LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

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-BP102522.M
Title : SVOA CALIBRATION

Signal : TIC: BP012414.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	4	7	20	rVB	1862434	2765185	24.65%	2.065%
2	3.240	39	43	46	rBV	62454	84653	0.75%	0.063%
3	3.663	112	115	126	rBV	296539	536615	4.78%	0.401%
4	3.875	145	151	159	rBV2	59695	109122	0.97%	0.082%
5	3.975	163	168	177	rVB	74983	116720	1.04%	0.087%
6	5.099	353	359	373	rVV	1298839	2106720	18.78%	1.574%
7	5.228	376	381	386	rVV	70215	104402	0.93%	0.078%
8	5.293	386	392	400	rVB	107741	173770	1.55%	0.130%
9	5.422	408	414	424	rBV	1242297	2668345	23.79%	1.993%
10	5.504	424	428	436	rVB	80615	127401	1.14%	0.095%
11	5.793	471	477	491	rBV	1123154	1760308	15.69%	1.315%
12	6.063	513	523	535	rBV5	25551	81066	0.72%	0.061%
13	6.269	549	558	564	rBV	381625	645060	5.75%	0.482%
14	6.457	583	590	595	rBV	164743	253012	2.26%	0.189%
15	6.510	595	599	611	rVB	85230	169202	1.51%	0.126%
16	6.763	636	642	647	rBV	146828	241932	2.16%	0.181%
17	6.869	654	660	666	rBV	79218	127619	1.14%	0.095%
18	6.963	671	676	681	rBV	260047	425068	3.79%	0.317%
19	7.004	681	683	688	rVB2	86402	109310	0.97%	0.082%
20	7.128	696	704	710	rBV	1325793	2147543	19.15%	1.604%
21	7.322	730	737	750	rVV2	1234990	2467887	22.00%	1.843%
22	7.434	750	756	761	rVV	268335	423498	3.78%	0.316%
23	7.487	761	765	771	rVV3	56776	106592	0.95%	0.080%
24	7.792	805	817	829	rBV	1362706	2390857	21.32%	1.786%
25	7.898	829	835	840	rVV	124424	211855	1.89%	0.158%
26	7.987	846	850	856	rVB2	145477	216056	1.93%	0.161%
27	8.151	871	878	883	rBV2	81954	184181	1.64%	0.138%
28	8.504	932	938	949	rBV	928178	1569144	13.99%	1.172%
29	8.787	981	986	995	rBV3	64575	122337	1.09%	0.091%
30	8.892	995	1004	1008	rBV2	51688	98758	0.88%	0.074%
31	8.957	1008	1015	1023	rVV	1521388	2671279	23.82%	1.995%
32	9.022	1023	1026	1036	rVV4	120309	265001	2.36%	0.198%
33	9.104	1036	1040	1046	rVB	66960	99276	0.89%	0.074%
34	9.410	1086	1092	1098	rVB4	41986	79832	0.71%	0.060%
35	9.581	1116	1121	1125	rBV	90387	138013	1.23%	0.103%
36	9.681	1131	1138	1151	rVB2	1357772	2487547	22.18%	1.858%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BHAC2

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-BP102522.M
Title : SVOA CALIBRATION

37	9.834	1158	1164	1168	rBV3	49061	85206	0.76%	0.064%
38	9.957	1180	1185	1189	rBV	107726	187328	1.67%	0.140%
39	10.010	1189	1194	1202	rVB4	126454	270480	2.41%	0.202%
40	10.086	1202	1207	1211	rBV2	70122	115297	1.03%	0.086%
41	10.134	1211	1215	1221	rVB2	46211	80145	0.71%	0.060%
42	10.216	1221	1229	1239	rBV	1929330	3395393	30.27%	2.536%
43	10.404	1257	1261	1266	rVB2	65179	95763	0.85%	0.072%
44	10.592	1286	1293	1299	rVV	2053185	3536580	31.53%	2.641%
45	10.639	1299	1301	1305	rVV	129543	201708	1.80%	0.151%
46	10.686	1305	1309	1311	rVV2	158612	258775	2.31%	0.193%
47	10.739	1311	1318	1331	rVB2	2005450	4105047	36.60%	3.066%
48	10.951	1347	1354	1360	rBV5	68003	201740	1.80%	0.151%
49	11.004	1360	1363	1368	rVB3	50538	82975	0.74%	0.062%
50	11.186	1390	1394	1400	rVB2	57536	101646	0.91%	0.076%
51	11.286	1405	1411	1420	rVB2	197381	357800	3.19%	0.267%
52	11.616	1462	1467	1475	rBV5	56715	108955	0.97%	0.081%
53	11.945	1513	1523	1531	rBV	1444190	2449985	21.84%	1.830%
54	12.416	1595	1603	1610	rVV5	48681	121907	1.09%	0.091%
55	12.598	1629	1634	1639	rVV	50624	83353	0.74%	0.062%
56	12.698	1647	1651	1656	rVV4	69275	113951	1.02%	0.085%
57	12.980	1685	1699	1705	rBV2	463458	1044400	9.31%	0.780%
58	13.298	1747	1753	1757	rBV	323982	507321	4.52%	0.379%
59	13.857	1841	1848	1856	rBV	3599175	5307104	47.32%	3.964%
60	13.922	1856	1859	1863	rVV	124146	152669	1.36%	0.114%
61	13.986	1868	1870	1878	rVV	73942	107988	0.96%	0.081%
62	14.133	1889	1895	1904	rVV	4183502	6438054	57.40%	4.809%
63	14.233	1907	1912	1926	rVB2	94041	217440	1.94%	0.162%
64	14.445	1940	1948	1963	rBV	3285470	5226387	46.60%	3.904%
65	14.674	1981	1987	1997	rVV	232143	482684	4.30%	0.361%
66	15.022	2040	2046	2061	rVB	124096	367624	3.28%	0.275%
67	15.192	2071	2075	2080	rBV	179871	241112	2.15%	0.180%
68	15.439	2111	2117	2131	rVB	5501340	8072351	71.97%	6.029%
69	15.574	2134	2140	2148	rBV	1893403	2630912	23.46%	1.965%
70	15.963	2200	2206	2218	rBV4	158573	356113	3.18%	0.266%
71	16.086	2222	2227	2233	rVV	316075	525692	4.69%	0.393%
72	16.145	2233	2237	2244	rVB	139944	241337	2.15%	0.180%
73	16.504	2290	2298	2303	rBV3	398495	678838	6.05%	0.507%
74	16.586	2309	2312	2322	rVB8	41140	111838	1.00%	0.084%
75	16.945	2370	2373	2383	rVB4	59685	109957	0.98%	0.082%
76	17.168	2406	2411	2413	rBV	563516	790149	7.04%	0.590%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

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-BP102522.M
Title : SVOA CALIBRATION

77	17.204	2413	2417	2428	rVB	4296684	6076360	54.18%	4.538%
78	17.304	2428	2434	2442	rBV2	6479995	9145536	81.54%	6.831%
79	17.433	2452	2456	2461	rBV	59737	92581	0.83%	0.069%
80	17.839	2520	2525	2529	rBV	191723	265493	2.37%	0.198%
81	17.892	2529	2534	2539	rVV	241433	386180	3.44%	0.288%
82	18.092	2565	2568	2577	rVB	124539	171218	1.53%	0.128%
83	18.245	2589	2594	2602	rVB	295056	398284	3.55%	0.297%
84	18.339	2606	2610	2613	rVV	257703	345712	3.08%	0.258%
85	18.386	2613	2618	2623	rVB	471061	788358	7.03%	0.589%
86	18.539	2640	2644	2647	rBV	94768	128134	1.14%	0.096%
87	18.762	2678	2682	2690	rBV3	75034	143812	1.28%	0.107%
88	19.121	2739	2743	2749	rBV2	125607	176923	1.58%	0.132%
89	19.192	2751	2755	2764	rVV	105676	177859	1.59%	0.133%
90	19.268	2764	2768	2772	rVV3	131877	175550	1.57%	0.131%
91	19.327	2775	2778	2792	rVB3	149176	267583	2.39%	0.200%
92	19.545	2809	2815	2820	rBV	8310509	11124948	99.19%	8.309%
93	19.592	2820	2823	2832	rVB	313168	458714	4.09%	0.343%
94	19.762	2849	2852	2857	rVB2	66043	107181	0.96%	0.080%
95	19.839	2862	2865	2870	rVB	69343	86551	0.77%	0.065%
96	19.986	2887	2890	2895	rVV	138387	185422	1.65%	0.138%
97	20.039	2896	2899	2906	rVB6	74204	110679	0.99%	0.083%
98	21.298	3108	3113	3124	rBV	5852766	7250027	64.64%	5.415%
99	23.533	3486	3493	3505	rVB2	5547955	11215947	100.00%	8.377%
100	23.686	3513	3519	3529	rVB2	3495258	6759229	60.26%	5.048%

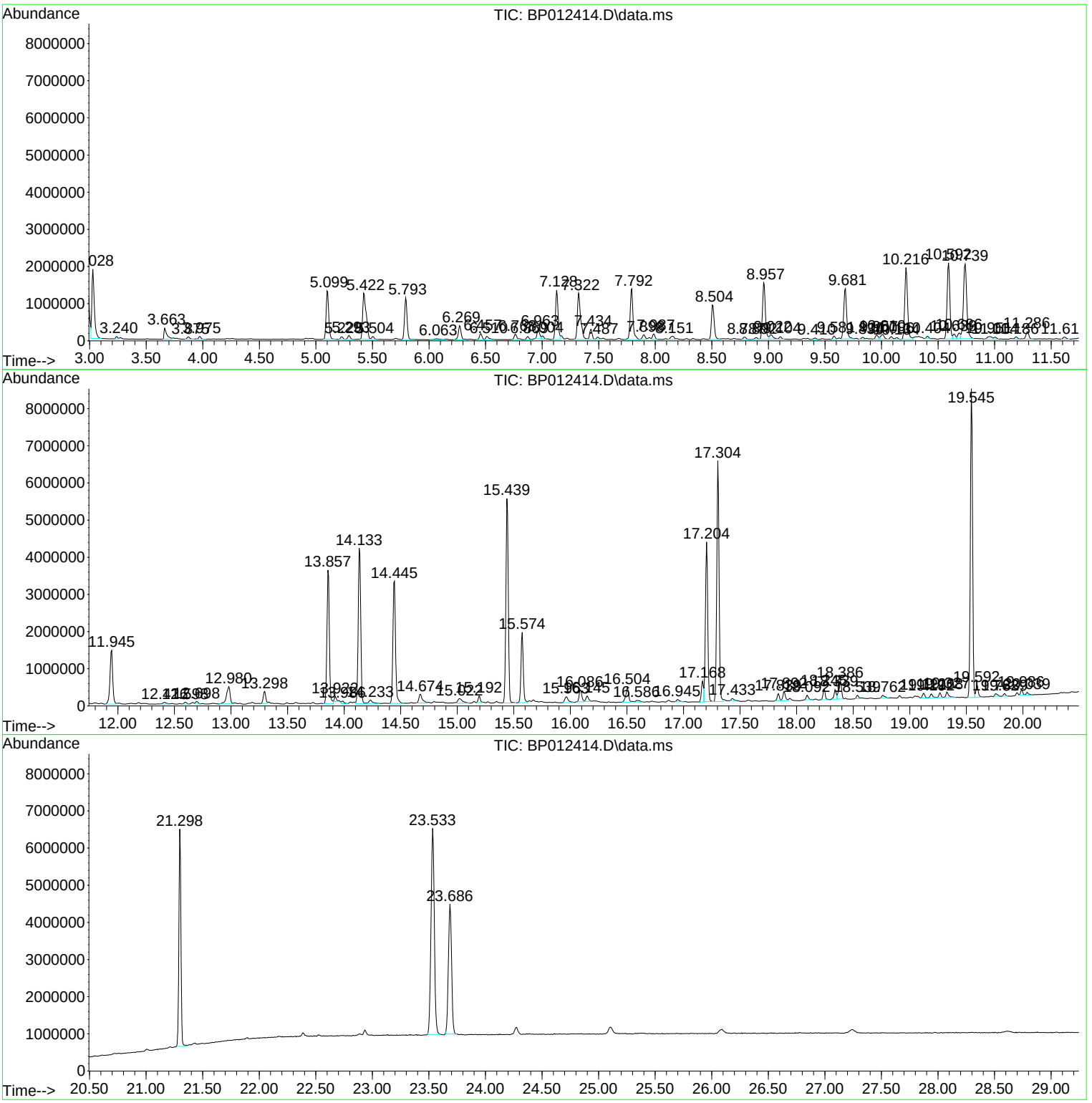
Sum of corrected areas: 133887451

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

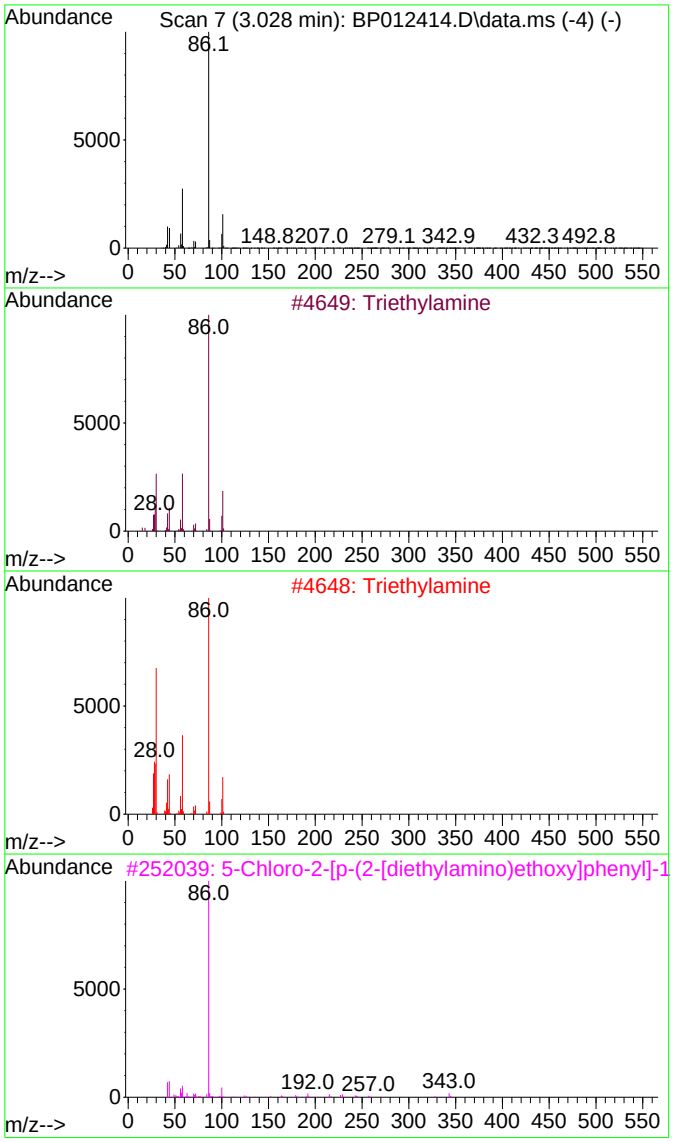
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

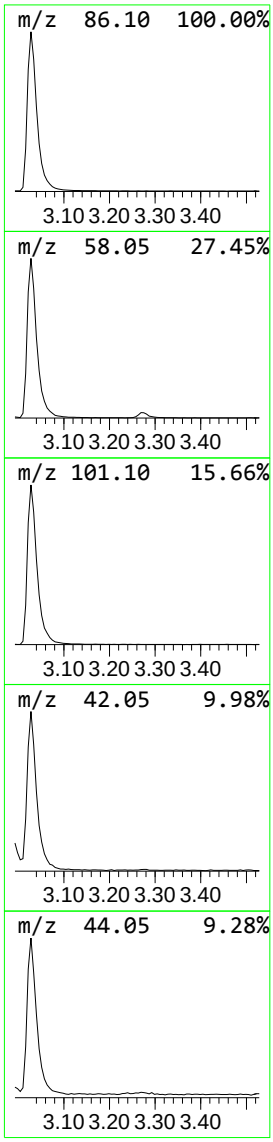
****

Peak Number 1 Triethylamine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	23.13 ng/ul	2765190	1,4-Dichlorobenzene-d4	7.792

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethylamine	101	C6H15N	000121-44-8	91
2	Triethylamine	101	C6H15N	000121-44-8	70
3	5-Chloro-2-[p-(2-[diethylamino)e...	343	C19H22ClN3O	1000251-32-3	64
4	Tetraethyl ammonium fluoride	149	C8H20FN	000665-46-3	56
5	1,2-Ethanediamine, N,N-diethyl-	116	C6H16N2	000100-36-7	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

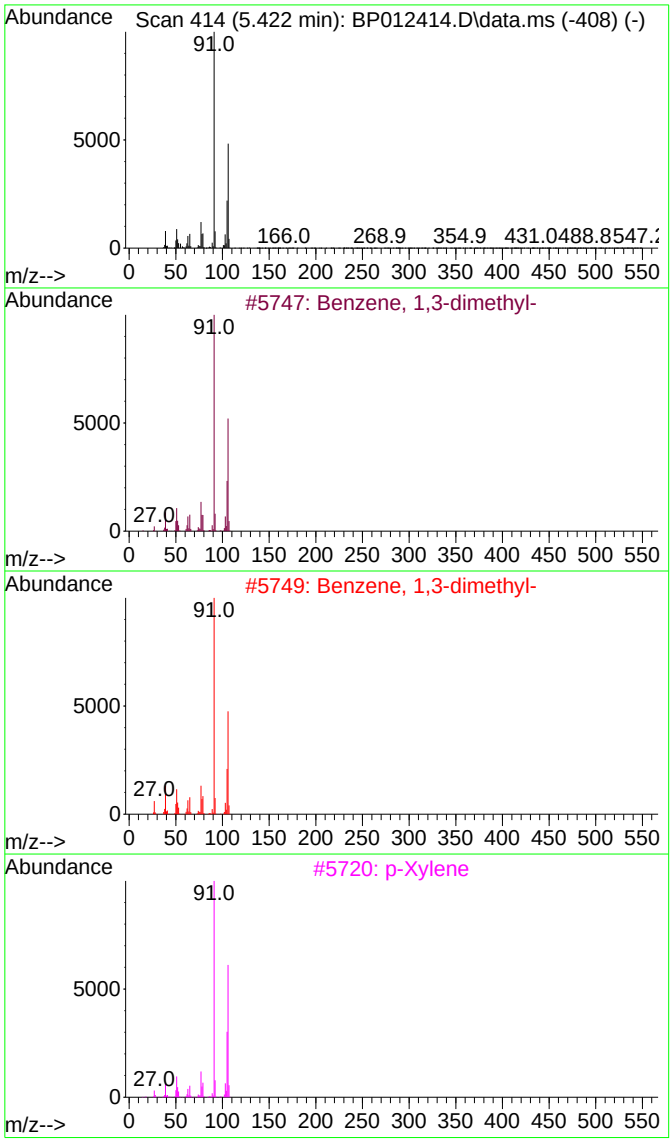
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

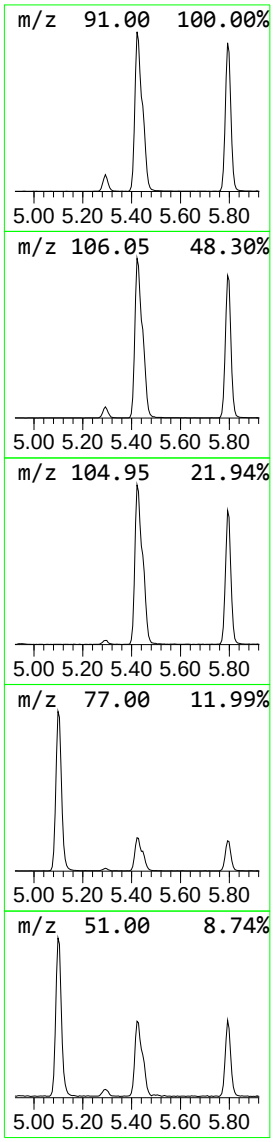
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.422	22.32 ng/ul	2668350	1,4-Dichlorobenzene-d4	7.792

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	p-Xylene	106	C8H10	000106-42-3	97
4	o-Xylene	106	C8H10	000095-47-6	97
5	p-Xylene	106	C8H10	000106-42-3	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

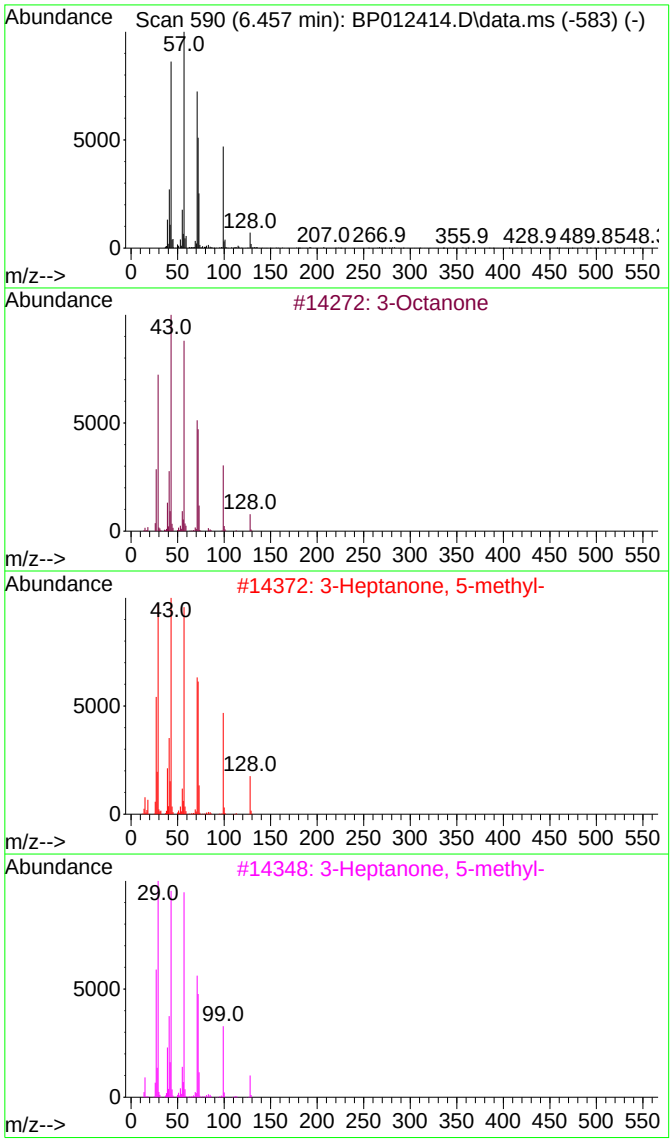
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

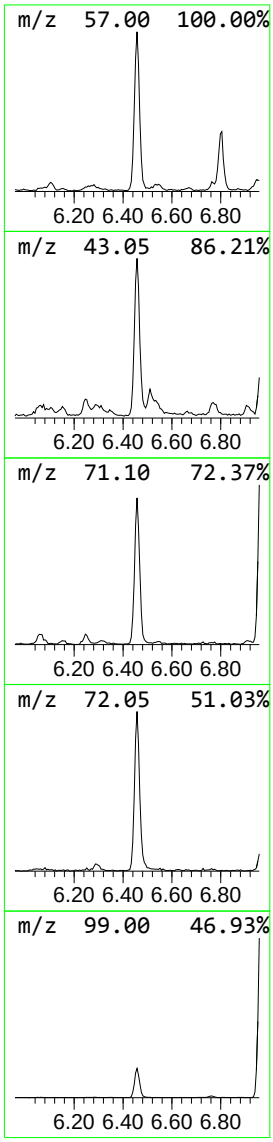
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Octanone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.457	2.12 ng/ul	253012	1,4-Dichlorobenzene-d4	7.792

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanone	128	C8H16O	000106-68-3	72
2	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	64
3	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	64
4	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	64
5	3-Octanone	128	C8H16O	000106-68-3	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

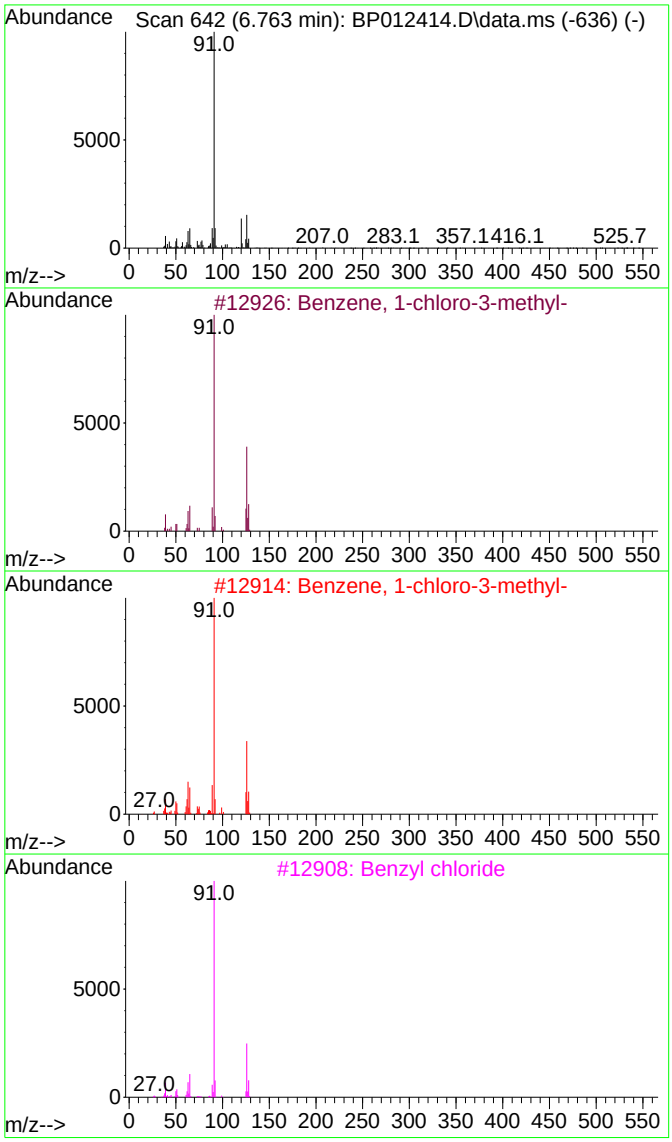
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

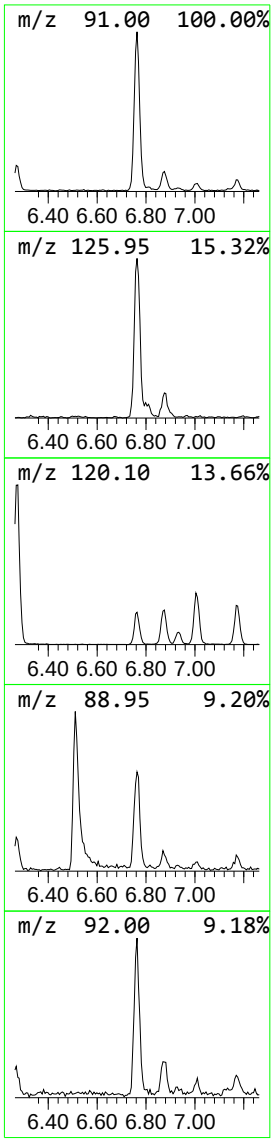
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-chloro-3-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.763	2.02 ng/ul	241932	1,4-Dichlorobenzene-d4	7.792

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	83
2	Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	81
3	Benzyl chloride	126	C7H7Cl	000100-44-7	74
4	Benzyl chloride	126	C7H7Cl	000100-44-7	74
5	Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

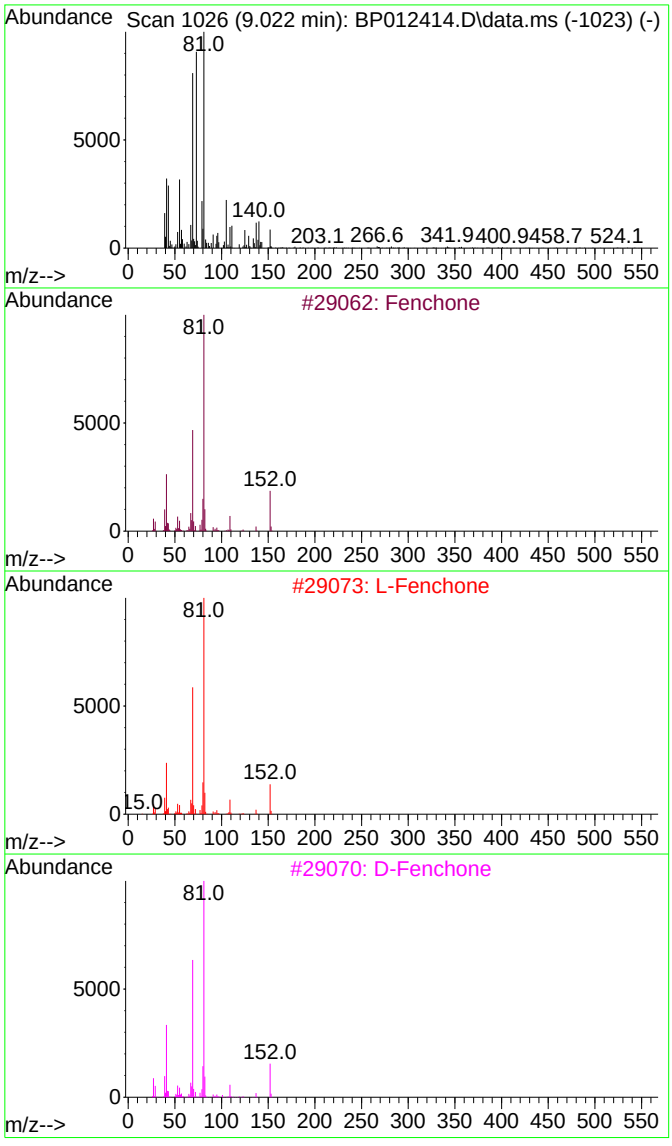
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

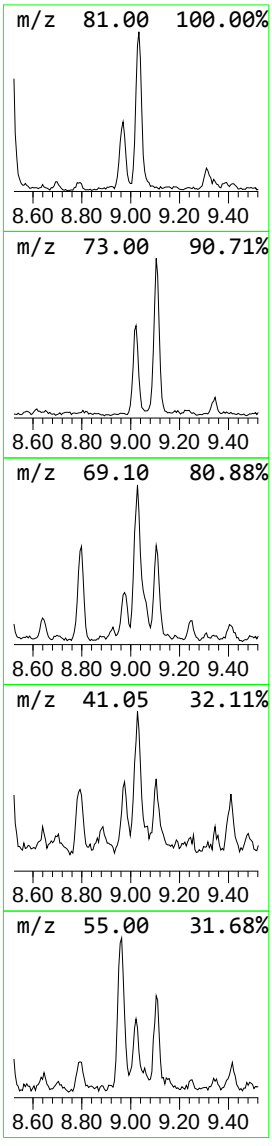
****

Peak Number 9 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.022	2.22 ng/ul	265001	1,4-Dichlorobenzene-d4	7.792

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fenchone	152	C10H16O	001195-79-5	38
2	L-Fenchone	152	C10H16O	007787-20-4	38
3	D-Fenchone	152	C10H16O	004695-62-9	38
4	Fenchone	152	C10H16O	001195-79-5	30
5	Fenchone	152	C10H16O	001195-79-5	25





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Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

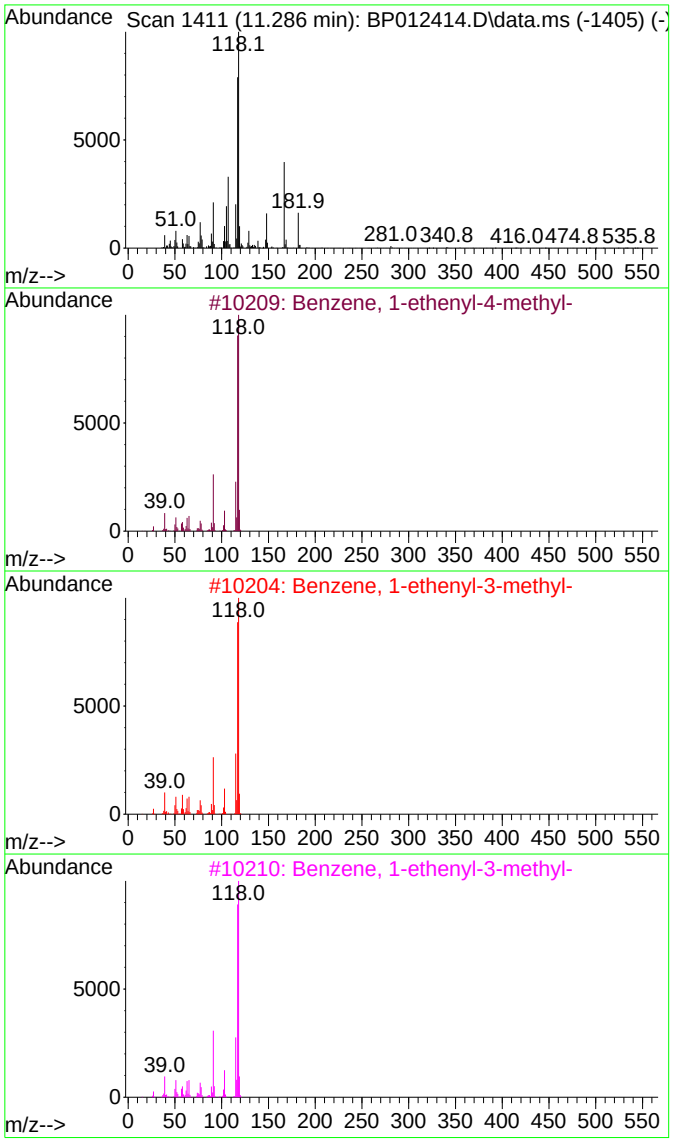
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

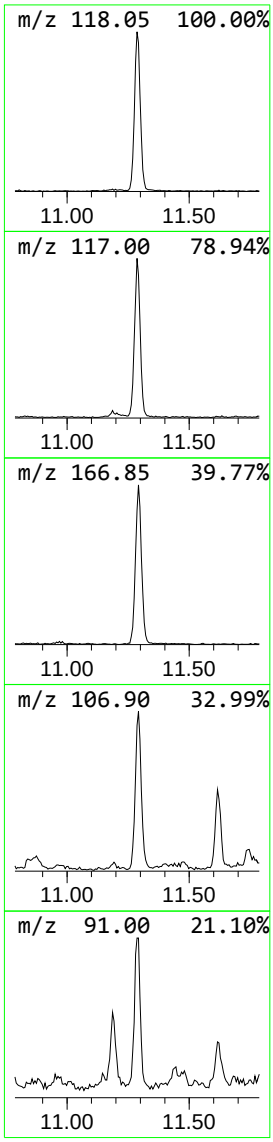
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethenyl-4-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.286	2.02 ng/ul	357800	Naphthalene-d8	10.592

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	60
2	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	53
3	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	53
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	49
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	49





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Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

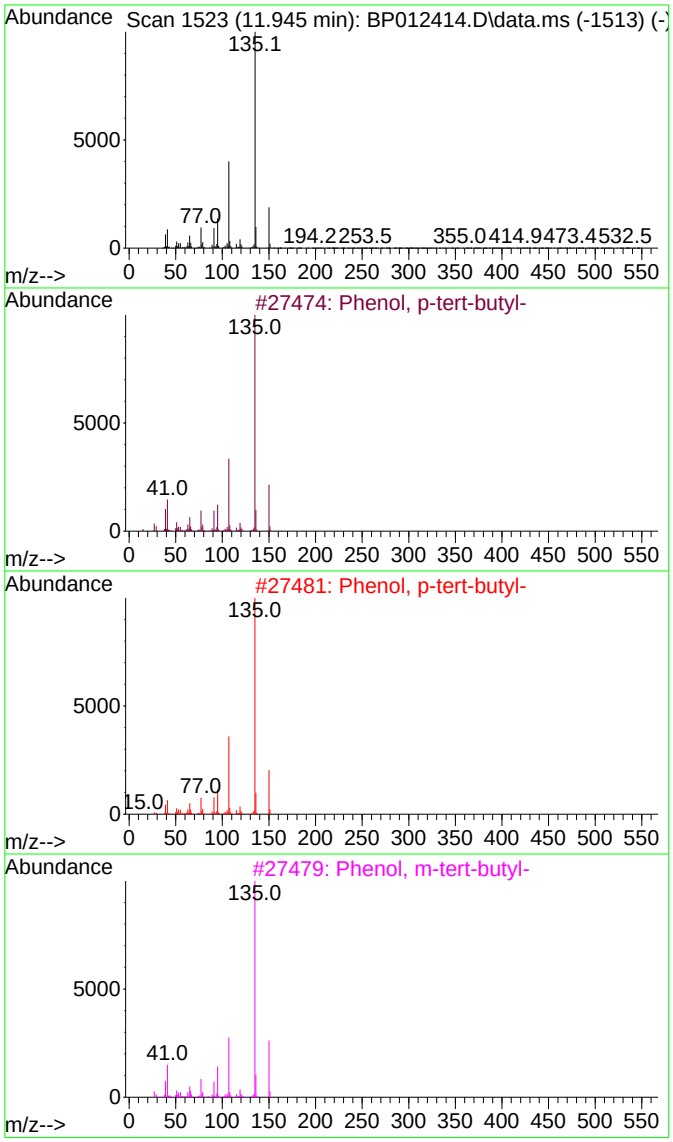
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

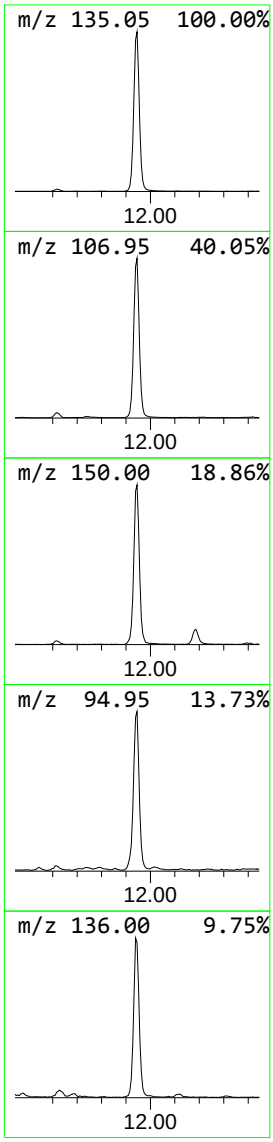
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Phenol, p-tert-butyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.945	13.86 ng/ul	2449990	Naphthalene-d8	10.592

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91





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Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

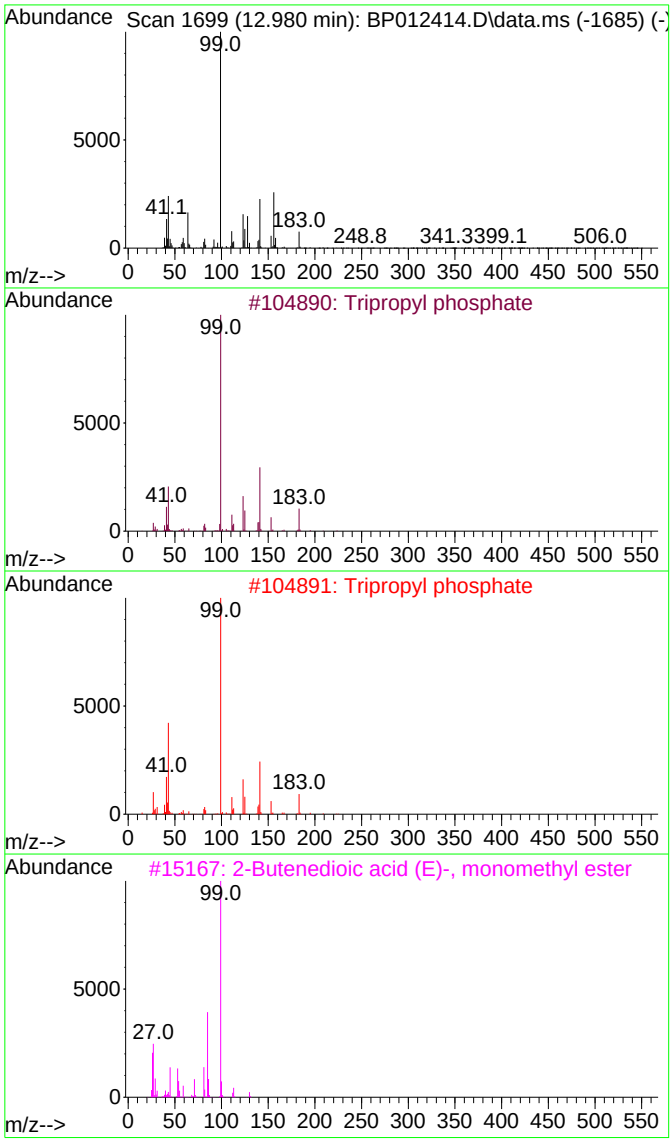
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

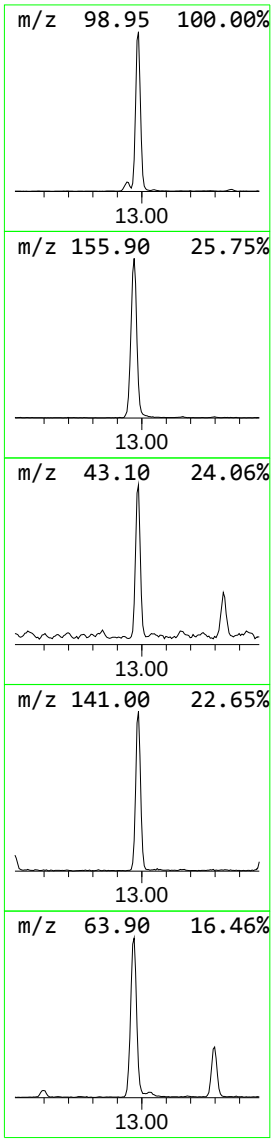
****

Peak Number 12 Tripropyl phosphate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	4.00 ng/ul	1044400	Acenaphthene-d10	14.445

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tripropyl phosphate	224	C9H21O4P	000513-08-6	58
2	Tripropyl phosphate	224	C9H21O4P	000513-08-6	35
3	2-Butenedioic acid (E)-, monomet...	130	C5H6O4	002756-87-8	25
4	Spiro[1,3-dioxolane-2,2'(1'H)-na...	196	C12H20O2	006857-86-9	23
5	4,4'-Bi-1,3,2-dioxaborolane, 2,2...	198	C8H16B2O4	062337-94-4	17





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Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

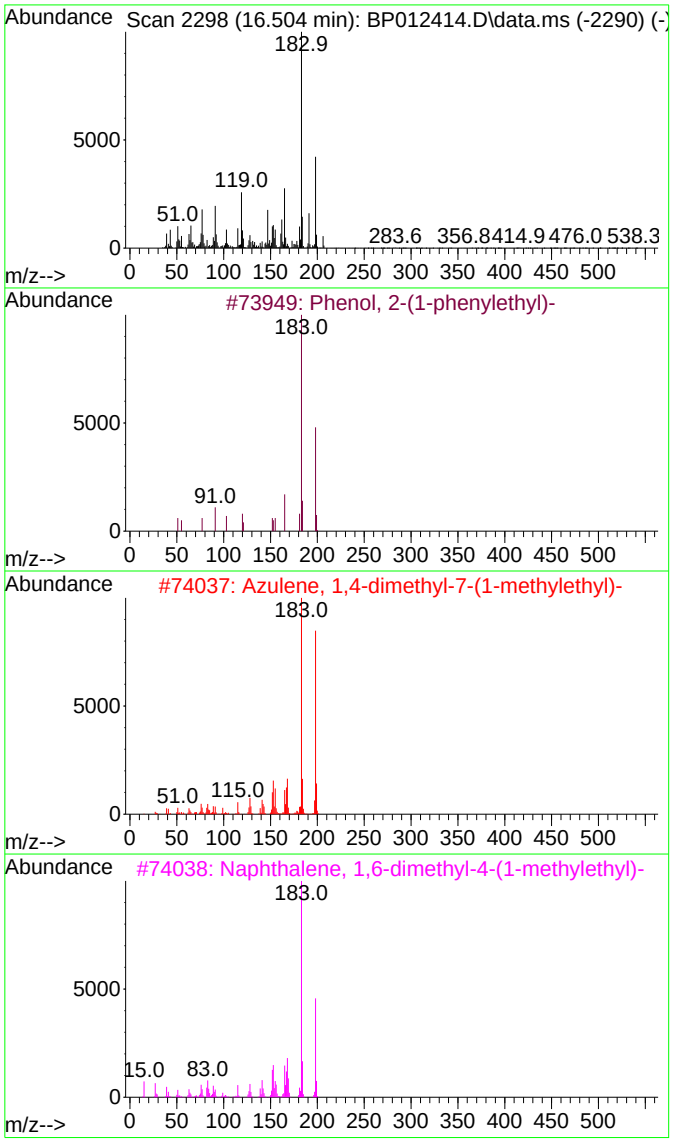
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

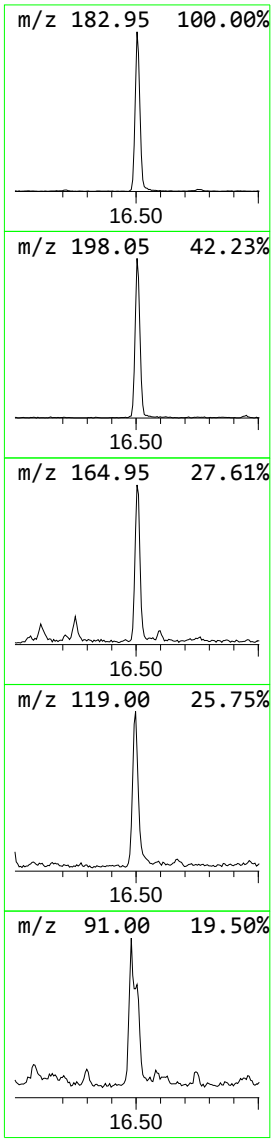
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenol, 2-(1-phenylethyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.504	2.23 ng/ul	678838	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-phenylethyl)-	198	C14H14O	004237-44-9	93
2	Azulene, 1,4-dimethyl-7-(1-methy...	198	C15H18	000489-84-9	70
3	Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	70
4	Naphthalene, 1,6-dimethyl-4-(1-m...	198	C15H18	000483-78-3	64
5	Phenol, 4-(1-phenylethyl)-	198	C14H14O	001988-89-2	62





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Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

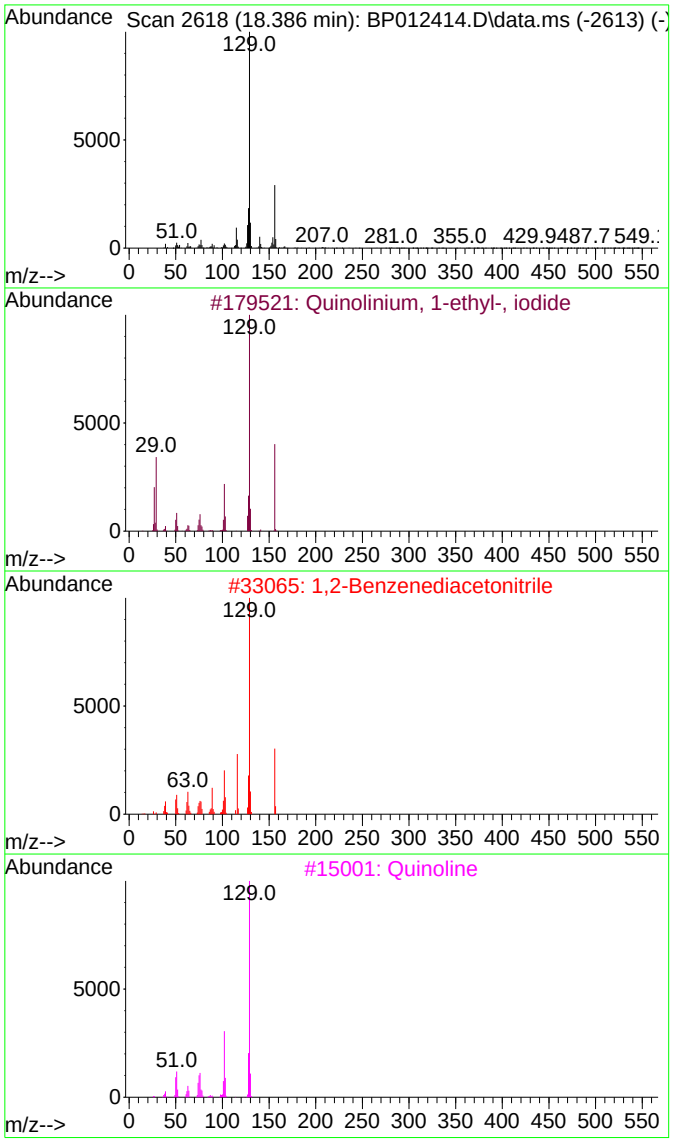
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION

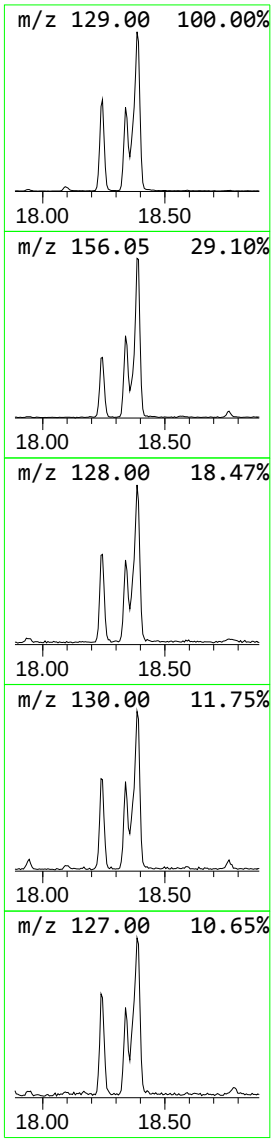
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Quinolinium, 1-ethyl-, iodide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.386	2.59 ng/ul	788358	Phenanthrene-d10	17.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinolinium, 1-ethyl-, iodide	285	C11H12IN	000634-35-5	64
2	1,2-Benzenediacetonitrile	156	C10H8N2	000613-73-0	59
3	Quinoline	129	C9H7N	000091-22-5	52
4	Isoquinoline	129	C9H7N	000119-65-3	52
5	Benzeneacetonitrile, .alpha.-met...	129	C9H7N	000495-10-3	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012414.D
Acq On : 01 Nov 2022 07:33
Operator : CG/JU
Sample : N5216-13
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHAC2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.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
Triethylamine	3.028	23.1	ng/ul	2765190	1	7.792	2390860	20.0
Benzene, 1,3-di...	5.422	22.3	ng/ul	2668350	1	7.792	2390860	20.0
3-Octanone	6.457	2.1	ng/ul	253012	1	7.792	2390860	20.0
Benzene, 1-chlo...	6.763	2.0	ng/ul	241932	1	7.792	2390860	20.0
unknown-01	9.022	2.2	ng/ul	265001	1	7.792	2390860	20.0
Benzene, 1-ethe...	11.286	2.0	ng/ul	357800	2	10.592	3536580	20.0
Phenol, p-tert-...	11.945	13.9	ng/ul	2449990	2	10.592	3536580	20.0
Tripropyl phosp...	12.980	4.0	ng/ul	1044400	3	14.445	5226390	20.0
Phenol, 2-(1-ph...	16.504	2.2	ng/ul	678838	4	17.204	6076360	20.0
Quinolinium, 1-...	18.386	2.6	ng/ul	788358	4	17.204	6076360	20.0