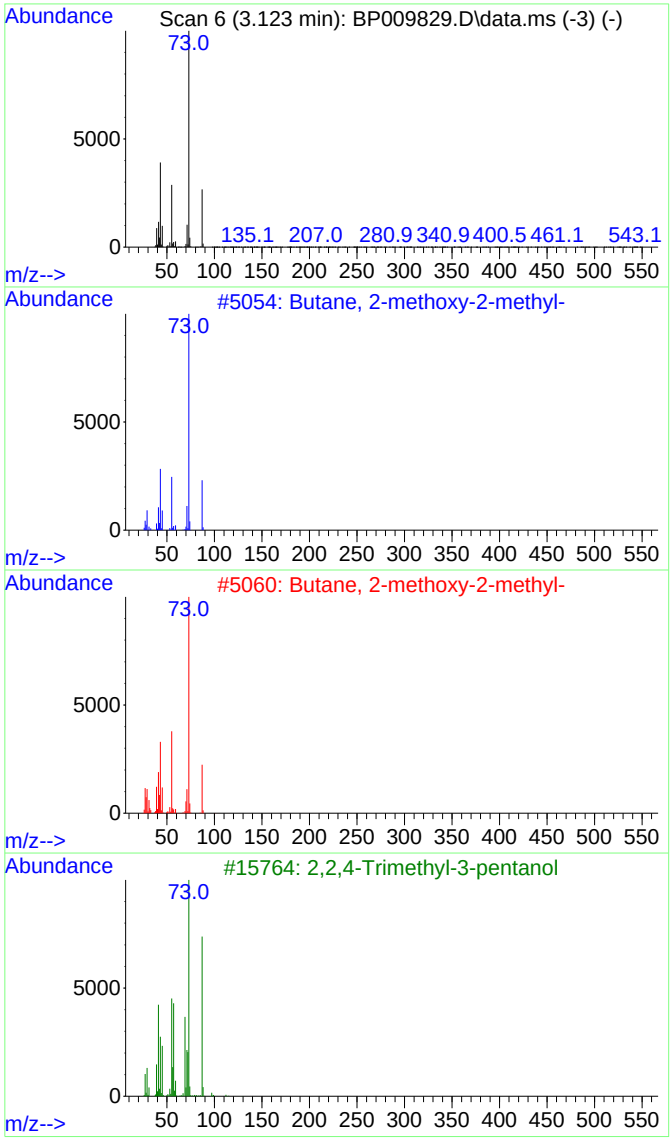
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

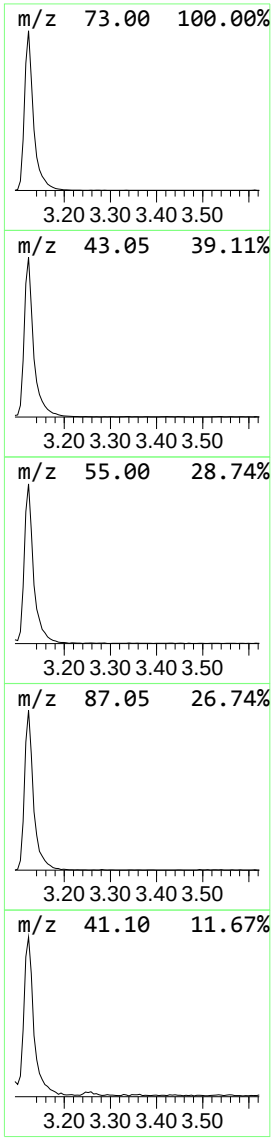
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.123	23.01 ng/ul	955775	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
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Operator : CG/JU
Sample : N2293-07
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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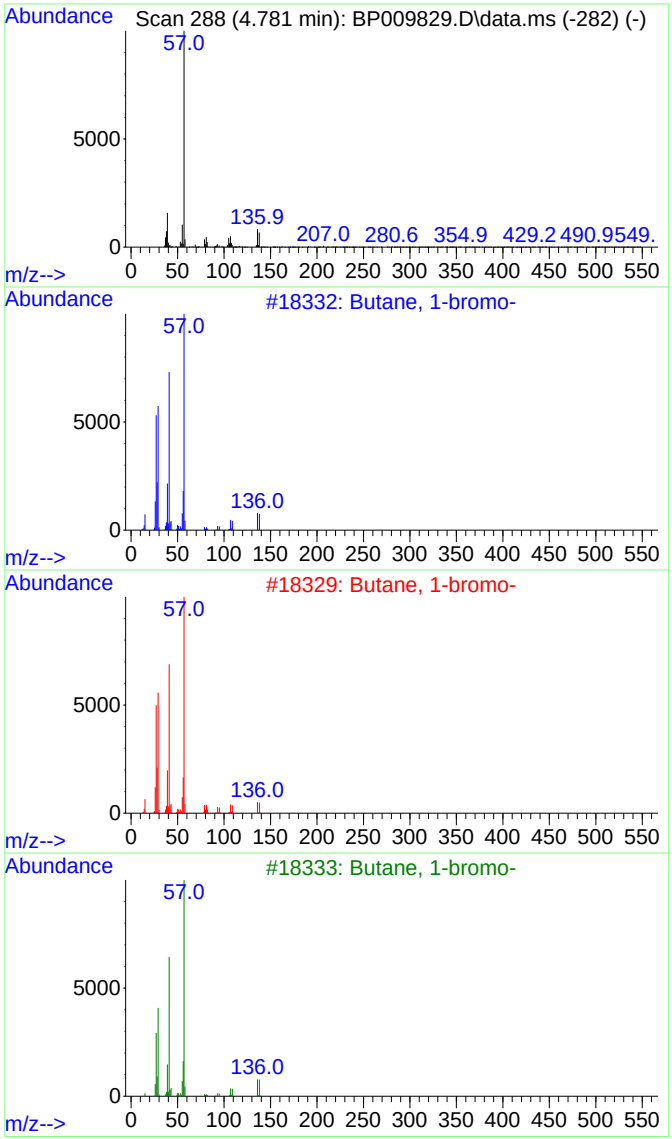
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

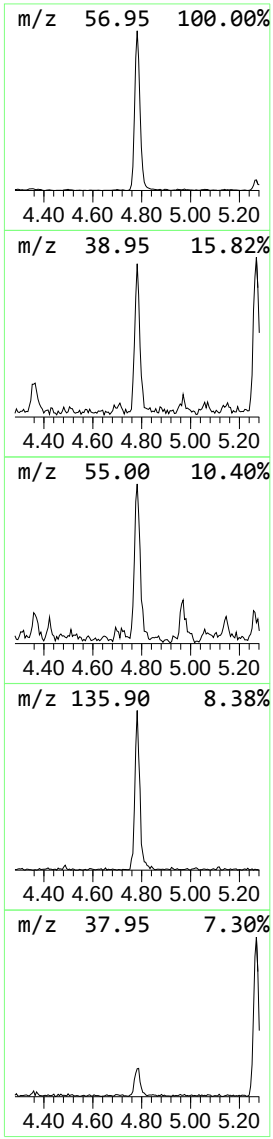
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 1-bromo- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.781	2.13 ng/ul	88589	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1-bromo-	136	C4H9Br	000109-65-9	50
2	Butane, 1-bromo-	136	C4H9Br	000109-65-9	49
3	Butane, 1-bromo-	136	C4H9Br	000109-65-9	25
4	Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	9
5	Oxirane, (bromomethyl)-	136	C3H5BrO	003132-64-7	9





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Instrument :
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ClientSampleId :
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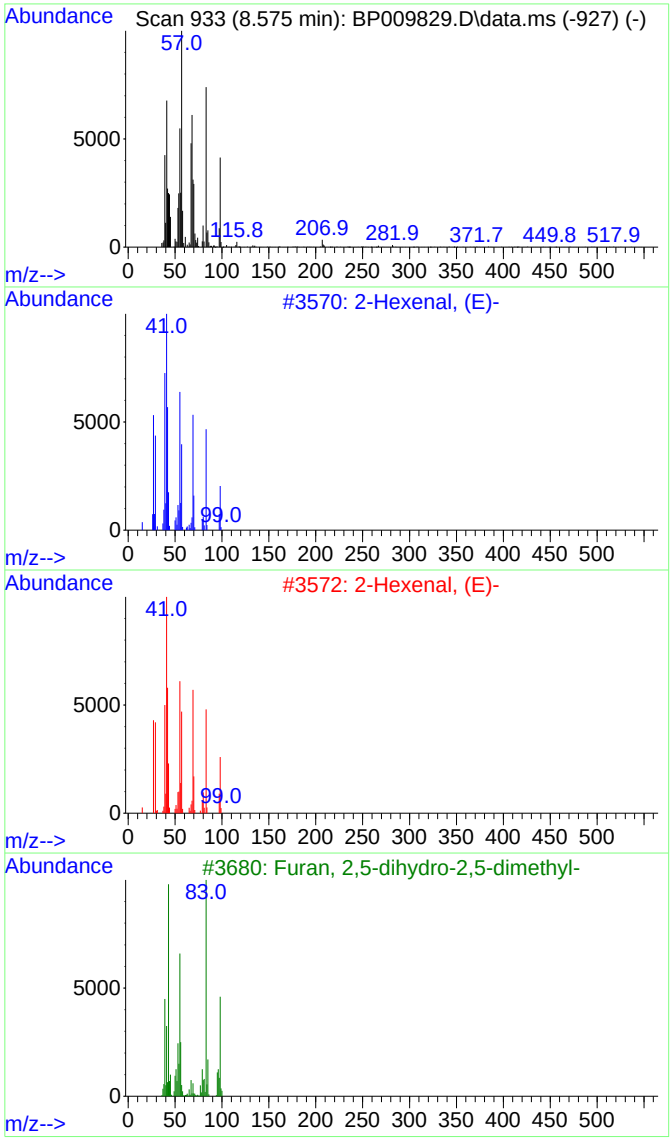
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

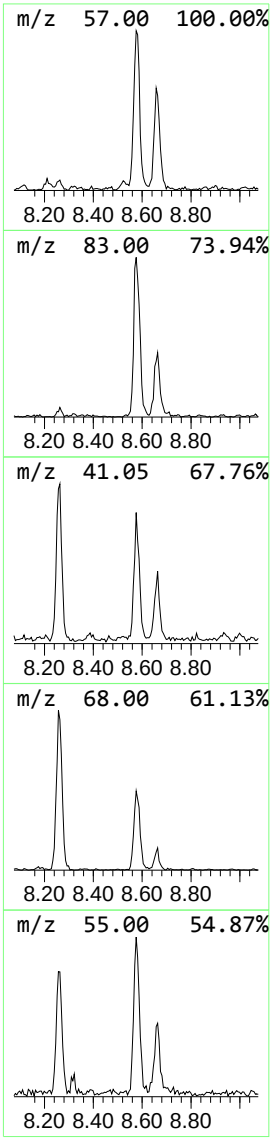
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Hexenal, (E)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.575	2.66 ng/ul	110493	1,4-Dichlorobenzene-d4	7.963

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexenal, (E)-	98	C6H10O	006728-26-3	50
2	2-Hexenal, (E)-	98	C6H10O	006728-26-3	50
3	Furan, 2,5-dihydro-2,5-dimethyl-	98	C6H10O	059242-27-2	50
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	46
5	Cyclohexane, methyl-	98	C7H14	000108-87-2	46





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Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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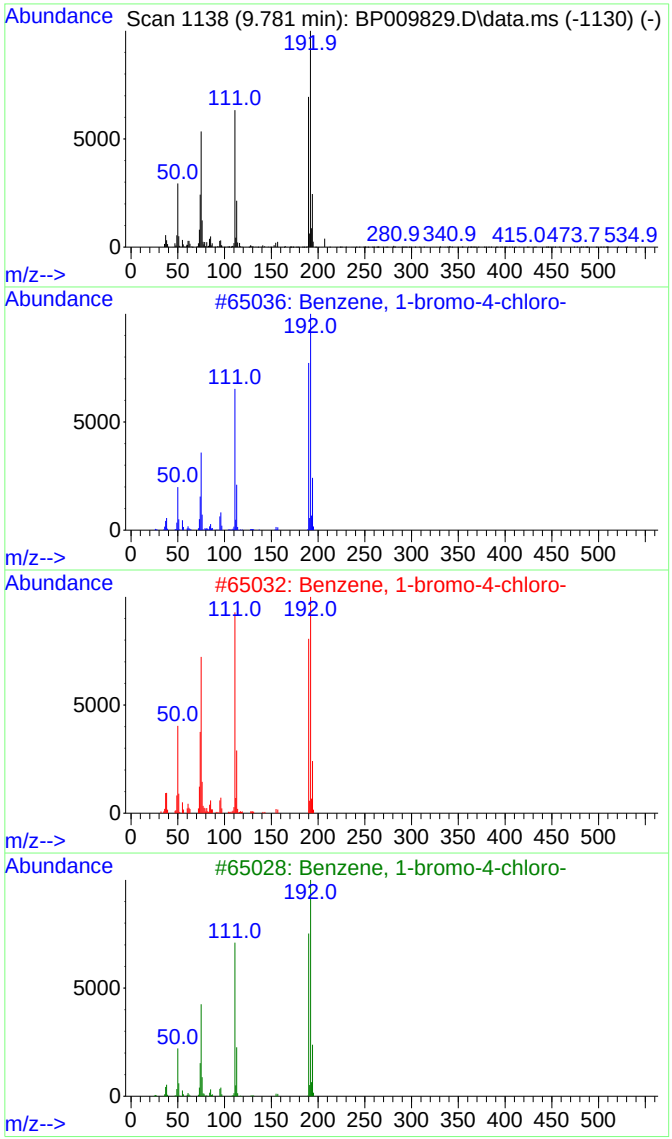
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

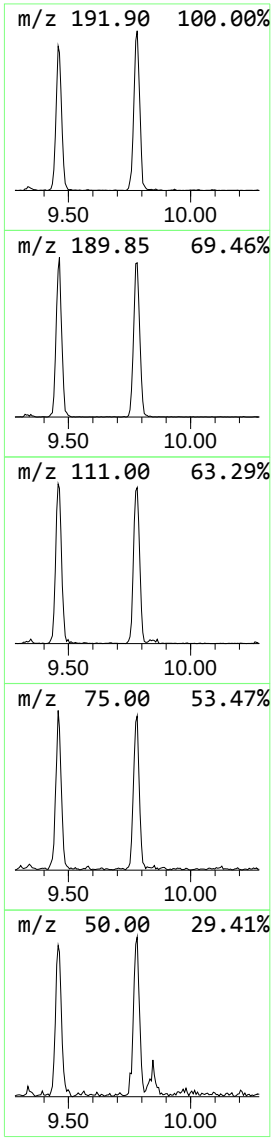
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.781	2.05 ng/ul	130643	Naphthalene-d8	10.769

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
3	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
4	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	96
5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

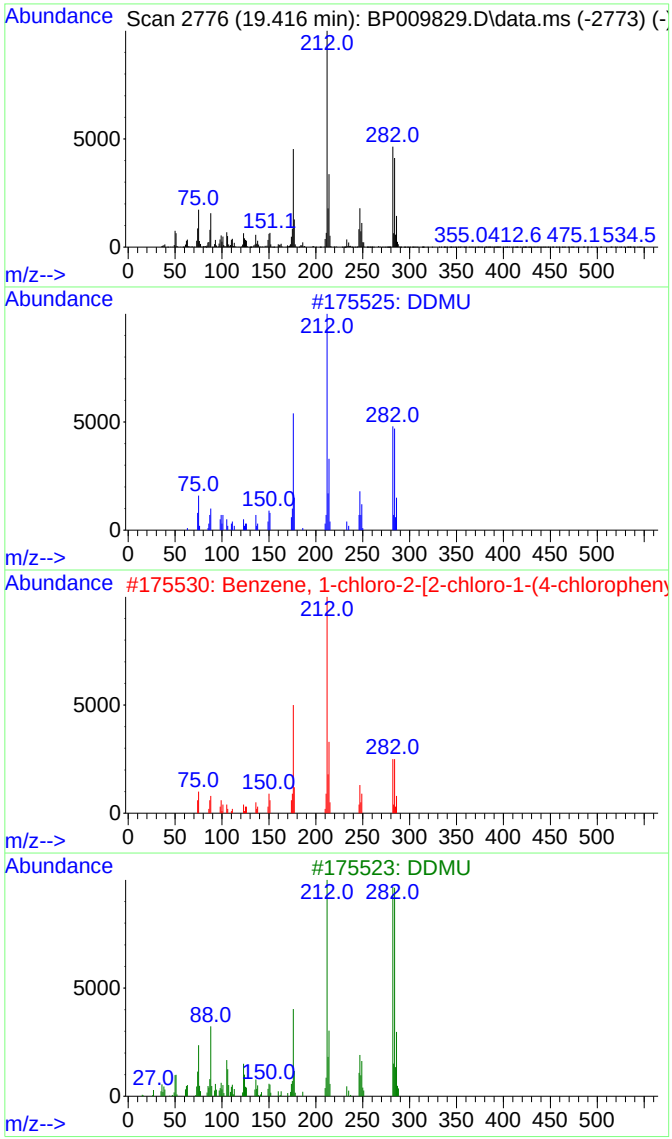
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

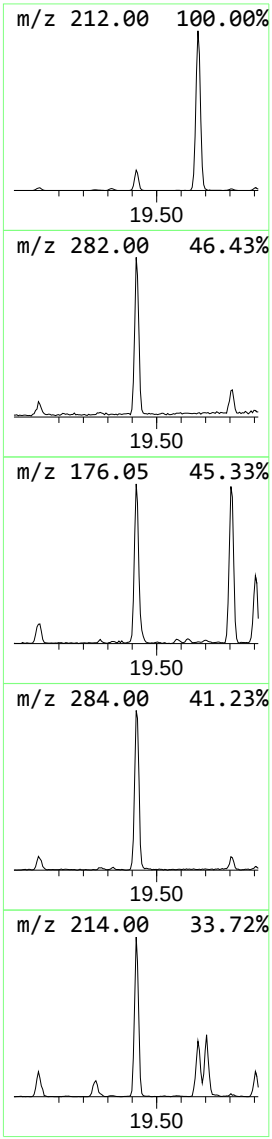
****

Peak Number 6 DDMU Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.416	4.84 ng/ul	652256	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	DDMU	282	C14H9Cl3	001022-22-6	99
2	Benzene, 1-chloro-2-[2-chloro-1-...	282	C14H9Cl3	014835-94-0	98
3	DDMU	282	C14H9Cl3	001022-22-6	95
4	9-Chlorophenanthrene	212	C14H9Cl	000947-72-8	94
5	Benzene, 1-chloro-2-[2-chloro-1-...	282	C14H9Cl3	014835-94-0	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

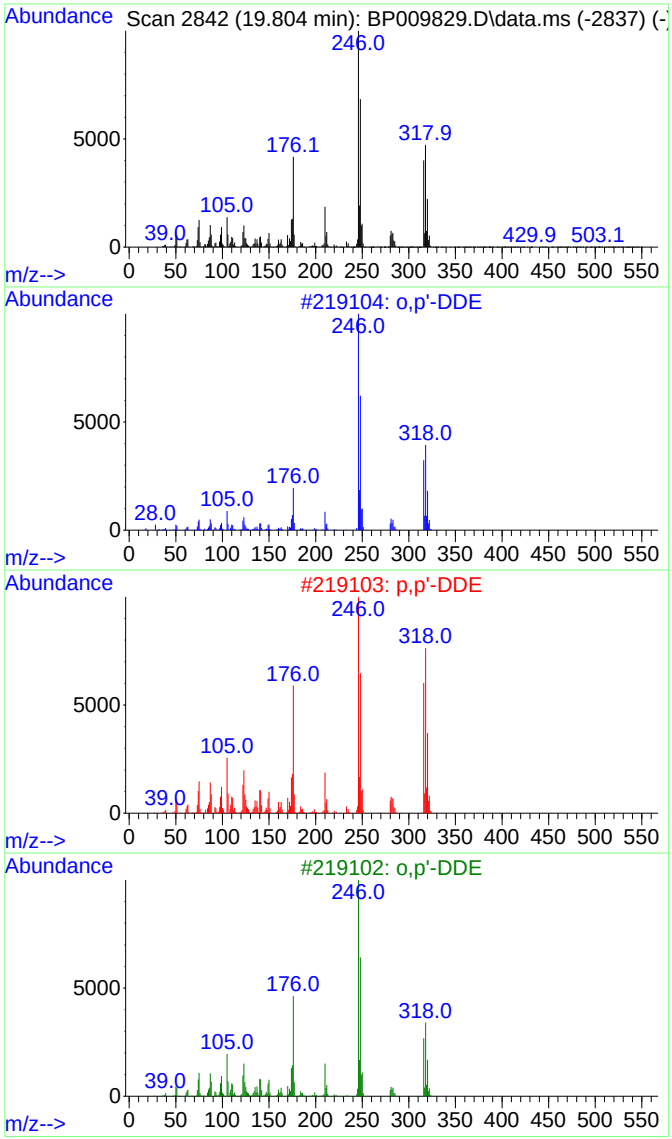
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

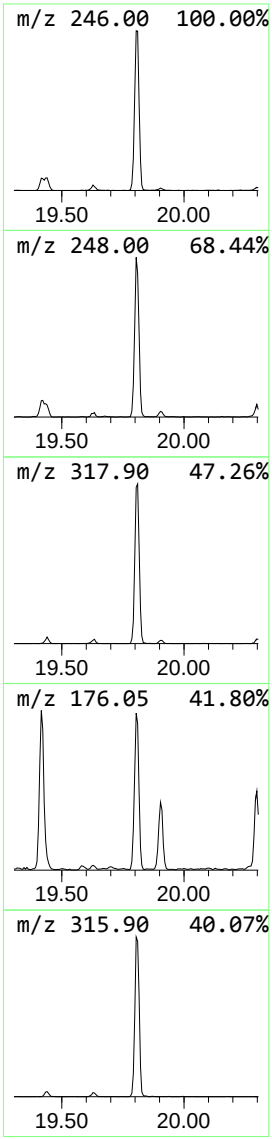
****

Peak Number 7 o,p'-DDE Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.804	6.02 ng/ul	810353	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
2	p,p'-DDE	316	C14H8Cl4	000072-55-9	99
3	o,p'-DDE	316	C14H8Cl4	003424-82-6	99
4	p,p'-DDE	316	C14H8Cl4	000072-55-9	99
5	p,p'-DDE	316	C14H8Cl4	000072-55-9	99





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Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

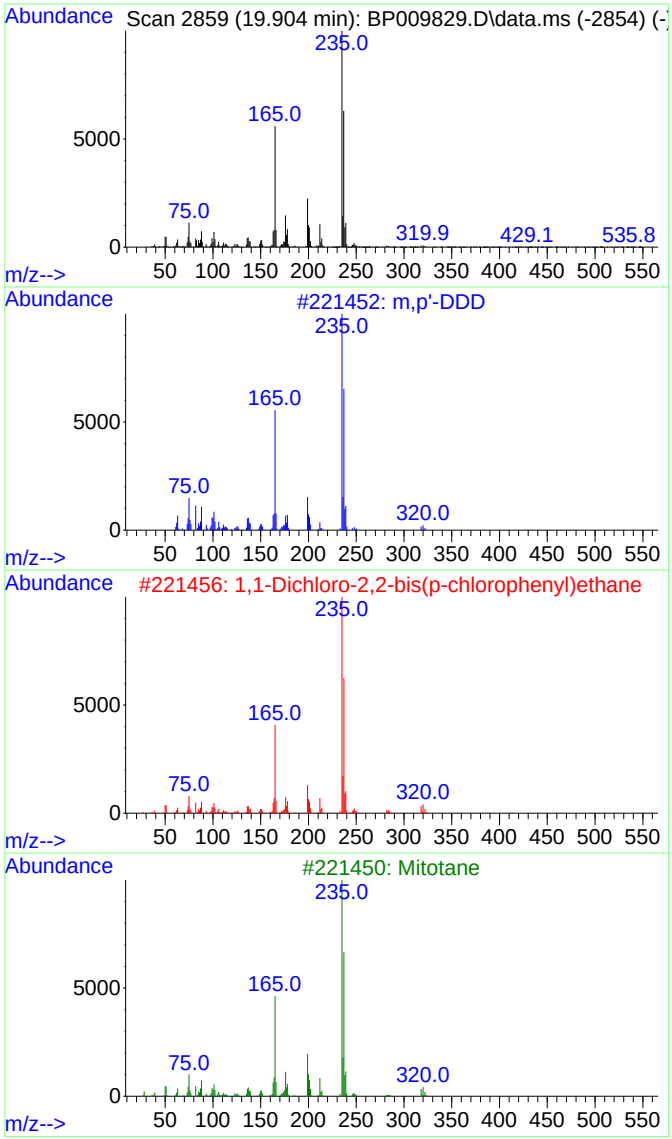
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

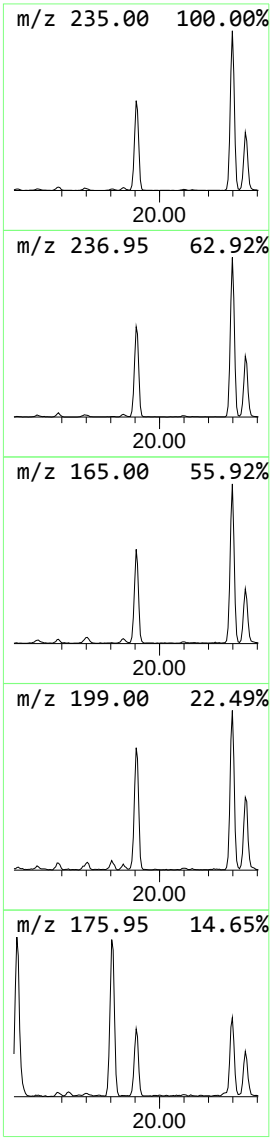
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 m,p'-DDD Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.904	5.10 ng/ul	686357	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m,p'-DDD	318	C14H10Cl4	004329-12-8	98
2	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	96
3	Mitotane	318	C14H10Cl4	000053-19-0	96
4	Mitotane	318	C14H10Cl4	000053-19-0	90
5	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	90





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Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

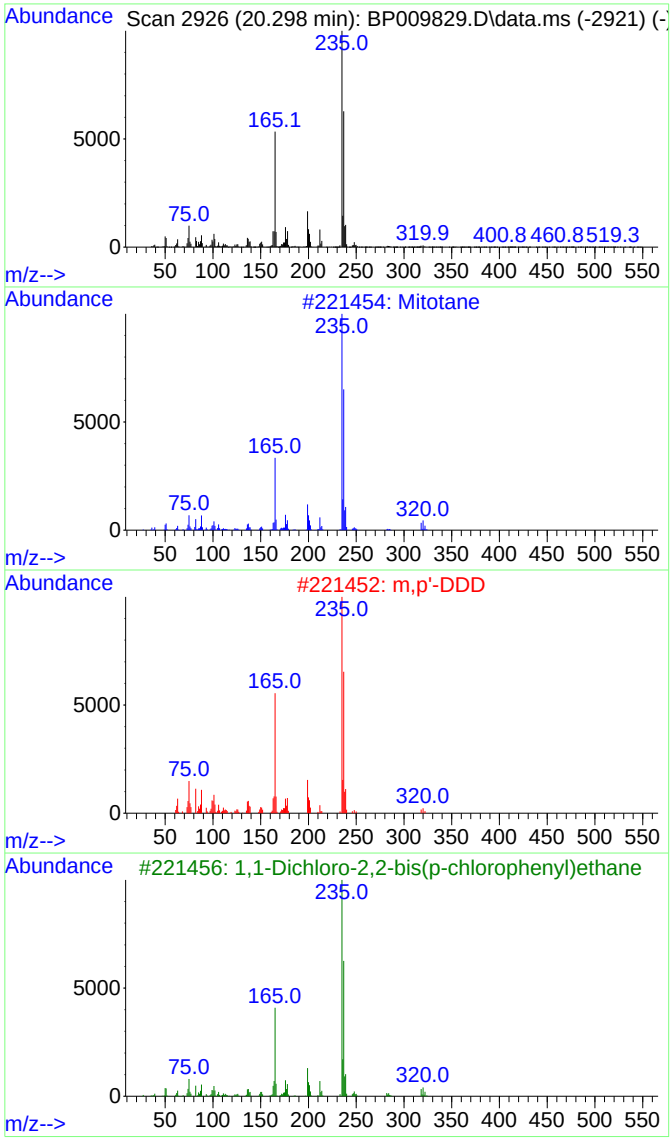
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

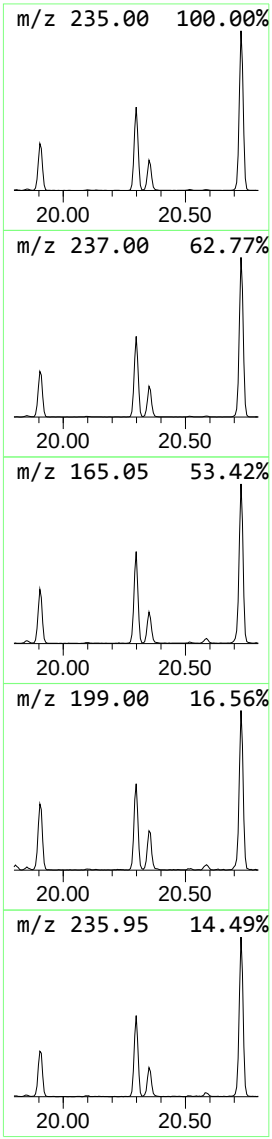
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Mitotane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.298	7.80 ng/ul	1051210	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mitotane	318	C14H10Cl4	000053-19-0	99
2	m,p'-DDD	318	C14H10Cl4	004329-12-8	99
3	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	92
4	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	90
5	p,p'-DDT	352	C14H9Cl5	000050-29-3	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

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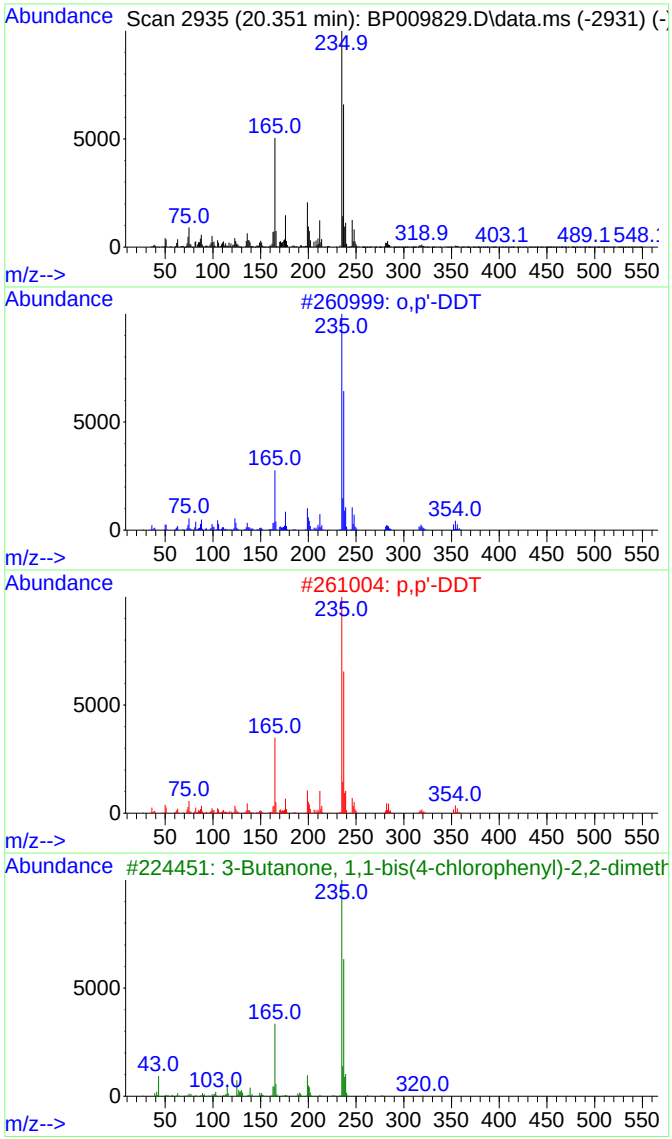
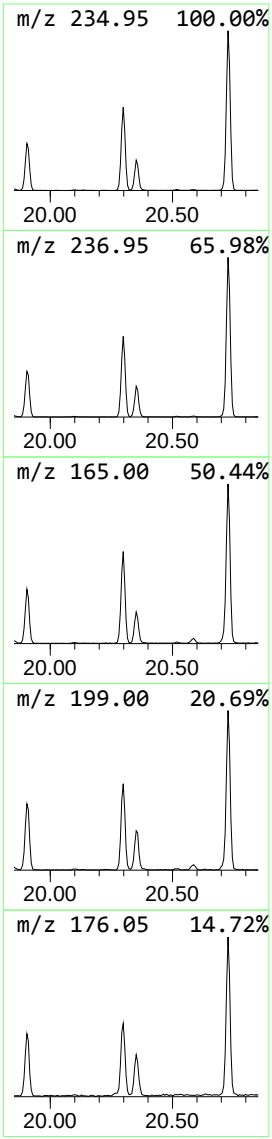
Peak Number 10 o,p'-DDT Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.351	3.38 ng/ul	455136	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o,p'-DDT	352	C14H9Cl5	000789-02-6	98
2	p,p'-DDT	352	C14H9Cl5	000050-29-3	95
3	3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	93
4	o,p'-DDT	352	C14H9Cl5	000789-02-6	91
5	o,p'-DDT	352	C14H9Cl5	000789-02-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

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

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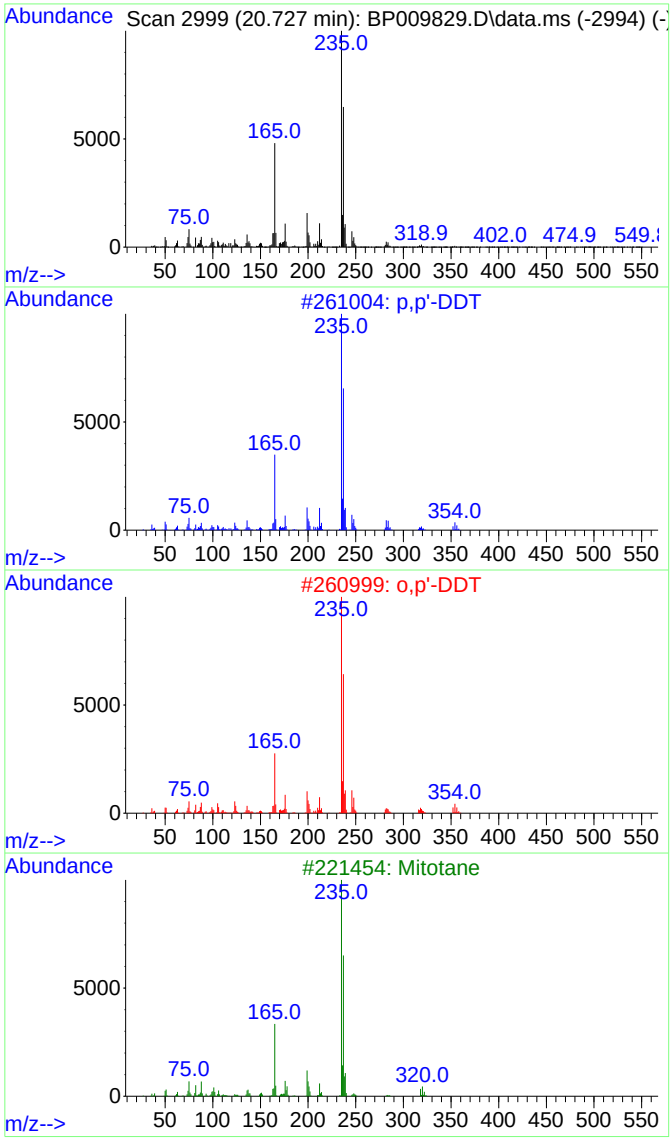
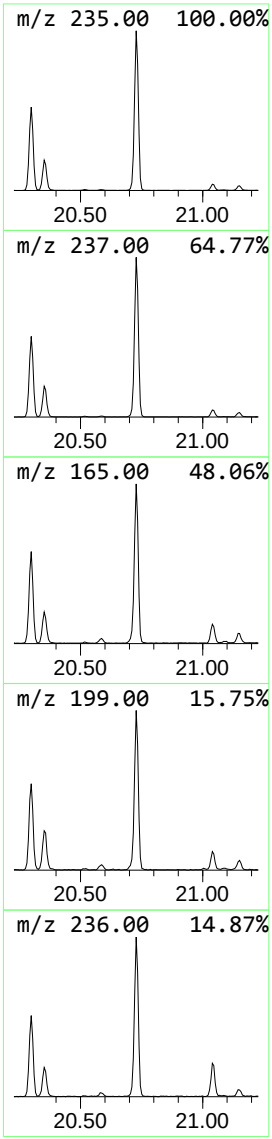
Peak Number 11 p,p'-DDT Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.727	15.39 ng/ul	2073390	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p,p'-DDT	352	C14H9Cl5	000050-29-3	98
2	o,p'-DDT	352	C14H9Cl5	000789-02-6	97
3	Mitotane	318	C14H10Cl4	000053-19-0	97
4	1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	96
5	p,p'-DDT	352	C14H9Cl5	000050-29-3	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

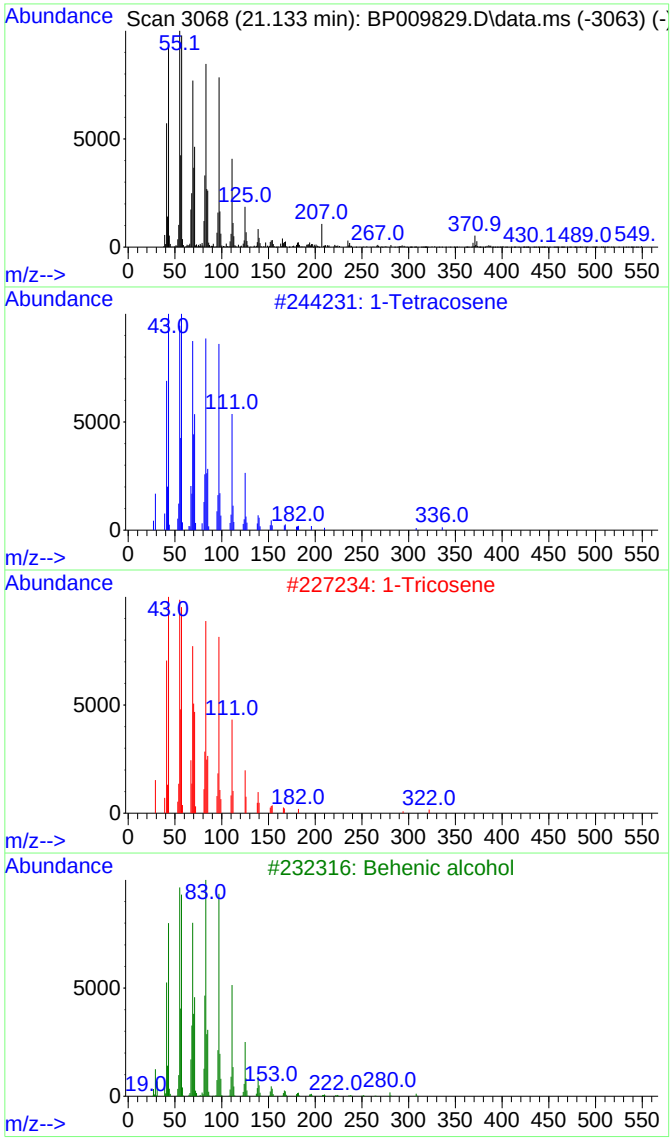
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

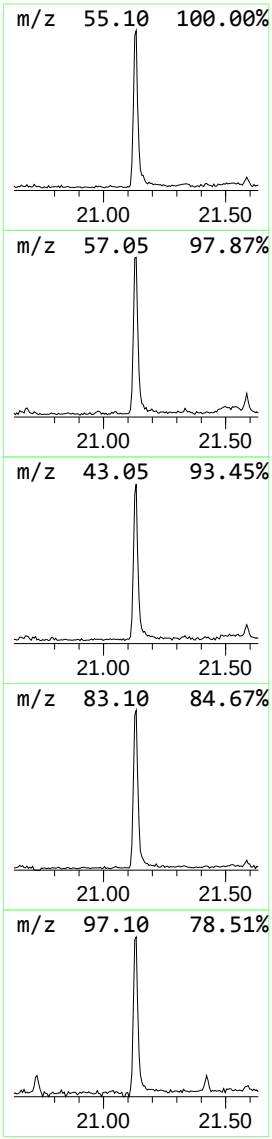
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.133	5.05 ng/ul	680660	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	95
2	1-Tricosene	322	C23H46	018835-32-0	95
3	Behenic alcohol	326	C22H46O	000661-19-8	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	94
5	Pentacos-1-ene	350	C25H50	016980-85-1	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

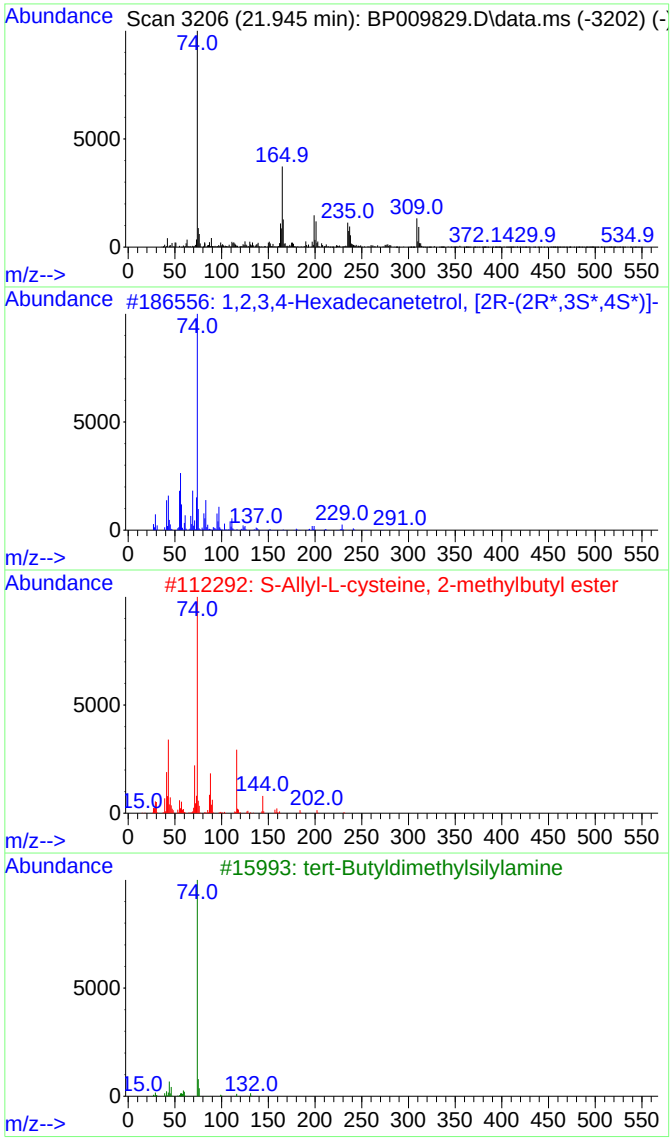
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION

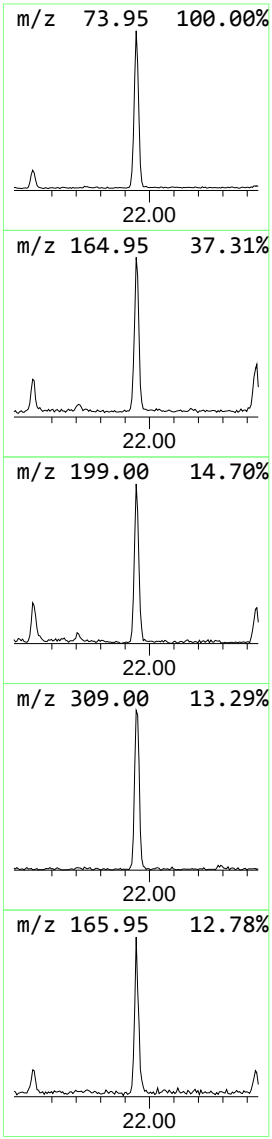
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.945	2.89 ng/ul	389180	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,4-Hexadecanetetrol, [2R-(2...	290	C16H34O4	136953-06-5	17
2	S-Allyl-L-cysteine, 2-methylbuty...	231	C11H21NO2S	1000453-77-6	17
3	tert-Butyldimethylsilylamine	131	C6H17NSi	041879-37-2	9
4	2-(4-Chlorophenoxy)thioacetamide	201	C8H8ClNOS	035368-44-6	9
5	Pentanoic acid, 3-methyl-, methy...	130	C7H14O2	002177-78-8	9





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009829.D
Acq On : 14 Apr 2022 15:49
Operator : CG/JU
Sample : N2293-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.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
Butane, 2-metho...	3.123	23.0	ng/ul	955775	1	7.963	830595	20.0
Butane, 1-bromo-	4.781	2.1	ng/ul	88589	1	7.963	830595	20.0
2-Hexenal, (E)-	8.575	2.7	ng/ul	110493	1	7.963	830595	20.0
Benzene, 1-brom...	9.781	2.0	ng/ul	130643	2	10.769	1272730	20.0
DDMU	19.416	4.8	ng/ul	652256	5	21.422	2693780	20.0
o,p'-DDE	19.804	6.0	ng/ul	810353	5	21.422	2693780	20.0
m,p'-DDD	19.904	5.1	ng/ul	686357	5	21.422	2693780	20.0
Mitotane	20.298	7.8	ng/ul	1051210	5	21.422	2693780	20.0
o,p'-DDT	20.351	3.4	ng/ul	455136	5	21.422	2693780	20.0
p,p'-DDT	20.727	15.4	ng/ul	2073390	5	21.422	2693780	20.0
(DEL) Alkane: S...	21.133	5.0	ng/ul	680660	5	21.422	2693780	20.0
unknown-01	21.945	2.9	ng/ul	389180	5	21.422	2693780	20.0