LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

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-BP112624.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP023365.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.105	16	20	25	rBV	19184	24056	0.74%	0.079%
2	3.534	90	93	109	rBV	35051	81656	2.50%	0.267%
3	3.993	165	171	180	rBV5	8109	20212	0.62%	0.066%
4	4.246	210	214	220	rVB2	12257	18001	0.55%	0.059%
5	4.752	297	300	307	rVV3	7333	14028	0.43%	0.046%
6	5.140	359	366	372	rBV	24261	41308	1.26%	0.135%
7	5.646	445	452	455	rBV2	18891	32375	0.99%	0.106%
8	5.687	455	459	473	rVB	34815	75181	2.30%	0.246%
9	6.634	615	620	625	rBV	37280	67512	2.06%	0.221%
10	6.763	636	642	658	rBV2	76704	201826	6.17%	0.660%
11	6.893	658	664	676	rVB	285828	451763	13.82%	1.478%
12	7.093	689	698	707	rBV	318903	577901	17.67%	1.890%
13	7.158	707	709	717	rVB4	11674	21348	0.65%	0.070%
14	7.358	738	743	753	rVB2	7865	16638	0.51%	0.054%
15	7.540	764	774	782	rBV	350306	563714	17.24%	1.844%
16	7.605	782	785	794	rVV2	17736	33989	1.04%	0.111%
17	7.822	816	822	834	rBV4	21178	64480	1.97%	0.211%
18	8.093	864	868	875	rBV	12379	20919	0.64%	0.068%
19	8.163	875	880	890	rBV3	47136	114794	3.51%	0.376%
20	8.258	890	896	913	rVV	242161	533992	16.33%	1.747%
21	8.375	913	916	924	rVB6	8282	15090	0.46%	0.049%
22	8.622	949	958	964	rBV2	90039	175964	5.38%	0.576%
23	8.687	964	969	982	rVB	338614	637603	19.50%	2.086%
24	9.399	1084	1090	1105	rBV2	300173	579389	17.72%	1.895%
25	9.557	1110	1117	1124	rVV2	22062	55585	1.70%	0.182%
26	9.675	1126	1137	1149	rVV4	46152	157875	4.83%	0.516%
27	9.763	1149	1152	1153	rVV3	9406	11063	0.34%	0.036%
28	9.793	1153	1157	1167	rVB2	22128	43645	1.33%	0.143%
29	9.940	1174	1182	1198	rBV	399705	837152	25.60%	2.739%
30	10.093	1201	1208	1221	rVV2	52320	134005	4.10%	0.438%
31	10.204	1221	1227	1233	rVV7	14953	32080	0.98%	0.105%
32	10.281	1234	1240	1249	rVB	405042	709020	21.68%	2.319%
33	10.446	1259	1268	1284	rBV	212622	477107	14.59%	1.561%
34	10.593	1288	1293	1301	rVB7	15173	28416	0.87%	0.093%
35	10.828	1327	1333	1339	rBV2	18681	37560	1.15%	0.123%
36	10.887	1339	1343	1348	rVV	22671	40557	1.24%	0.133%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

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-BP112624.MA.M
Title : SVOA CALIBRATION

37	10.969	1352	1357	1365	rVB2	11717	27287	0.83%	0.089%
38	11.069	1368	1374	1378	rBV9	5799	14070	0.43%	0.046%
39	11.128	1378	1384	1391	rVV2	31045	66791	2.04%	0.218%
40	11.210	1394	1398	1405	rVV	16721	33587	1.03%	0.110%
41	11.293	1408	1412	1421	rVB5	8730	21507	0.66%	0.070%
42	11.546	1449	1455	1460	rBV8	6674	12530	0.38%	0.041%
43	11.646	1466	1472	1482	rBV6	16327	41634	1.27%	0.136%
44	12.216	1565	1569	1573	rBV2	10879	16808	0.51%	0.055%
45	12.369	1590	1595	1601	rBV7	6919	14133	0.43%	0.046%
46	12.434	1601	1606	1611	rBV3	19196	35624	1.09%	0.117%
47	12.493	1611	1616	1632	rVB3	21258	64471	1.97%	0.211%
48	12.622	1632	1638	1639	rBV	27657	43935	1.34%	0.144%
49	12.745	1653	1659	1668	rVB	40631	69215	2.12%	0.226%
50	12.993	1697	1701	1705	rVV6	6681	13084	0.40%	0.043%
51	13.040	1705	1709	1715	rVB2	16570	30778	0.94%	0.101%
52	13.110	1715	1721	1734	rBV2	57944	134231	4.11%	0.439%
53	13.263	1740	1747	1755	rBV7	9899	25986	0.79%	0.085%
54	13.340	1755	1760	1762	rVV6	8225	16030	0.49%	0.052%
55	13.381	1765	1767	1771	rVV5	8956	12067	0.37%	0.039%
56	13.428	1773	1775	1784	rVB9	7317	11380	0.35%	0.037%
57	13.569	1791	1799	1805	rBV	995098	1404444	42.95%	4.594%
58	13.675	1813	1817	1821	rVB6	17804	23138	0.71%	0.076%
59	13.745	1824	1829	1837	rVB6	7305	14906	0.46%	0.049%
60	13.845	1837	1846	1855	rBV2	795140	1365980	41.77%	4.468%
61	13.934	1856	1861	1870	rVV	33529	62498	1.91%	0.204%
62	14.034	1873	1878	1889	rVV	207904	301026	9.21%	0.985%
63	14.151	1889	1898	1908	rVB	697774	1057149	32.33%	3.458%
64	14.463	1945	1951	1964	rBV4	36694	129541	3.96%	0.424%
65	14.681	1981	1988	1997	rVV	111919	231645	7.08%	0.758%
66	14.751	1997	2000	2012	rVB6	24129	52803	1.61%	0.173%
67	14.863	2012	2019	2026	rVB4	14286	30655	0.94%	0.100%
68	14.928	2026	2030	2034	rVB	10951	14256	0.44%	0.047%
69	14.998	2035	2042	2050	rVB	234425	304918	9.33%	0.997%
70	15.163	2063	2070	2079	rBV	1274755	1942660	59.41%	6.355%
71	15.275	2086	2089	2093	rBV5	10277	19229	0.59%	0.063%
72	15.328	2093	2098	2110	rVV	440085	716144	21.90%	2.343%
73	15.557	2132	2137	2144	rBV	20743	35314	1.08%	0.116%
74	15.798	2174	2178	2184	rVB8	8138	13742	0.42%	0.045%
75	15.969	2199	2207	2214	rBV	11682	31615	0.97%	0.103%
76	16.039	2215	2219	2222	rBV3	18598	25732	0.79%	0.084%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

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-BP112624.MA.M
Title : SVOA CALIBRATION

77	16.075	2222	2225	2229	rBV3	12953	19047	0.58%	0.062%
78	16.551	2301	2306	2311	rVB6	9478	14626	0.45%	0.048%
79	16.969	2370	2377	2387	rBV	1017356	1419582	43.41%	4.644%
80	17.063	2387	2393	2408	rVV	1503160	2323239	71.05%	7.600%
81	17.269	2423	2428	2434	rVB	17404	26116	0.80%	0.085%
82	17.463	2456	2461	2467	rVB2	21705	34912	1.07%	0.114%
83	17.569	2471	2479	2484	rBV6	12856	28286	0.87%	0.093%
84	17.657	2487	2494	2499	rBV	384079	506045	15.48%	1.655%
85	17.839	2519	2525	2531	rVB5	26739	56245	1.72%	0.184%
86	17.898	2531	2535	2541	rBV8	16047	23173	0.71%	0.076%
87	17.963	2543	2546	2547	rBV	11352	10091	0.31%	0.033%
88	17.998	2549	2552	2559	rVB3	26317	42821	1.31%	0.140%
89	18.110	2566	2571	2575	rBV	106609	141218	4.32%	0.462%
90	18.186	2582	2584	2590	rVB6	12062	14427	0.44%	0.047%
91	18.380	2613	2617	2622	rBV2	84554	136105	4.16%	0.445%
92	18.486	2632	2635	2643	rVB6	20355	34271	1.05%	0.112%
93	18.710	2668	2673	2680	rBV2	32788	61179	1.87%	0.200%
94	18.992	2717	2721	2727	rVB3	21070	31625	0.97%	0.103%
95	19.080	2731	2736	2740	rBV	18972	30314	0.93%	0.099%
96	19.239	2760	2763	2769	rBV7	13214	19341	0.59%	0.063%
97	19.451	2792	2799	2817	rBV	1930473	2994236	91.57%	9.795%
98	21.392	3123	3129	3143	rBV	1111097	1851586	56.63%	6.057%
99	24.351	3624	3632	3651	rVB	1289368	3269887	100.00%	10.697%
100	24.563	3660	3668	3689	rVB	637042	2050956	62.72%	6.709%

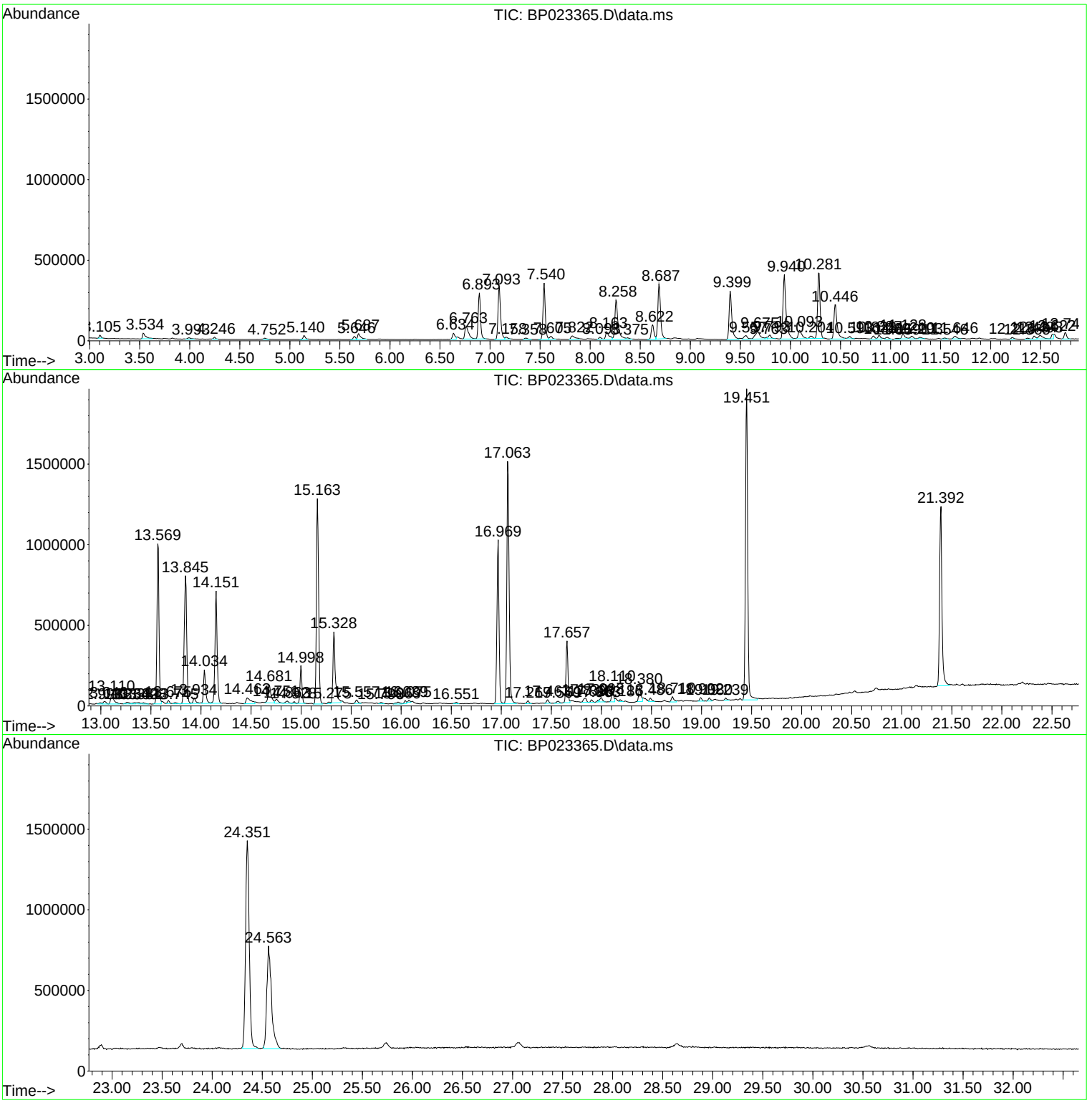
Sum of corrected areas: 30569675

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.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\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

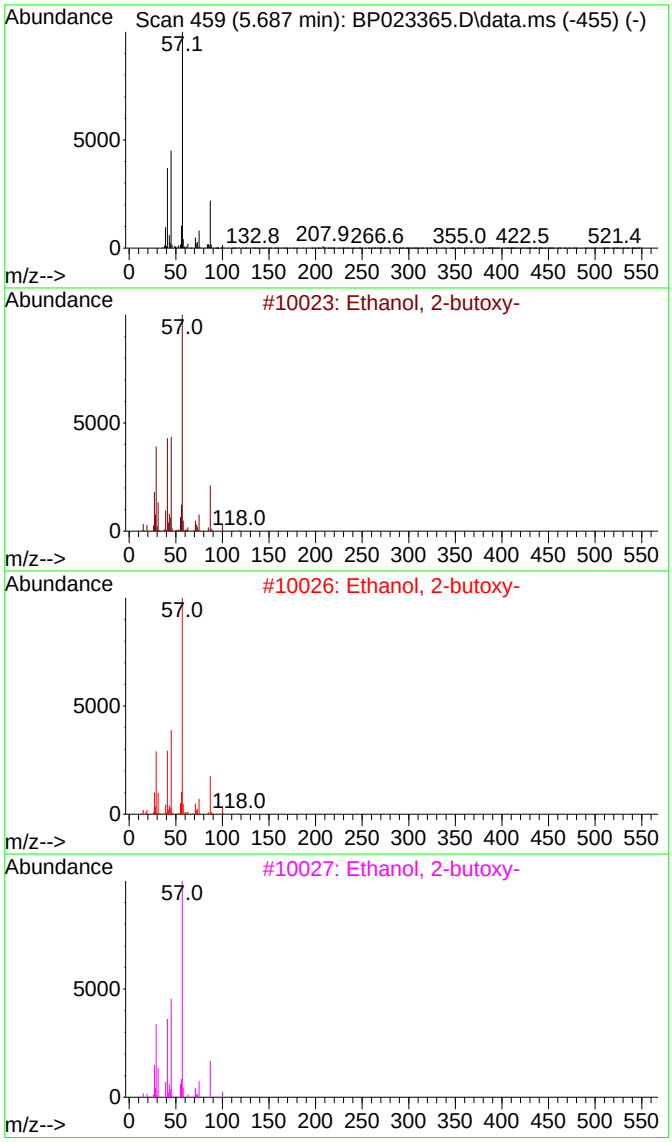
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

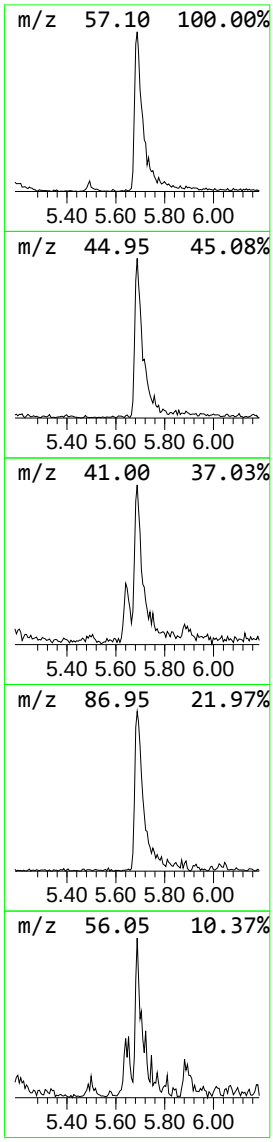
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.687	2.67 ng/ul	75181	1,4-Dichlorobenzene-d4	7.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	90
2	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	56
4	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	53
5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

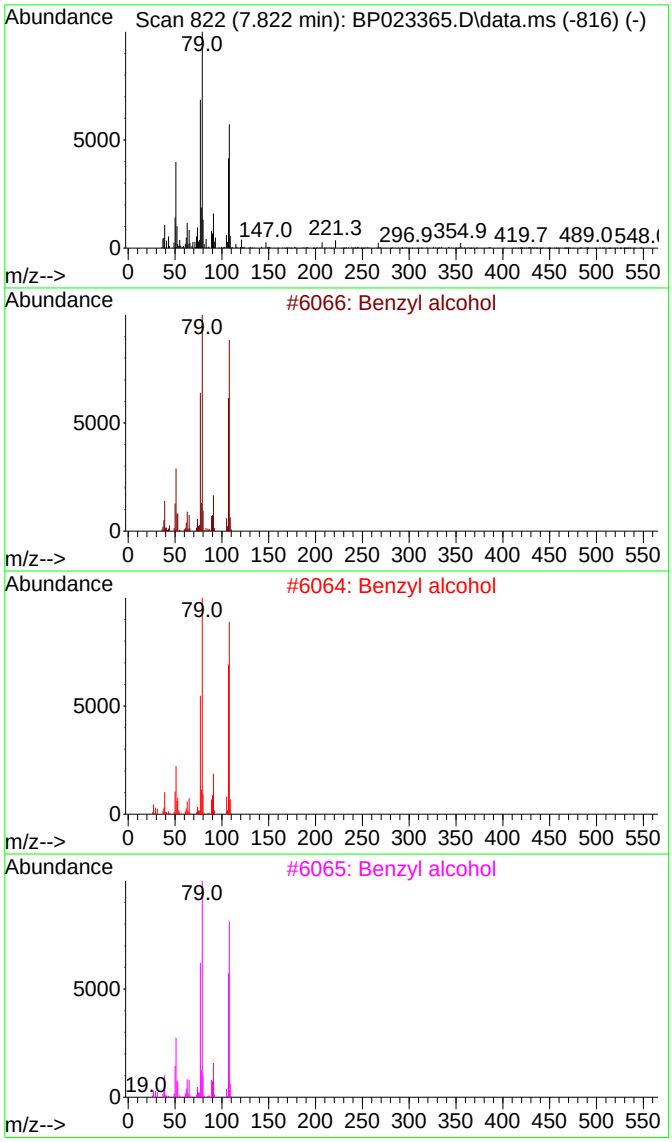
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

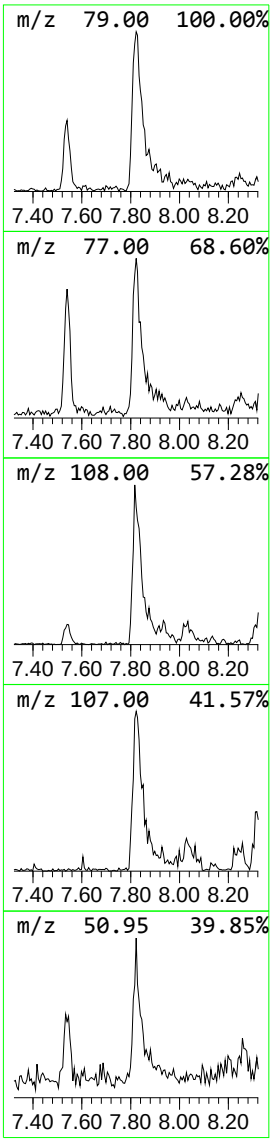
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzyl alcohol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.822	2.29 ng/ul	64480	1,4-Dichlorobenzene-d4	7.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl alcohol	108	C7H8O	000100-51-6	97
2	Benzyl alcohol	108	C7H8O	000100-51-6	96
3	Benzyl alcohol	108	C7H8O	000100-51-6	95
4	Benzyl alcohol	108	C7H8O	000100-51-6	94
5	N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

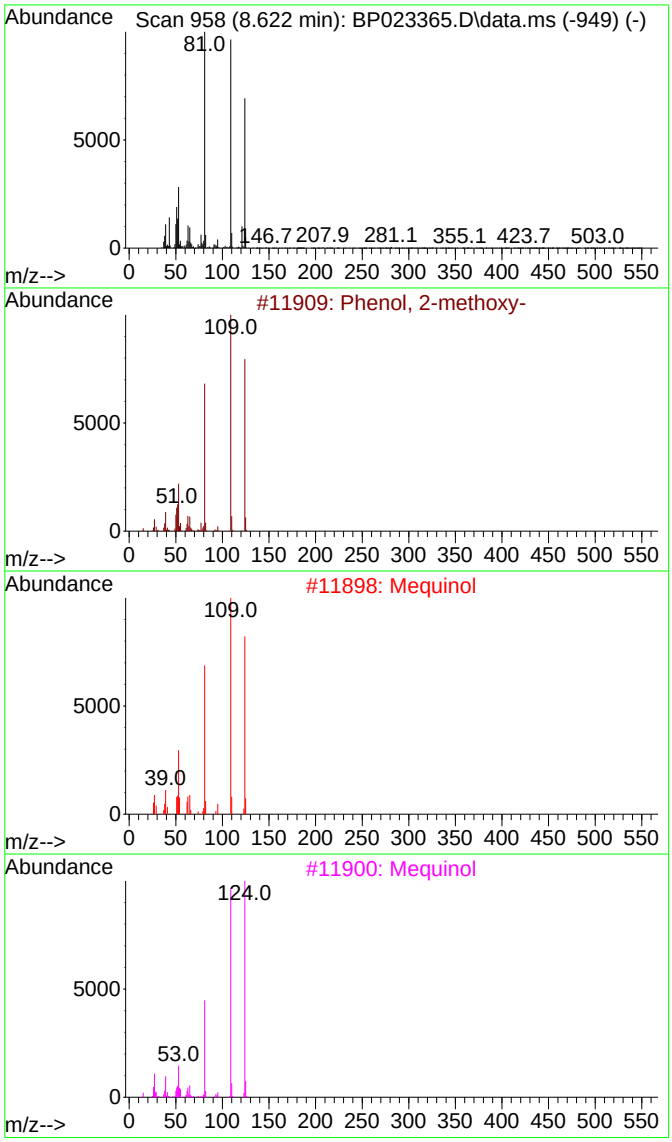
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

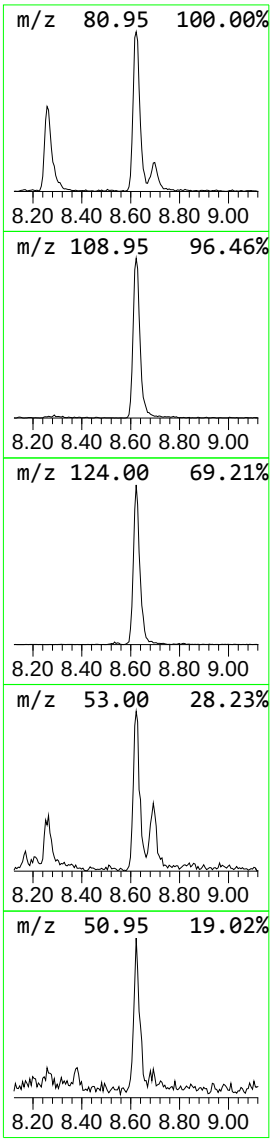
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenol, 2-methoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.622	6.24 ng/ul	175964	1,4-Dichlorobenzene-d4	7.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	95
2	Mequinol	124	C7H8O2	000150-76-5	91
3	Mequinol	124	C7H8O2	000150-76-5	91
4	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90
5	Mequinol	124	C7H8O2	000150-76-5	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

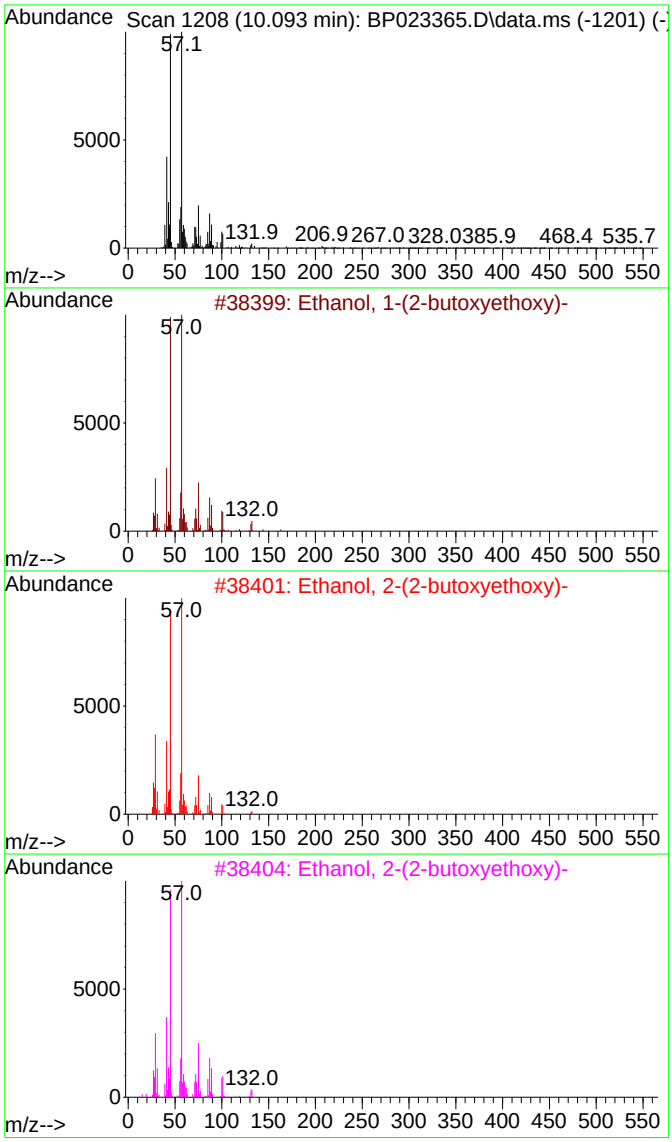
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

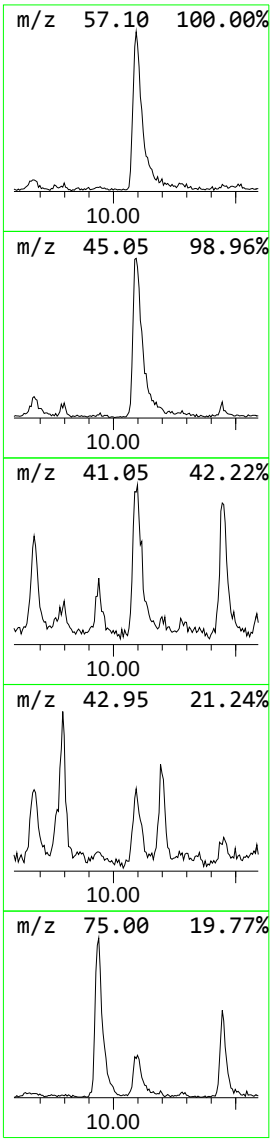
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.093	3.78 ng/ul	134005	Naphthalene-d8	10.281

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

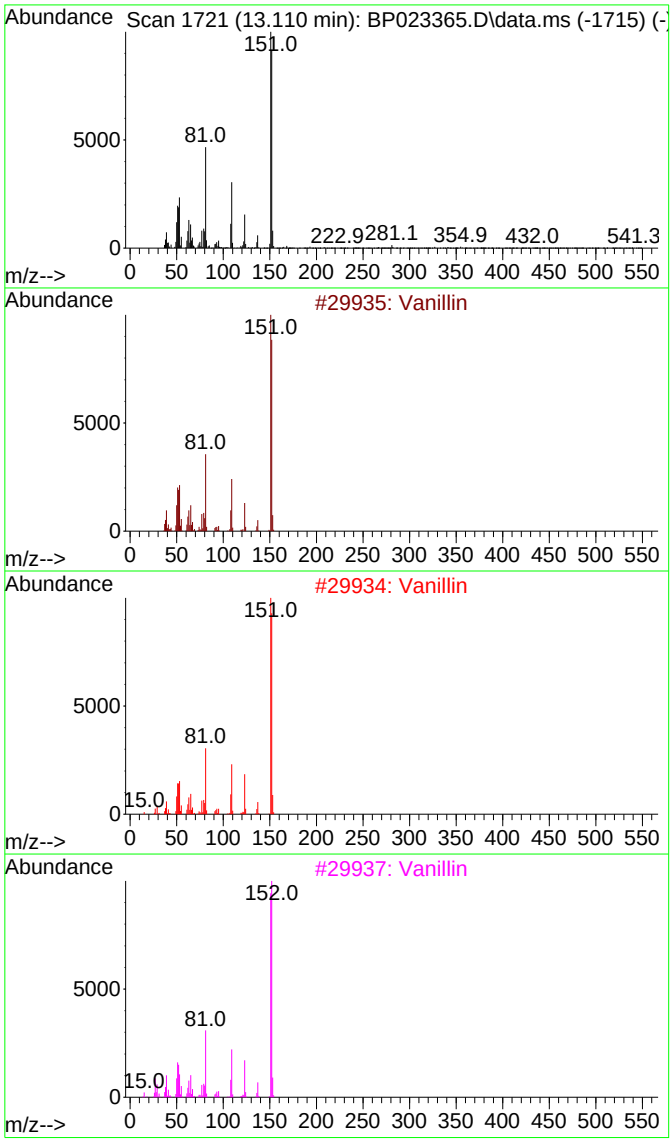
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

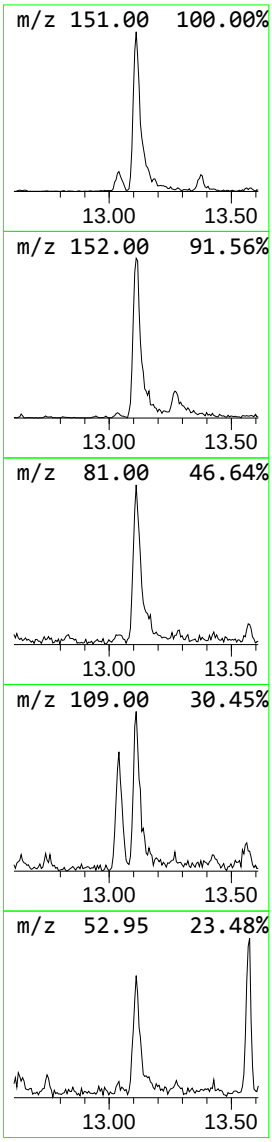
****

Peak Number 8 Vanillin Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.110	2.54 ng/ul	134231	Acenaphthene-d10	14.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	98
2	Vanillin	152	C8H8O3	000121-33-5	97
3	Vanillin	152	C8H8O3	000121-33-5	96
4	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96
5	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

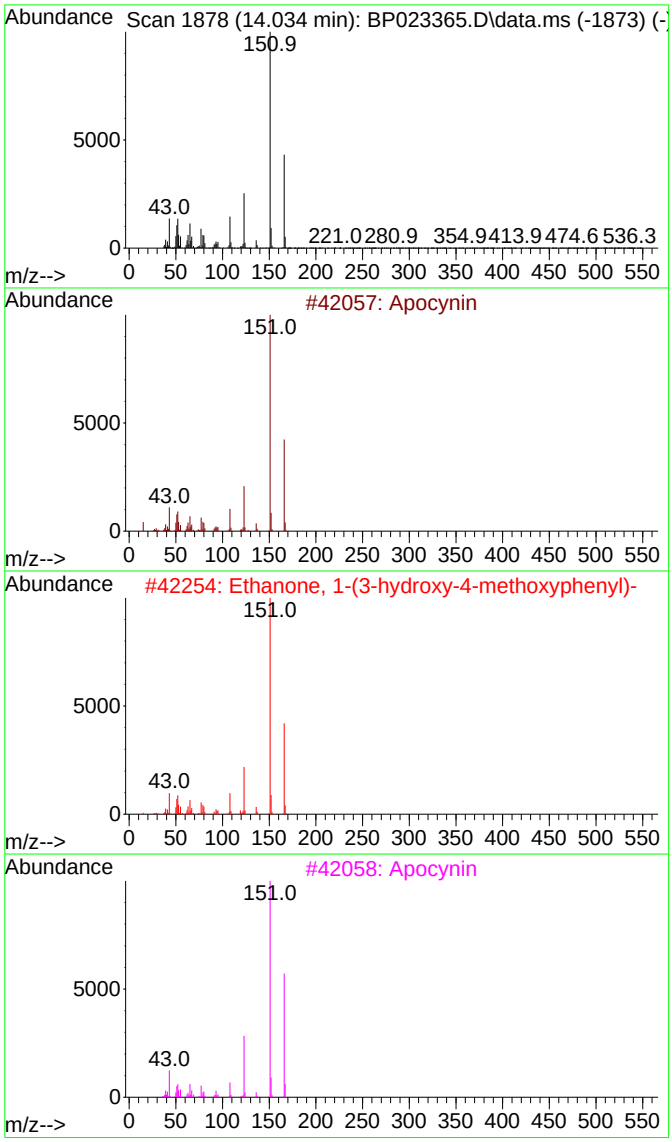
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

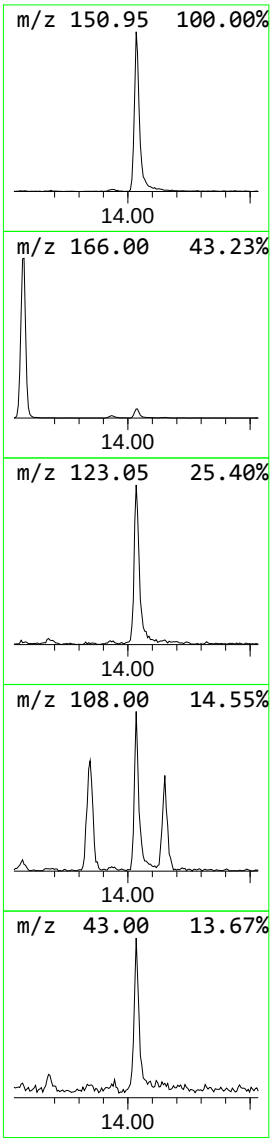
****

Peak Number 9 Apocynin Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.034	5.70 ng/ul	301026	Acenaphthene-d10	14.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Apocynin	166	C9H10O3	000498-02-2	96
2	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	95
3	Apocynin	166	C9H10O3	000498-02-2	94
4	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	93
5	Apocynin	166	C9H10O3	000498-02-2	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

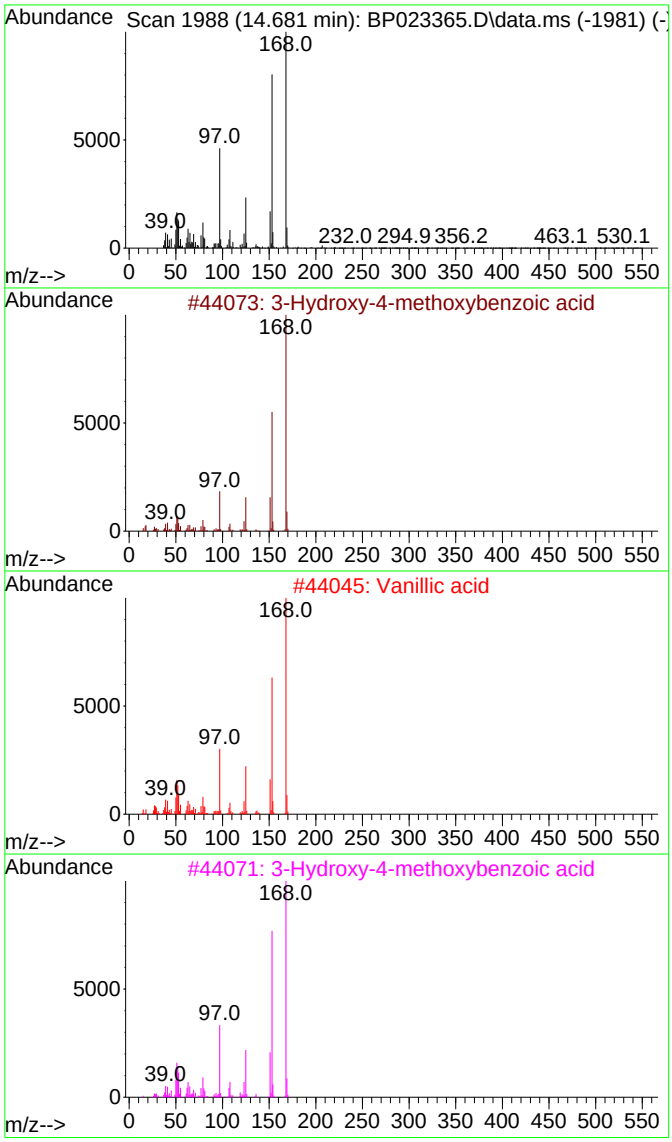
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

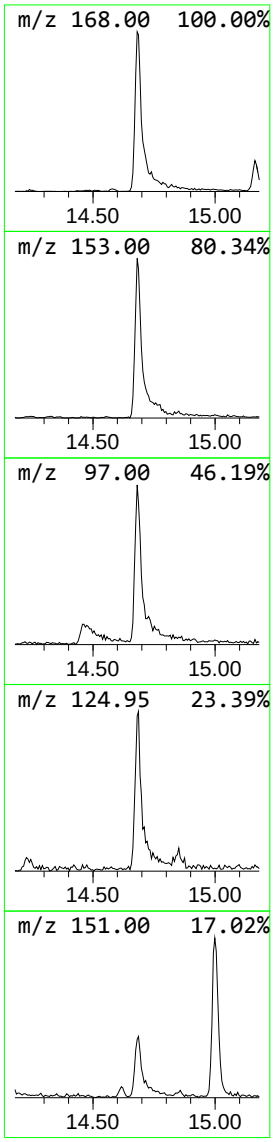
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 3-Hydroxy-4-methoxybenzoic ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.681	4.38 ng/ul	231645	Acenaphthene-d10	14.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hydroxy-4-methoxybenzoic acid	168	C8H8O4	000645-08-9	97
2	Vanillic acid	168	C8H8O4	000121-34-6	96
3	3-Hydroxy-4-methoxybenzoic acid	168	C8H8O4	000645-08-9	96
4	3-Hydroxy-4-methoxybenzoic acid	168	C8H8O4	000645-08-9	96
5	Vanillic acid	168	C8H8O4	000121-34-6	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

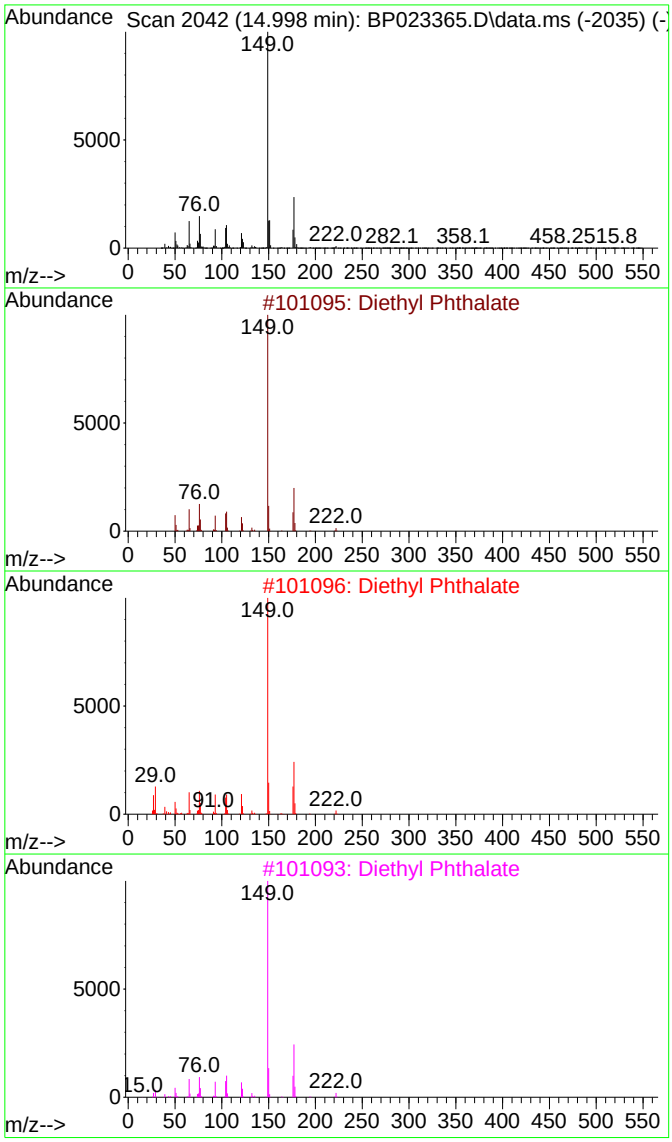
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

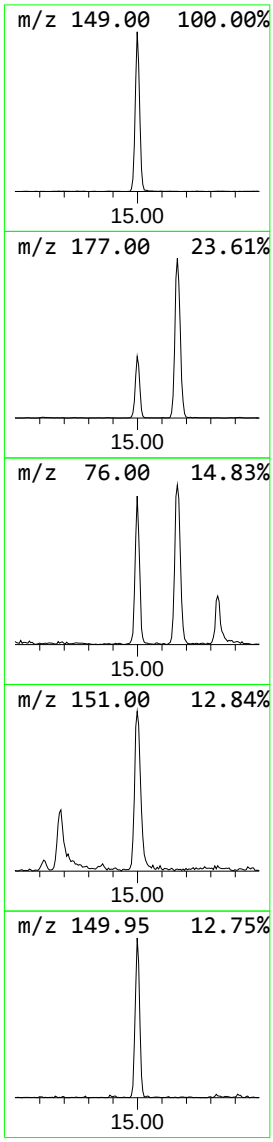
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Diethyl Phthalate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.998	5.77 ng/ul	304918	Acenaphthene-d10	14.151

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyl Phthalate	222	C12H14O4	000084-66-2	96
2	Diethyl Phthalate	222	C12H14O4	000084-66-2	95
3	Diethyl Phthalate	222	C12H14O4	000084-66-2	91
4	Diethyl Phthalate	222	C12H14O4	000084-66-2	87
5	Phthalic acid, allyl ethyl ester	234	C13H14O4	033672-94-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

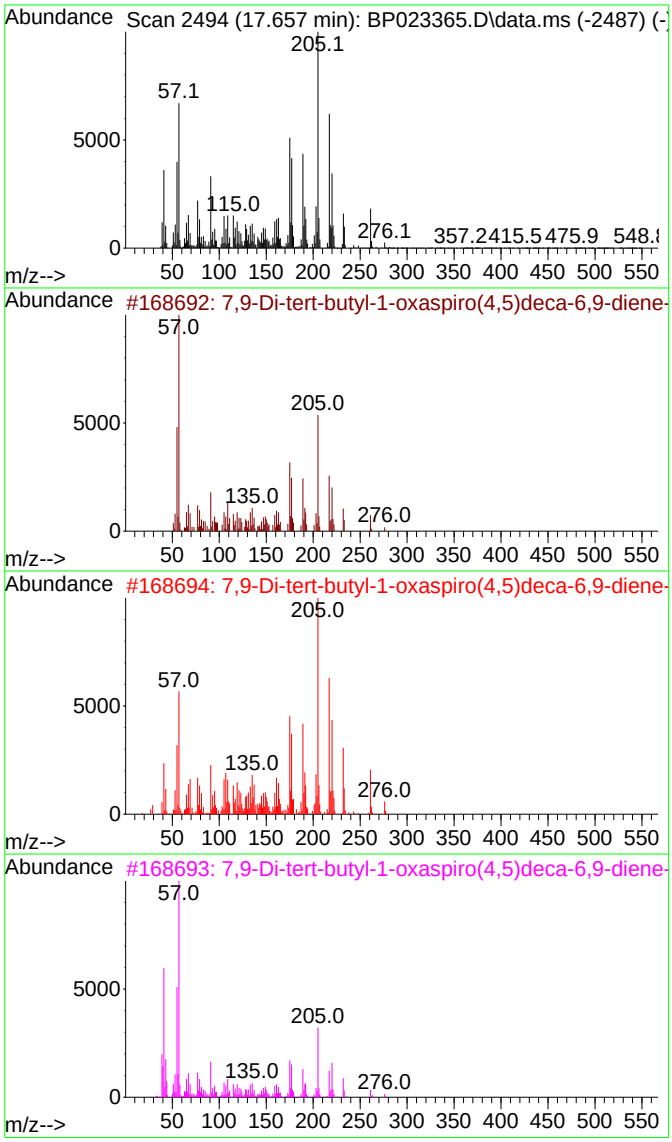
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

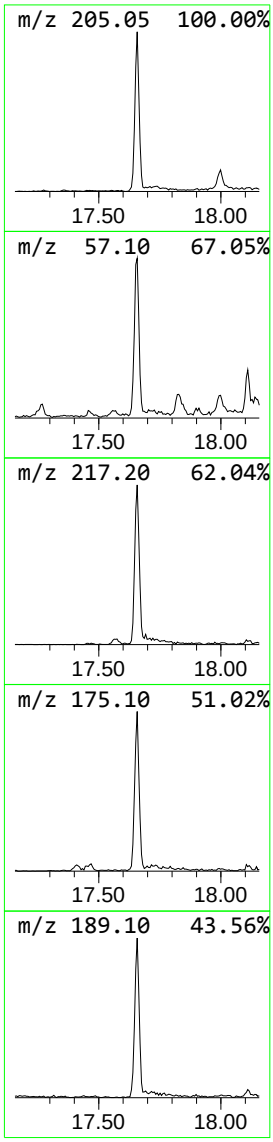
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.657	7.13 ng/ul	506045	Phenanthrene-d10	16.969

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	94
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	91
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	55
4	(1R,3aR,5aR,9aS)-1,4,4,7-Tetrame...	220	C15H24O	104188-25-2	49
5	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	45





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023365.D
Acq On : 11 Dec 2024 04:02
Operator : RC/JU
Sample : P5111-12
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG461

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.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
Ethanol, 2-butoxy-	5.687	2.7	ng/ul	75181	1	7.540	563714	20.0
Benzyl alcohol	7.822	2.3	ng/ul	64480	1	7.540	563714	20.0
Phenol, 2-methoxy-	8.622	6.2	ng/ul	175964	1	7.540	563714	20.0
Ethanol, 1-(2-b...	10.093	3.8	ng/ul	134005	2	10.281	709020	20.0
Vanillin	13.110	2.5	ng/ul	134231	3	14.151	1057150	20.0
Apocynin	14.034	5.7	ng/ul	301026	3	14.151	1057150	20.0
3-Hydroxy-4-met...	14.681	4.4	ng/ul	231645	3	14.151	1057150	20.0
Diethyl Phthalate	14.998	5.8	ng/ul	304918	3	14.151	1057150	20.0
7,9-Di-tert-but...	17.657	7.1	ng/ul	506045	4	16.969	1419580	20.0