

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022005.D
 Acq On : 24 Sep 2024 21:45
 Operator : RC/JU
 Sample : P4092-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B37

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-BP092424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022005.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.193	31	35	38	rBV	46572	58393	1.26%	0.176%
2	3.222	38	40	44	rVV	27289	31579	0.68%	0.095%
3	3.258	44	46	51	rVB3	6584	7276	0.16%	0.022%
4	3.505	84	88	94	rBV7	2349	5045	0.11%	0.015%
5	3.640	107	111	126	rBV	126386	193546	4.17%	0.583%
6	3.734	126	127	136	rVB9	3159	5190	0.11%	0.016%
7	3.964	161	166	172	rVB2	7583	12307	0.26%	0.037%
8	4.487	249	255	265	rVB	62901	100081	2.15%	0.301%
9	4.864	313	319	325	rBV6	3444	7381	0.16%	0.022%
10	5.393	404	409	419	rBV	9449	18983	0.41%	0.057%
11	5.758	465	471	477	rVB	8417	13006	0.28%	0.039%
12	6.705	625	632	640	rBV	140280	217879	4.69%	0.656%
13	6.893	657	664	672	rBV	129123	209744	4.52%	0.631%
14	7.046	684	690	699	rBV	354022	560273	12.06%	1.687%
15	7.252	716	725	734	rBV	437841	692086	14.90%	2.083%
16	7.363	738	744	750	rVB4	6539	14323	0.31%	0.043%
17	7.710	796	803	813	rBV	333376	549898	11.84%	1.655%
18	7.828	819	823	830	rVB2	5135	8789	0.19%	0.026%
19	8.252	890	895	902	rBV9	2114	4669	0.10%	0.014%
20	8.340	904	910	915	rBV8	3464	6934	0.15%	0.021%
21	8.405	915	921	931	rBV	370914	587369	12.65%	1.768%
22	8.540	938	944	948	rVB3	4046	7456	0.16%	0.022%
23	8.846	989	996	1007	rVB	467064	745400	16.05%	2.244%
24	9.422	1089	1094	1101	rVB8	2575	5558	0.12%	0.017%
25	9.563	1110	1118	1127	rBV	378202	664661	14.31%	2.001%
26	10.093	1200	1208	1221	rBV	640469	1069319	23.02%	3.219%
27	10.469	1264	1272	1281	rBV	410929	747092	16.09%	2.249%
28	10.604	1288	1295	1313	rBV	479675	819380	17.64%	2.467%
29	12.057	1535	1542	1547	rBV2	9545	15838	0.34%	0.048%
30	13.722	1817	1825	1835	rBV	1257943	1732855	37.31%	5.217%
31	14.004	1866	1873	1884	rBV2	1342584	1866625	40.19%	5.619%
32	14.316	1919	1926	1935	rBV	724402	1139435	24.53%	3.430%
33	14.522	1955	1961	1979	rBV	94127	212830	4.58%	0.641%
34	14.751	1994	2000	2008	rBV	26923	44043	0.95%	0.133%
35	15.316	2089	2096	2106	rBV	1715958	2505875	53.95%	7.544%
36	15.445	2111	2118	2131	rVV	629258	829965	17.87%	2.499%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Title : SVOA CALIBRATION

37	15.569	2135	2139	2146	rVB4	9404	16722	0.36%	0.050%
38	16.686	2326	2329	2335	rVB6	4707	7978	0.17%	0.024%
39	17.116	2395	2402	2413	rBV	1005626	1546051	33.29%	4.654%
40	17.216	2413	2419	2436	rVB	1974597	2897310	62.38%	8.722%
41	18.045	2556	2560	2567	rBV3	23945	34033	0.73%	0.102%
42	19.133	2741	2745	2751	rBV2	31978	47300	1.02%	0.142%
43	19.204	2753	2757	2761	rBV	31218	47920	1.03%	0.144%
44	19.386	2785	2788	2793	rVB6	15128	18628	0.40%	0.056%
45	19.586	2816	2822	2831	rVB	2852571	3910362	84.20%	11.772%
46	21.539	3148	3154	3161	rBV2	1209635	2060018	44.35%	6.201%
47	24.592	3664	3673	3688	rVB	1709012	4644388	100.00%	13.981%
48	24.821	3704	3712	3724	rVB	906312	2276561	49.02%	6.853%

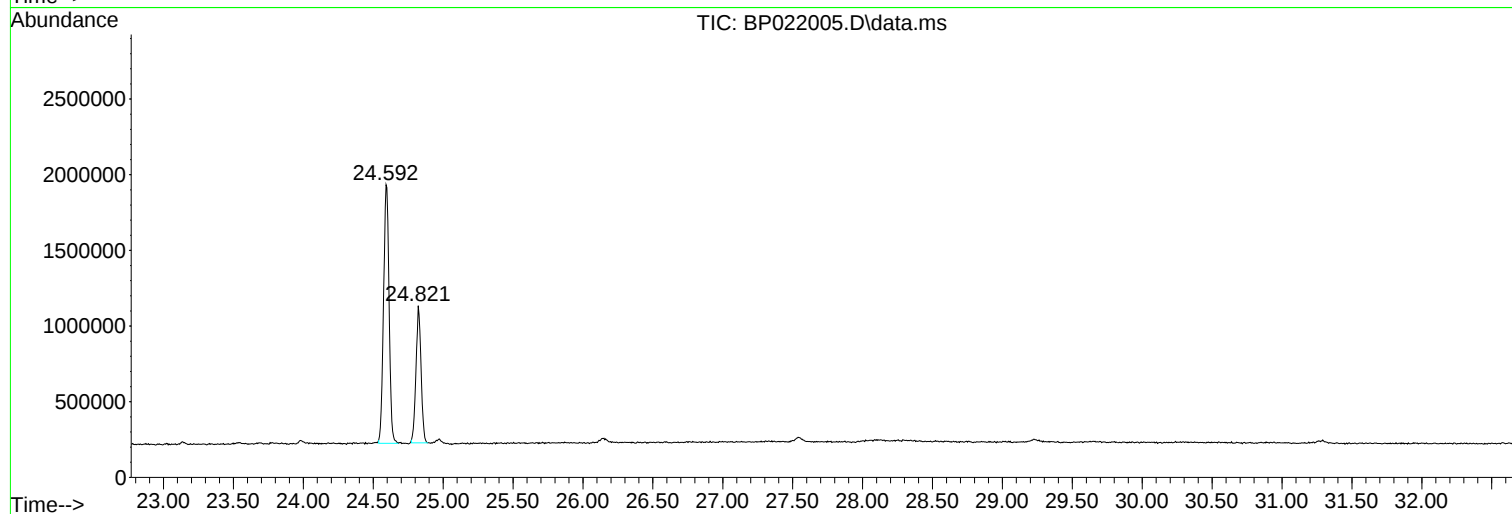
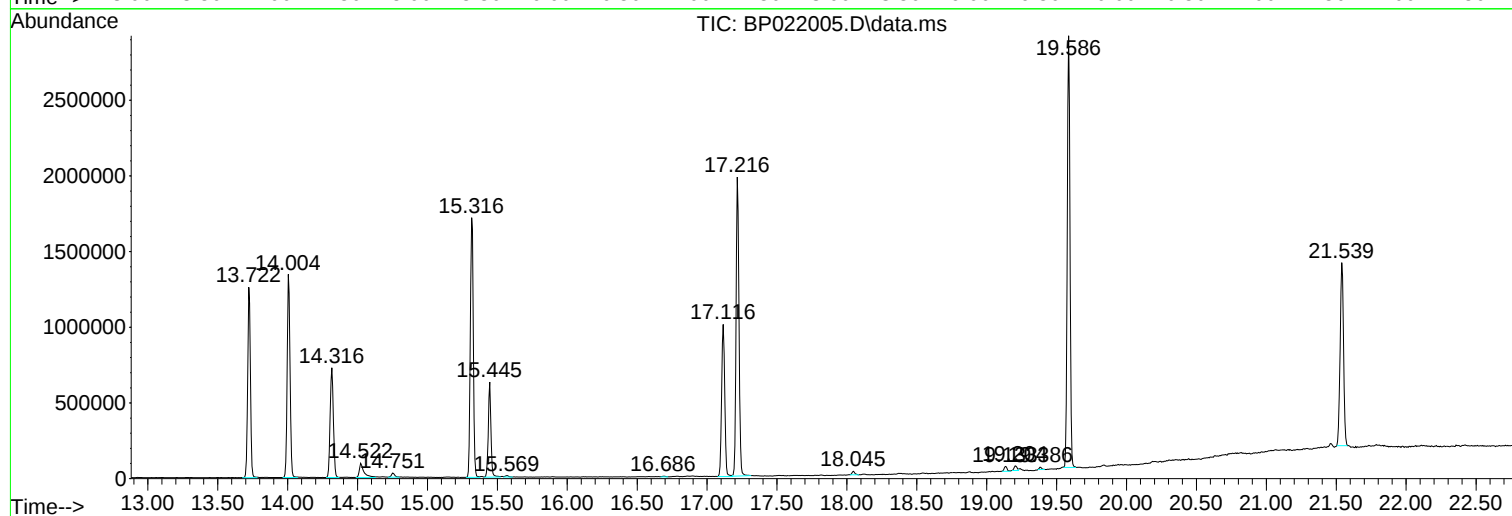
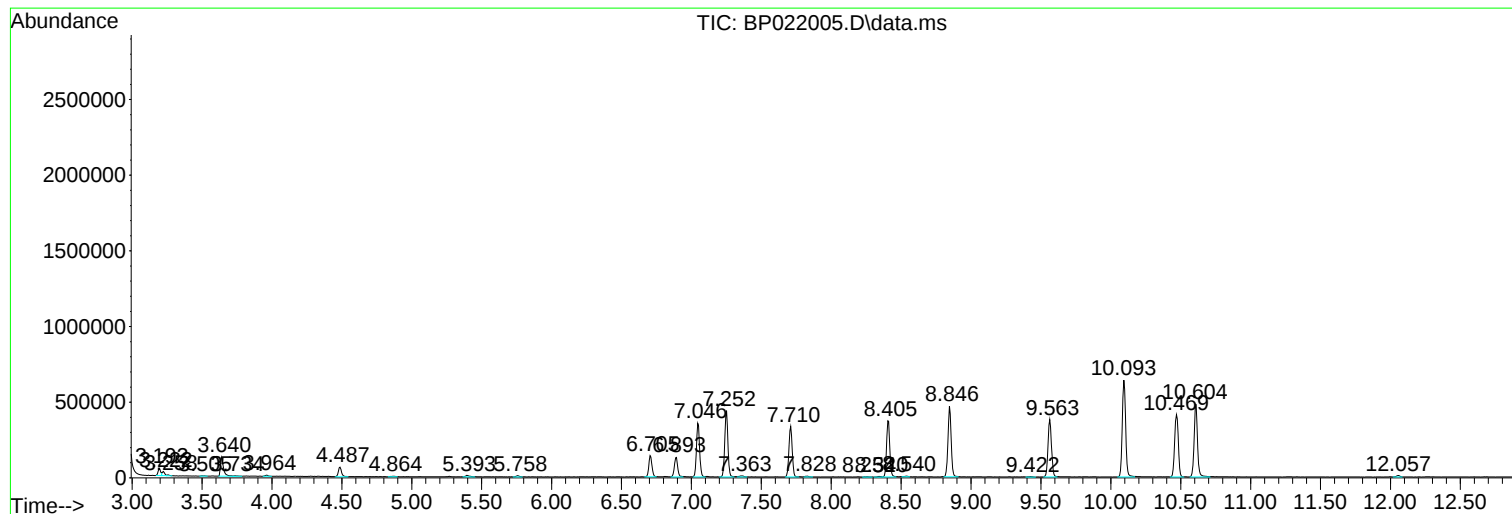
Sum of corrected areas: 33218354

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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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



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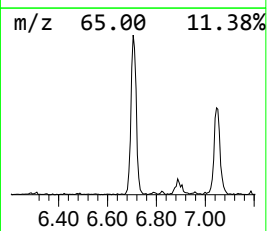
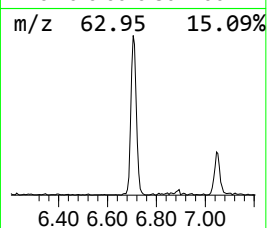
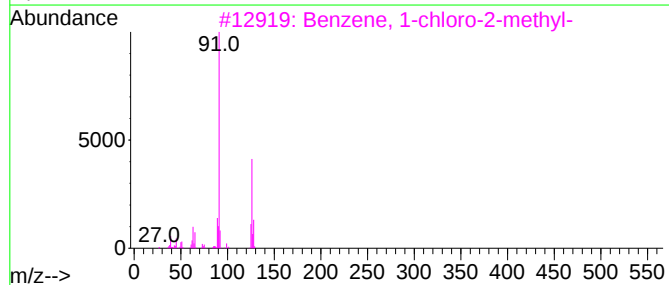
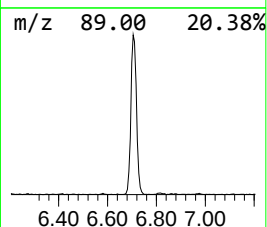
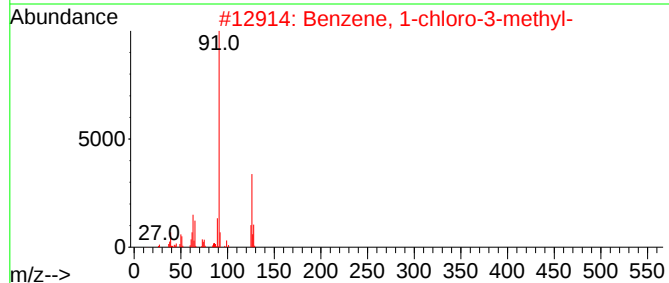
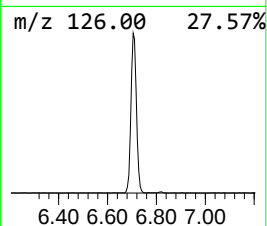
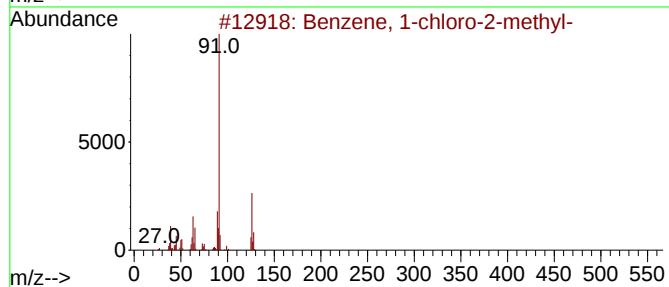
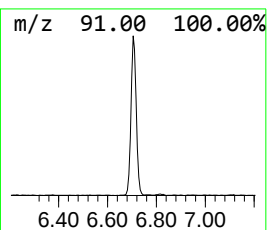
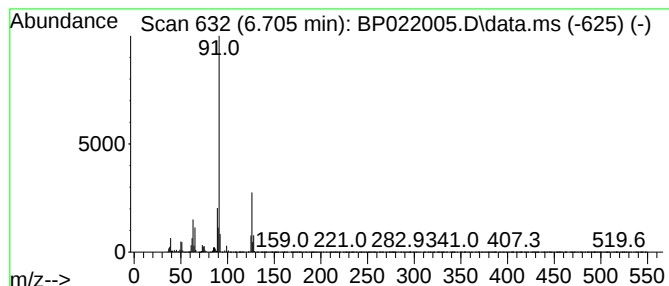
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1-chloro-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.705	7.92 ng/ul	217879	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	96
2			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	95
3			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	90
4			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	90
5			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	90



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-chlo...	6.705	7.9	ng/u1	217879	1	7.710	549898	20.0