LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014646.D
Acq On : 11 Apr 2023 09:28
Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB9D0DL

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-BP032823.M
Title : SVOA CALIBRATION

Signal : TIC: BP014646.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	32	35	44	rVB	115370	154288	8.75%	0.980%
2	3.840	109	111	119	rBV3	13198	30593	1.73%	0.194%
3	4.117	154	158	162	rBV3	9863	13387	0.76%	0.085%
4	4.428	206	211	212	rBV5	2604	4230	0.24%	0.027%
5	4.570	230	235	247	rVB	75289	120342	6.82%	0.764%
6	5.264	346	353	357	rBV3	12722	21551	1.22%	0.137%
7	6.446	548	554	558	rBV7	5843	11652	0.66%	0.074%
8	6.493	558	562	569	rVB8	4646	8177	0.46%	0.052%
9	6.840	619	621	626	rVB6	3541	5156	0.29%	0.033%
10	7.046	650	656	660	rBV7	2385	4348	0.25%	0.028%
11	7.311	696	701	712	rVV	33829	68306	3.87%	0.434%
12	7.446	719	724	725	rVV4	4747	6392	0.36%	0.041%
13	7.522	729	737	747	rVB4	18129	47989	2.72%	0.305%
14	7.628	753	755	760	rVB6	4001	5047	0.29%	0.032%
15	7.669	760	762	768	rVB7	3460	5783	0.33%	0.037%
16	7.828	785	789	791	rBV4	5307	7414	0.42%	0.047%
17	7.875	792	797	801	rVB6	5576	10071	0.57%	0.064%
18	7.928	801	806	808	rBV4	7387	9957	0.56%	0.063%
19	7.975	808	814	827	rBV	416675	707545	40.11%	4.492%
20	8.140	838	842	846	rVB7	5688	8082	0.46%	0.051%
21	8.187	846	850	852	rBV4	3807	6161	0.35%	0.039%
22	8.463	890	897	900	rBV6	9029	14292	0.81%	0.091%
23	8.640	920	927	932	rBV9	6810	17104	0.97%	0.109%
24	8.722	932	941	947	rBV9	14958	41354	2.34%	0.263%
25	8.793	952	953	958	rVB4	4750	6032	0.34%	0.038%
26	8.999	982	988	994	rBV4	11471	29197	1.66%	0.185%
27	9.081	997	1002	1010	rVB3	37013	68746	3.90%	0.436%
28	9.163	1010	1016	1026	rBV4	45245	99801	5.66%	0.634%
29	9.275	1032	1035	1038	rBV5	3404	4646	0.26%	0.029%
30	9.316	1038	1042	1054	rVB5	6895	24809	1.41%	0.158%
31	9.422	1057	1060	1065	rVB7	5363	7470	0.42%	0.047%
32	9.499	1069	1073	1082	rBV7	9249	20584	1.17%	0.131%
33	9.605	1086	1091	1100	rBV2	18386	37198	2.11%	0.236%
34	9.728	1109	1112	1116	rVB6	7341	10502	0.60%	0.067%
35	9.887	1132	1139	1146	rBV8	16375	36794	2.09%	0.234%
36	10.034	1162	1164	1169	rVB6	4222	5229	0.30%	0.033%

LSC Area Percent Report

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Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9D0DL

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-BP032823.M
Title : SVOA CALIBRATION

37	10.093	1171	1174	1181	rVB9	5409	10286	0.58%	0.065%
38	10.187	1181	1190	1197	rBV3	18726	45592	2.58%	0.289%
39	10.310	1202	1211	1215	rBV9	9754	21807	1.24%	0.138%
40	10.422	1224	1230	1231	rBV3	8461	12526	0.71%	0.080%
41	10.457	1231	1236	1240	rVV3	14088	28546	1.62%	0.181%
42	10.793	1284	1293	1306	rVV	571735	1064125	60.33%	6.757%
43	10.922	1307	1315	1319	rVV3	16936	36501	2.07%	0.232%
44	10.975	1319	1324	1333	rVB3	23540	62268	3.53%	0.395%
45	11.263	1370	1373	1378	rBV7	4093	6241	0.35%	0.040%
46	11.334	1380	1385	1386	rBV5	4228	5328	0.30%	0.034%
47	11.393	1393	1395	1400	rVB6	4347	5461	0.31%	0.035%
48	11.669	1436	1442	1443	rVV6	3307	4529	0.26%	0.029%
49	11.793	1458	1463	1468	rBV6	10584	17748	1.01%	0.113%
50	12.004	1493	1499	1504	rBV8	8755	17071	0.97%	0.108%
51	12.140	1514	1522	1532	rBV	59592	115411	6.54%	0.733%
52	12.322	1546	1553	1561	rBV9	18768	54797	3.11%	0.348%
53	12.516	1584	1586	1592	rVB6	7597	11466	0.65%	0.073%
54	12.798	1631	1634	1638	rBV6	7293	11178	0.63%	0.071%
55	12.887	1644	1649	1656	rVB6	5661	15313	0.87%	0.097%
56	13.016	1667	1671	1672	rBV4	4904	5921	0.34%	0.038%
57	13.140	1687	1692	1697	rBV2	15612	25617	1.45%	0.163%
58	13.228	1702	1707	1715	rVB5	33063	56198	3.19%	0.357%
59	13.687	1779	1785	1788	rBV8	15383	28111	1.59%	0.178%
60	13.793	1800	1803	1804	rBV3	8166	7010	0.40%	0.045%
61	13.881	1814	1818	1820	rBV5	6185	10658	0.60%	0.068%
62	13.951	1825	1830	1835	rVV2	17685	32597	1.85%	0.207%
63	14.040	1840	1845	1850	rVV	131373	199062	11.29%	1.264%
64	14.087	1850	1853	1857	rVB4	23822	29516	1.67%	0.187%
65	14.140	1860	1862	1866	rVB5	5018	6039	0.34%	0.038%
66	14.187	1866	1870	1875	rBV7	12028	27189	1.54%	0.173%
67	14.322	1888	1893	1900	rBV2	125637	202740	11.49%	1.287%
68	14.551	1927	1932	1938	rBV2	77491	114939	6.52%	0.730%
69	14.628	1939	1945	1952	rVV	915131	1349364	76.50%	8.568%
70	14.692	1952	1956	1962	rVB	28745	47731	2.71%	0.303%
71	15.192	2036	2041	2046	rBV2	55337	100525	5.70%	0.638%
72	15.310	2058	2061	2065	rBV5	12806	12498	0.71%	0.079%
73	15.381	2069	2073	2075	rBV2	9140	14656	0.83%	0.093%
74	15.628	2110	2115	2121	rBV	262852	380077	21.55%	2.413%
75	15.681	2121	2124	2130	rVB3	17736	28921	1.64%	0.184%
76	15.863	2151	2155	2158	rBV	16361	27841	1.58%	0.177%

LSC Area Percent Report

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Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB9D0DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Title : SVOA CALIBRATION

77	15.992	2173	2177	2189	rVB2	62230	115178	6.53%	0.731%
78	16.145	2197	2203	2219	rBV	356744	566830	32.14%	3.599%
79	16.339	2233	2236	2243	rVB2	42473	68204	3.87%	0.433%
80	16.481	2258	2260	2263	rBV3	10498	11085	0.63%	0.070%
81	16.663	2287	2291	2300	rBV	159069	252398	14.31%	1.603%
82	17.175	2374	2378	2380	rBV3	21208	30917	1.75%	0.196%
83	17.392	2409	2415	2420	rBV	1028633	1460692	82.81%	9.275%
84	17.434	2420	2422	2427	rVV	99628	124017	7.03%	0.787%
85	17.492	2427	2432	2441	rVV2	184178	291843	16.55%	1.853%
86	18.263	2559	2563	2569	rBV3	86931	141999	8.05%	0.902%
87	18.375	2579	2582	2585	rVB2	29562	32640	1.85%	0.207%
88	18.475	2596	2599	2608	rVB4	87950	121318	6.88%	0.770%
89	18.598	2618	2620	2623	rBV4	17869	20845	1.18%	0.132%
90	18.639	2623	2627	2633	rVB	304973	382453	21.68%	2.428%
91	18.728	2640	2642	2647	rBV6	18753	23761	1.35%	0.151%
92	18.816	2653	2657	2662	rBV	286171	365030	20.69%	2.318%
93	19.398	2753	2756	2759	rBV	72106	94785	5.37%	0.602%
94	19.728	2807	2812	2815	rBV	261295	380588	21.58%	2.417%
95	19.769	2815	2819	2827	rVB	450529	667544	37.85%	4.239%
96	20.145	2879	2883	2890	rBV	197821	252491	14.31%	1.603%
97	21.363	3086	3090	3095	rVB	598614	626580	35.52%	3.978%
98	21.480	3106	3110	3120	rBV	1122853	1559255	88.40%	9.900%
99	23.821	3505	3508	3522	rVB2	245629	461555	26.17%	2.931%
100	23.986	3530	3536	3550	rVB2	817206	1763871	100.00%	11.200%

Sum of corrected areas: 15749512

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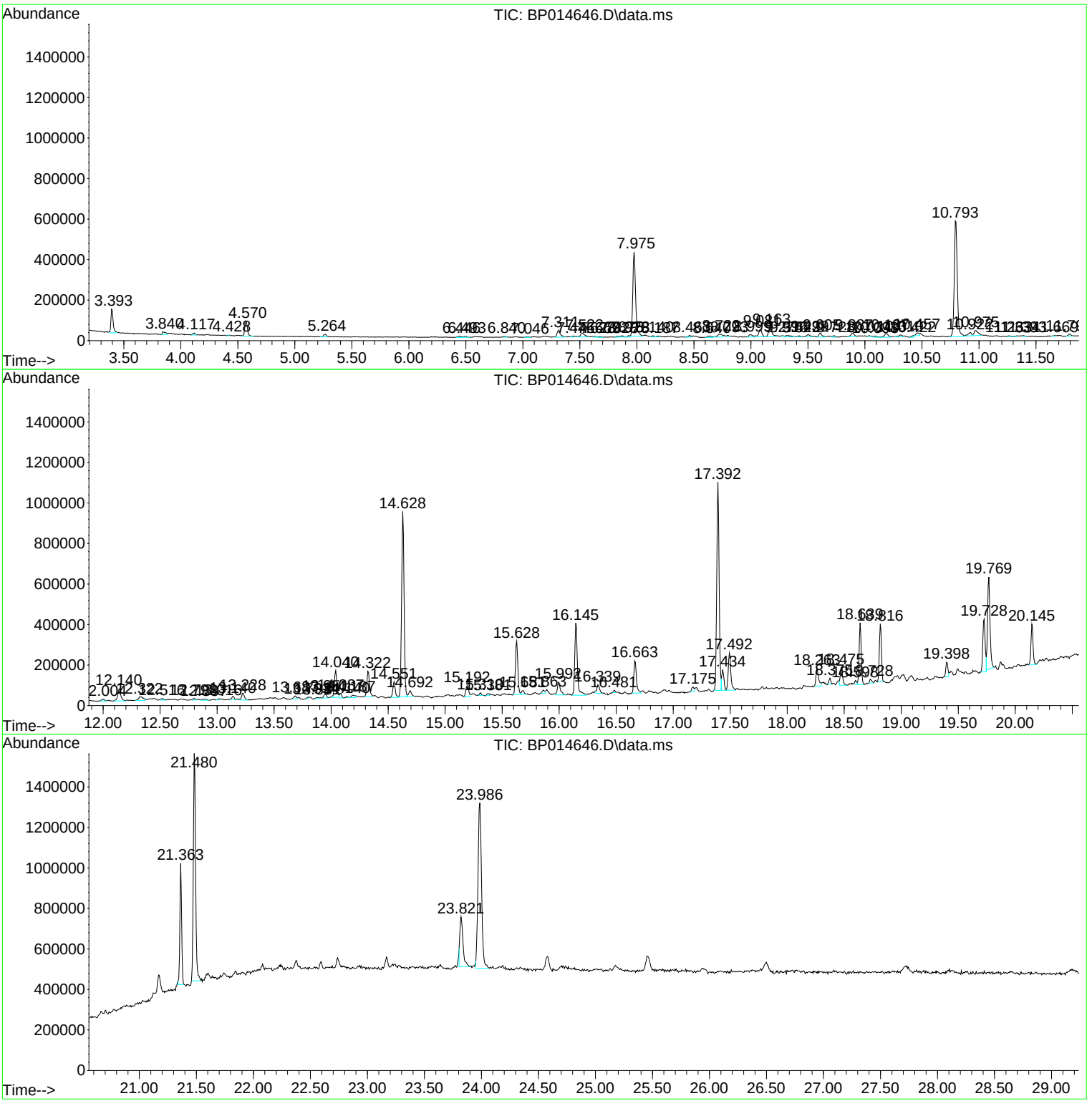
Instrument :
BNA_P
ClientSampleId :
CB9D0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014646.D
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

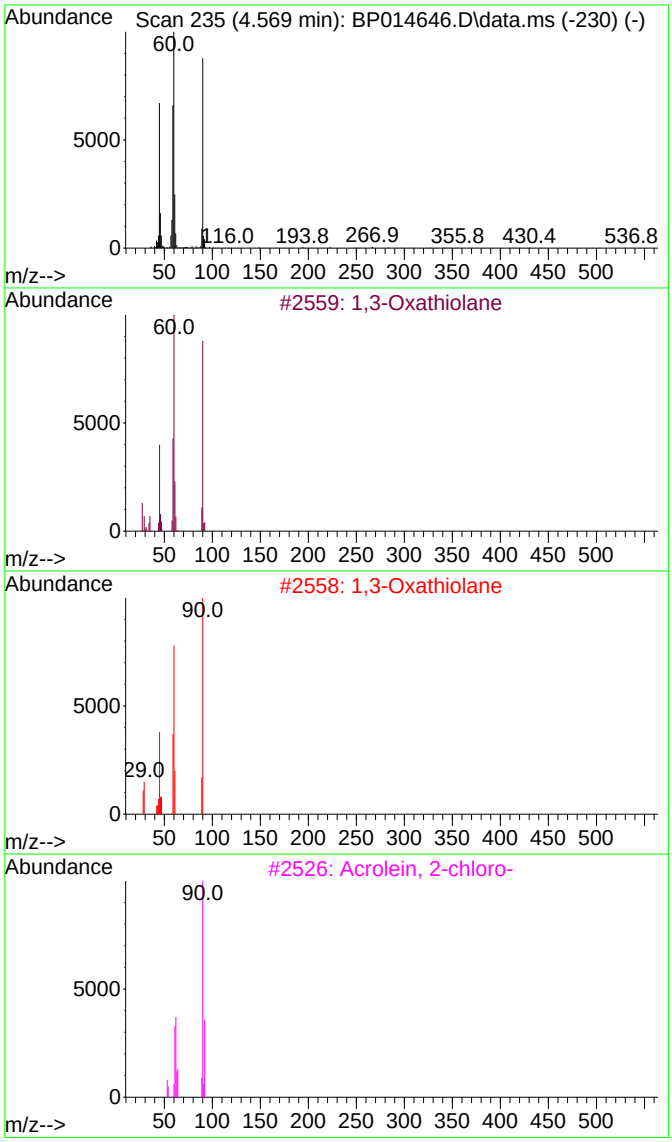
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

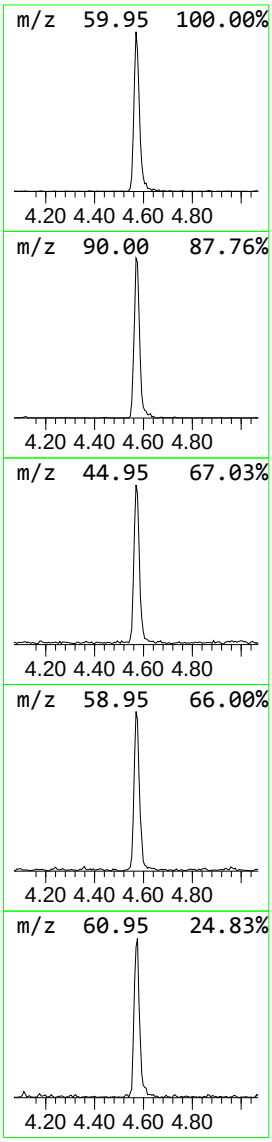
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.569	3.40 ng/ul	120342	1,4-Dichlorobenzene-d4	7.975

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	87
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	64
3	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	59
4	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	59
5	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014646.D
Acq On : 11 Apr 2023 09:28
Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9D0DL

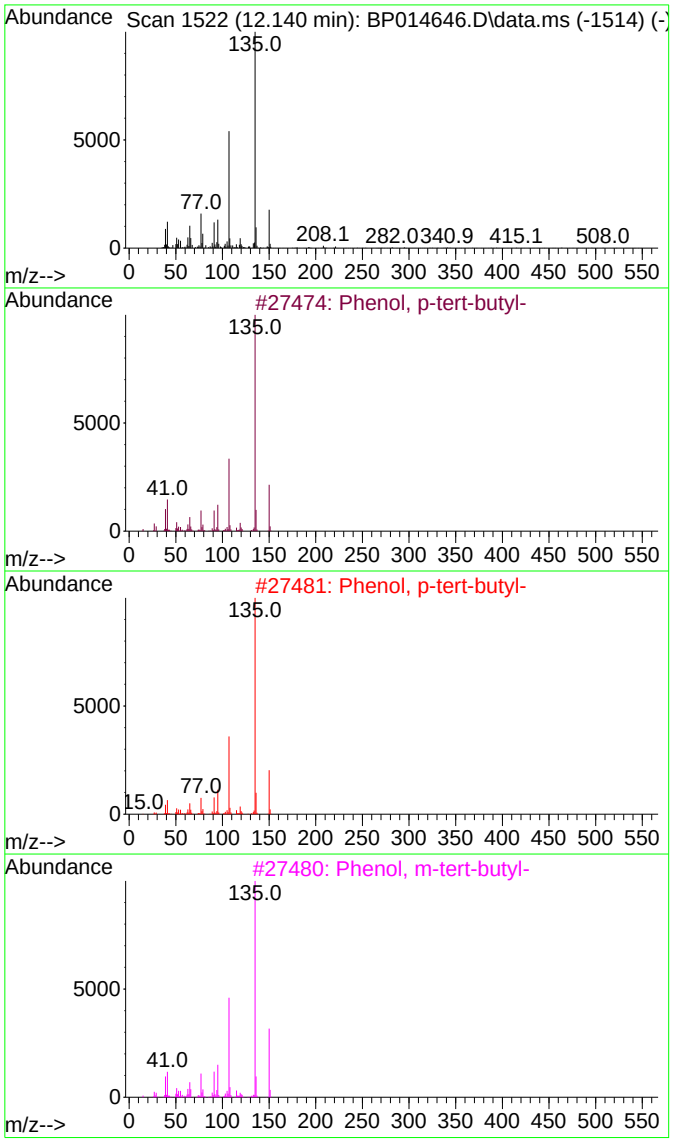
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

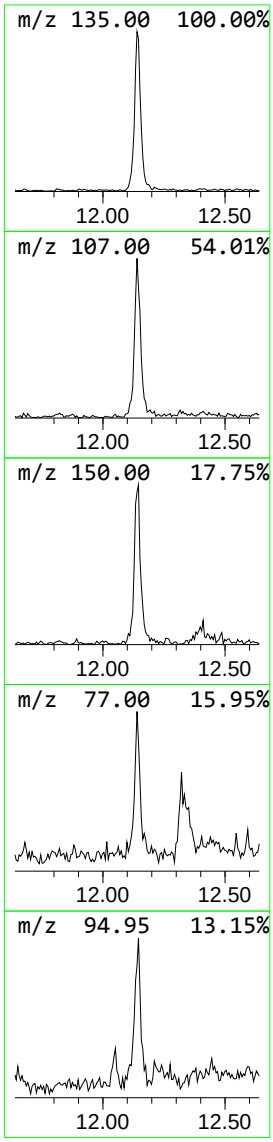
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenol, p-tert-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.140	2.17 ng/ul	115411	Naphthalene-d8	10.799

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	93
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	93
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	87
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
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Instrument :
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CB9D0DL

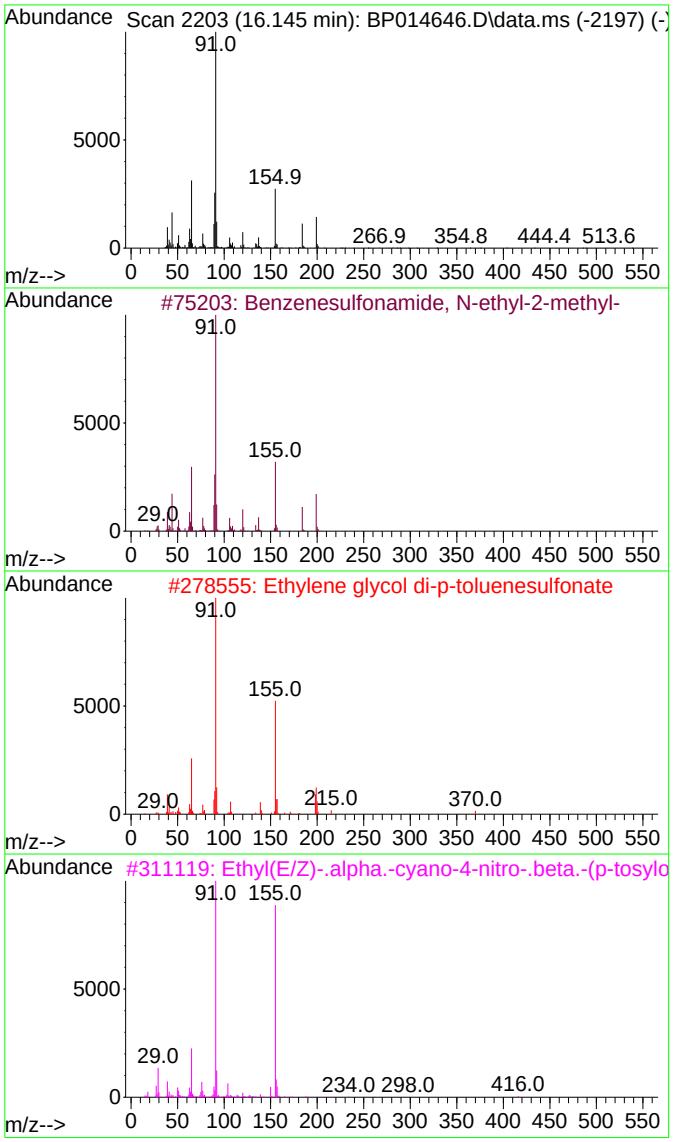
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

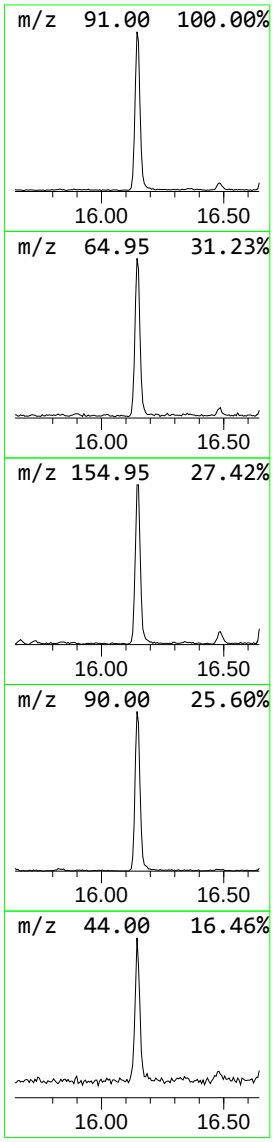
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.145	7.76 ng/ul	566830	Phenanthrene-d10	17.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Ethylene glycol di-p-toluenesulf...	370	C16H18O6S2	006315-52-2	59
3	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	58
4	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47
5	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	46





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ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB9D0DL

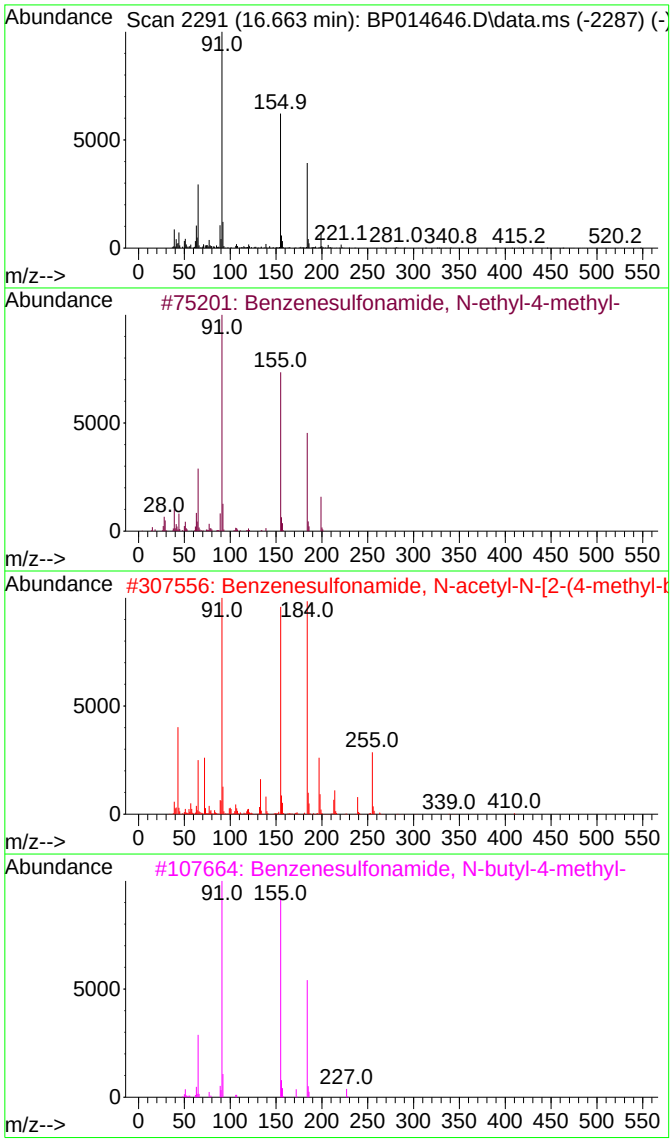
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

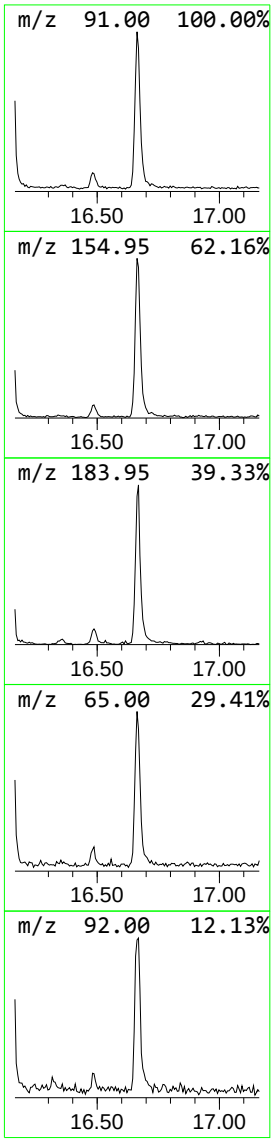
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzenesulfonamide, N-ethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.663	3.46 ng/ul	252398	Phenanthrene-d10	17.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	91
2	Benzenesulfonamide, N-acetyl-N-[...	410	C18H22N2O5S2	1000301-10-0	90
3	Benzenesulfonamide, N-butyl-4-me...	227	C11H17NO2S	001907-65-9	86
4	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	80
5	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	72





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Data File : BP014646.D
Acq On : 11 Apr 2023 09:28
Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CB9D0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

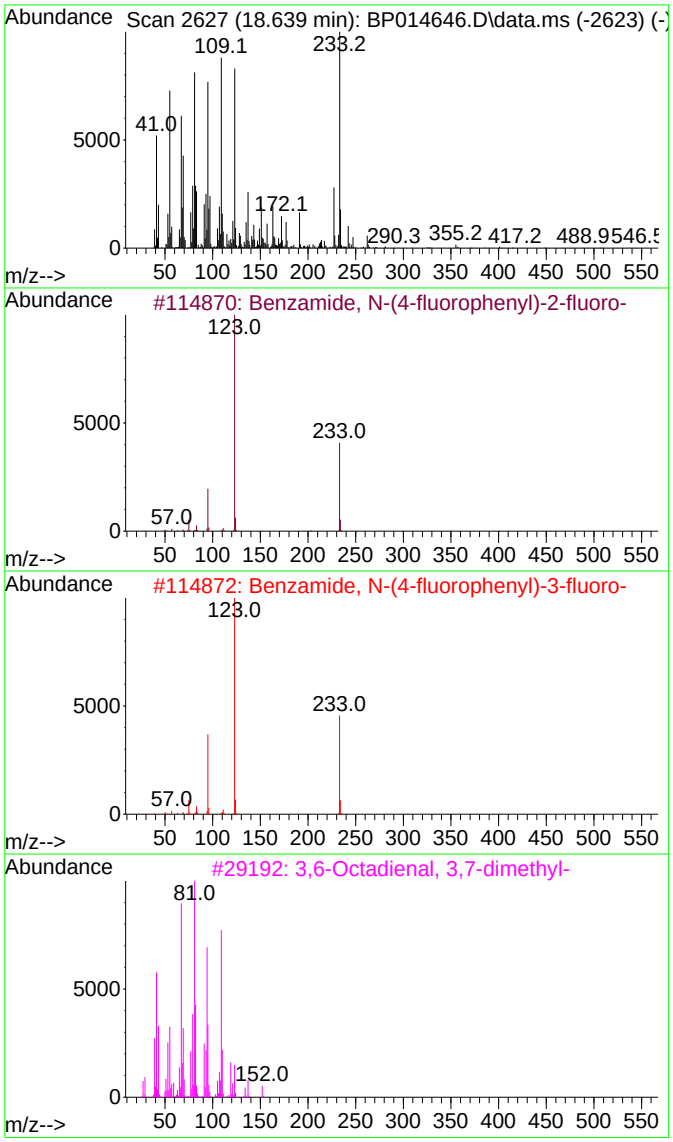
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

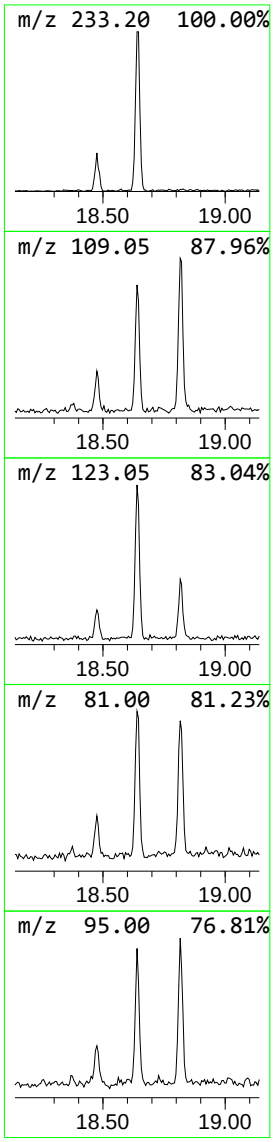
****

Peak Number 5 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.639	5.24 ng/ul	382453	Phenanthrene-d10	17.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzamide, N-(4-fluorophenyl)-2-...	233	C13H9F2NO	1000307-07-9	45
2	Benzamide, N-(4-fluorophenyl)-3-...	233	C13H9F2NO	1010307-16-3	38
3	3,6-Octadienal, 3,7-dimethyl-	152	C10H16O	055722-59-3	35
4	3-Octadecyne	250	C18H34	061886-64-4	30
5	Methyl ethyl ketone, N-(3-methyl...	233	C12H15N3S	1000292-93-3	27





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Acq On : 11 Apr 2023 09:28
Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9D0DL

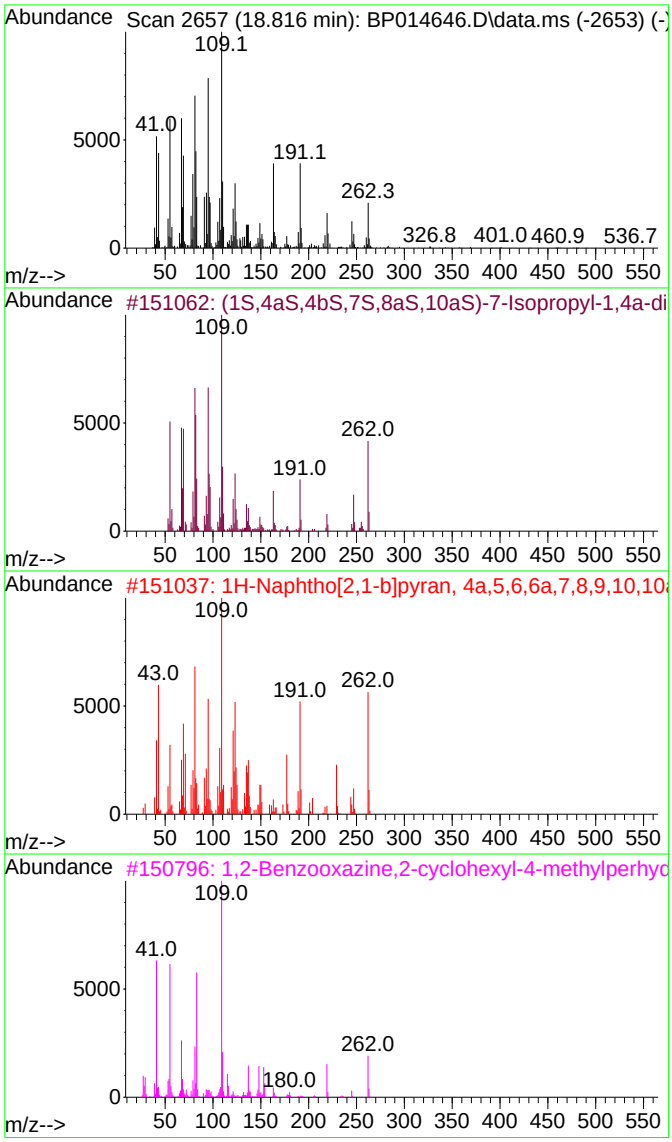
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

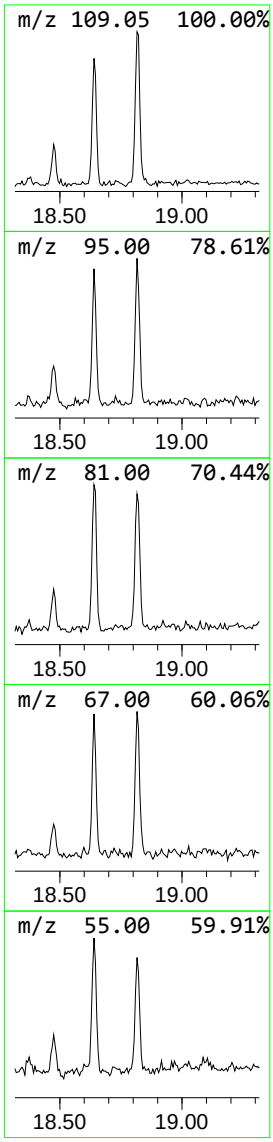
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (1S,4aS,4bS,7S,8aS,10aS)-7-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.816	5.00 ng/ul	365030	Phenanthrene-d10	17.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S,4aS,4bS,7S,8aS,10aS)-7-Isopr...	262	C19H34	002221-95-6	90
2	1H-Naphtho[2,1-b]pyran, 4a,5,6,6...	262	C18H30O	005153-92-4	38
3	1,2-Benzooxazine,2-cyclohexyl-4-...	262	C16H26N2O	037898-39-8	25
4	7-Hexadecyne	222	C16H30	074685-28-2	22
5	3-[(4-Fluorophenyl)methyl]-1,2,3...	192	C9H9FN4	1000435-39-8	18





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Data File : BP014646.D
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Operator : CG/JU
Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9D0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

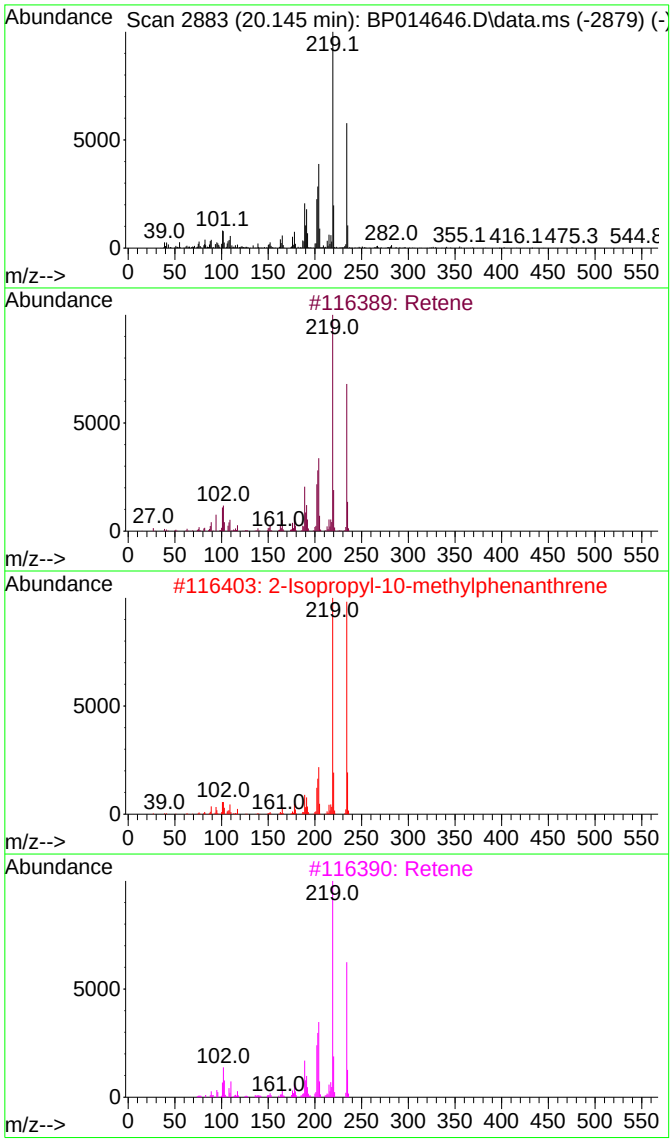
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

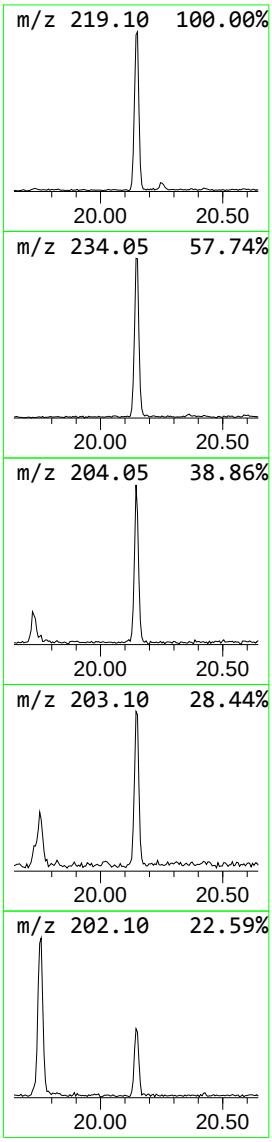
****

Peak Number 7 Retene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.145	3.24 ng/ul	252491	Chrysene-d12	21.480

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Retene	234	C18H18	000483-65-8	97
2	2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	97
3	Retene	234	C18H18	000483-65-8	94
4	Retene	234	C18H18	000483-65-8	91
5	1,4-Di-(4'-methylphenyl)buta-1,3...	234	C18H18	1000202-17-5	83





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Sample : 02226-01DL 5X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9D0DL

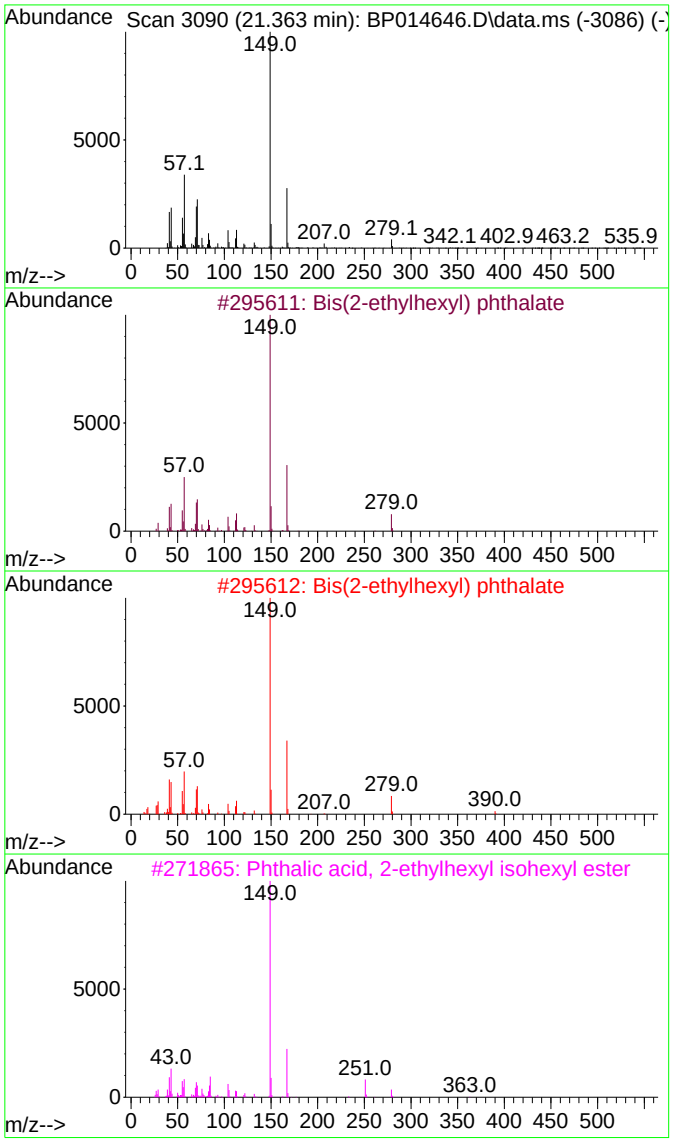
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

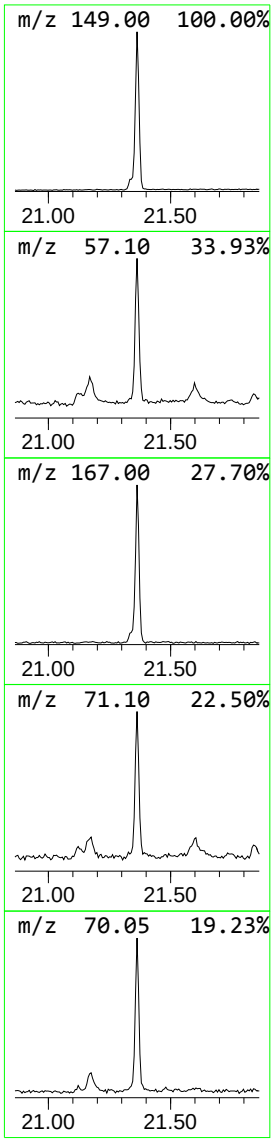
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Bis(2-ethylhexyl) phthalate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.363	8.04 ng/ul	626580	Chrysene-d12	21.480

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	90
2	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	90
3	Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	59
4	Phthalic acid, di(oct-3-yl) ester	390	C24H38O4	1000377-72-3	59
5	Phthalic acid, di(2-methylbutyl)...	306	C18H26O4	1000322-82-5	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014646.D
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BNA_P
ClientSampleId :
CB9D0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.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
1,3-Oxathiolane	4.569	3.4	ng/ul	120342	1	7.975	707545	20.0
Phenol, p-tert-...	12.140	2.2	ng/ul	115411	2	10.799	1064130	20.0
Benzenesulfonam...	16.145	7.8	ng/ul	566830	4	17.392	1460690	20.0
Benzenesulfonam...	16.663	3.5	ng/ul	252398	4	17.392	1460690	20.0
unknown-01	18.639	5.2	ng/ul	382453	4	17.392	1460690	20.0
(1S,4aS,4bS,7S,...	18.816	5.0	ng/ul	365030	4	17.392	1460690	20.0
Retene	20.145	3.2	ng/ul	252491	5	21.480	1559260	20.0
Bis(2-ethylhexy...	21.363	8.0	ng/ul	626580	5	21.480	1559260	20.0