

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012991.D
 Acq On : 08 Dec 2022 18:49
 Operator : CG/JU
 Sample : N5894-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA4

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012991.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.358	25	29	38	rVB	140982	212680	0.28%	0.069%
2	3.787	99	102	119	rBV	378159	658512	0.87%	0.213%
3	4.017	136	141	148	rVB2	68903	110599	0.15%	0.036%
4	4.117	154	158	164	rVB	59235	84129	0.11%	0.027%
5	5.269	346	354	372	rBV2	31976338	75788078	100.00%	24.565%
6	7.128	661	670	684	rBV2	545112	1182620	1.56%	0.383%
7	7.299	690	699	711	rBV	2342878	3698569	4.88%	1.199%
8	7.505	727	734	745	rBV	1432360	2463719	3.25%	0.799%
9	7.693	761	766	774	rVB	92980	158292	0.21%	0.051%
10	7.863	787	795	803	rBV	973637	1518838	2.00%	0.492%
11	8.016	815	821	841	rVB	21926178	38774442	51.16%	12.568%
12	8.334	866	875	889	rBV	19849968	35656509	47.05%	11.557%
13	8.681	926	934	943	rBV	1231976	1955952	2.58%	0.634%
14	9.140	1005	1012	1028	rBV	2315407	3818193	5.04%	1.238%
15	9.810	1119	1126	1130	rBV	64430	119043	0.16%	0.039%
16	9.869	1130	1136	1144	rVB	1350383	2177685	2.87%	0.706%
17	10.405	1219	1227	1234	rBV	2070212	3298967	4.35%	1.069%
18	10.652	1262	1269	1282	rBV3	252079	569703	0.75%	0.185%
19	10.793	1285	1293	1304	rBV	3299392	5563651	7.34%	1.803%
20	10.922	1304	1315	1328	rBV	1976264	3843612	5.07%	1.246%
21	12.369	1556	1561	1568	rBV	60374	103201	0.14%	0.033%
22	14.022	1835	1842	1857	rBV	5514664	7811370	10.31%	2.532%
23	14.310	1883	1891	1906	rBV	7256960	10303922	13.60%	3.340%
24	14.616	1935	1943	1953	rBV	5561715	8131199	10.73%	2.636%
25	14.798	1967	1974	1988	rBV	322366	596685	0.79%	0.193%
26	15.028	2007	2013	2020	rVB5	60558	93151	0.12%	0.030%
27	15.610	2104	2112	2124	rBV	9502721	13866462	18.30%	4.495%
28	15.722	2124	2131	2142	rVV	2016540	2685102	3.54%	0.870%
29	15.845	2147	2152	2159	rVB	47568	80784	0.11%	0.026%
30	17.369	2404	2411	2418	rBV	7139785	10042598	13.25%	3.255%
31	17.469	2421	2428	2439	rVV2	10922700	15656887	20.66%	5.075%
32	17.557	2439	2443	2456	rVB5	73372	152810	0.20%	0.050%
33	19.275	2731	2735	2741	rVB2	142263	187899	0.25%	0.061%
34	19.339	2742	2746	2751	rBV	133463	197089	0.26%	0.064%
35	19.698	2801	2807	2822	rVV	14703047	19024021	25.10%	6.166%
36	21.451	3100	3105	3113	rBV	8045396	10630003	14.03%	3.445%

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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

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37	23.780	3494	3501	3516	rBV2	7866834	16962941	22.38%	5.498%
38	23.945	3522	3529	3542	rVB2	5076339	10339760	13.64%	3.351%

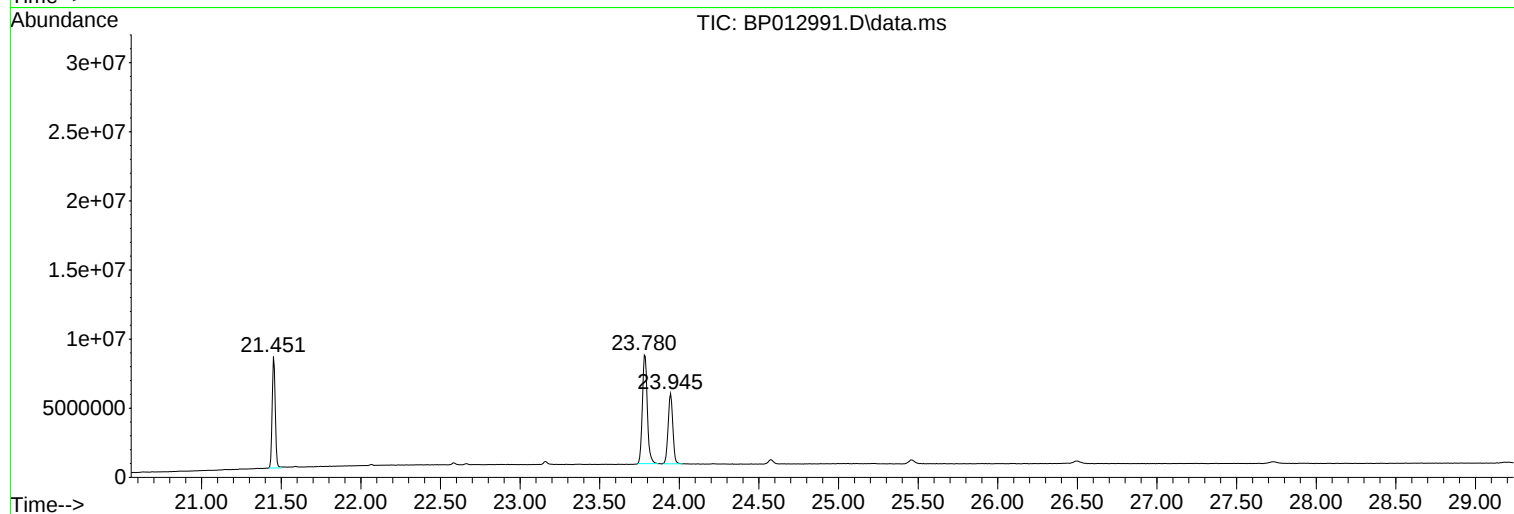
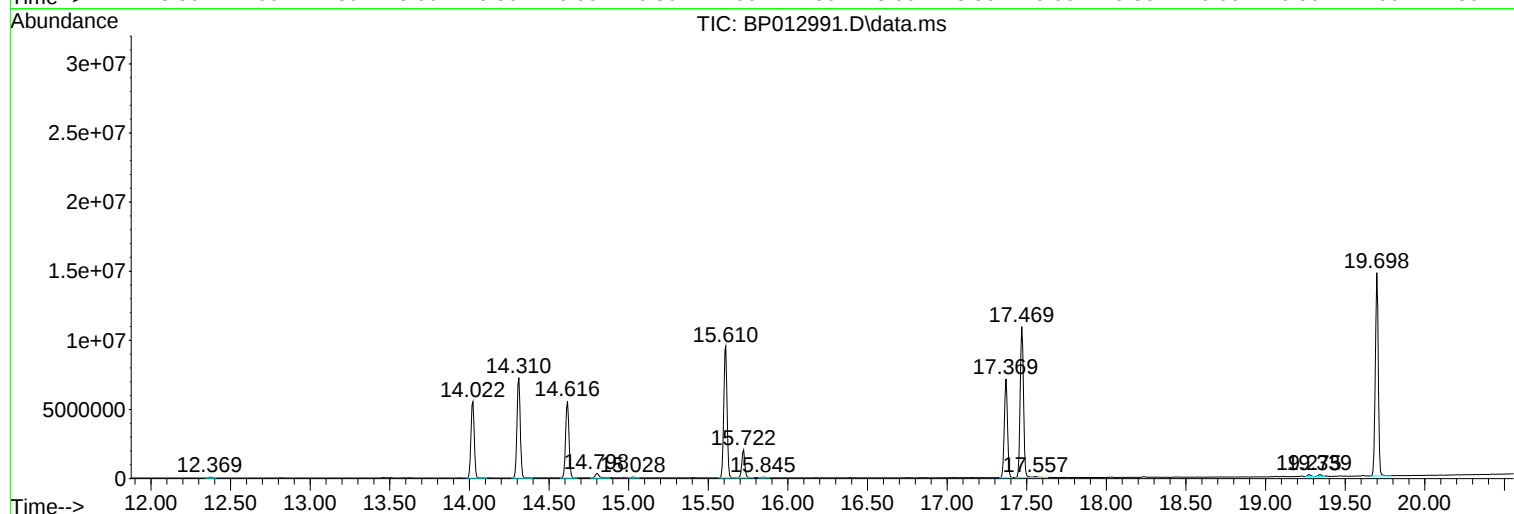
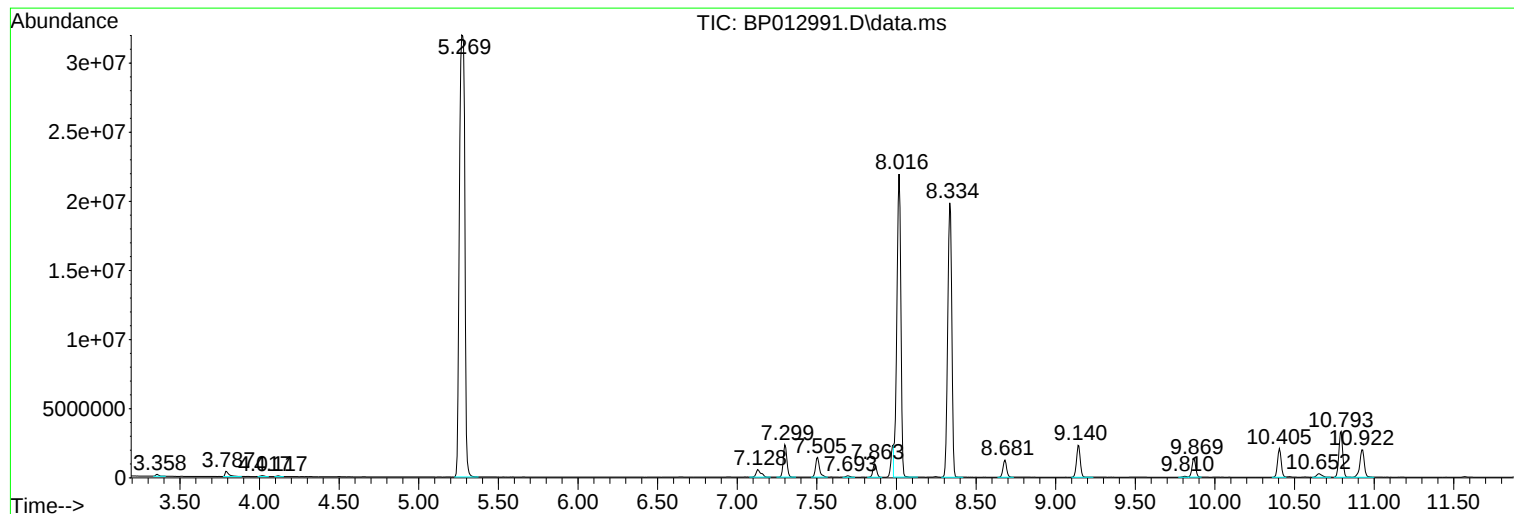
Sum of corrected areas: 308519677

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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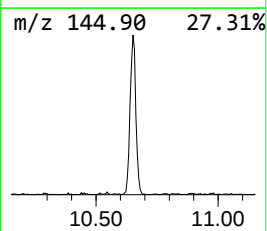
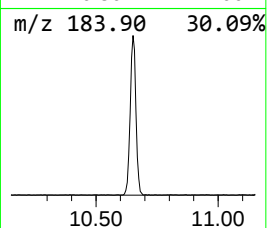
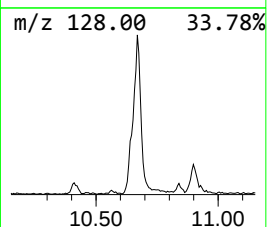
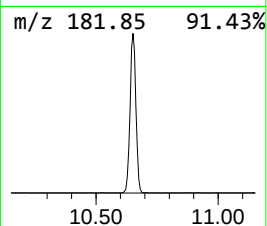
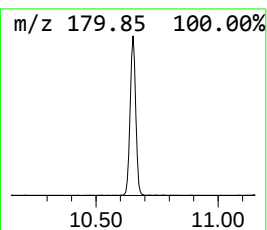
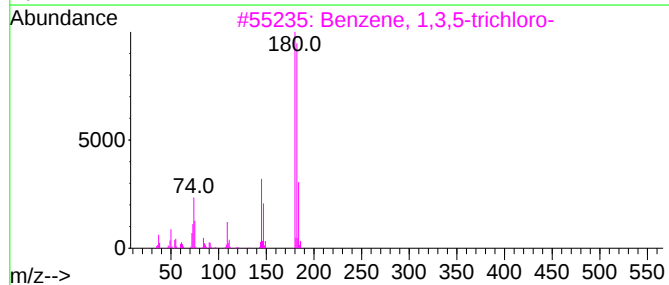
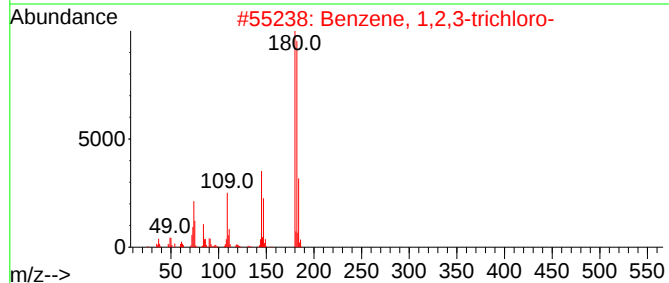
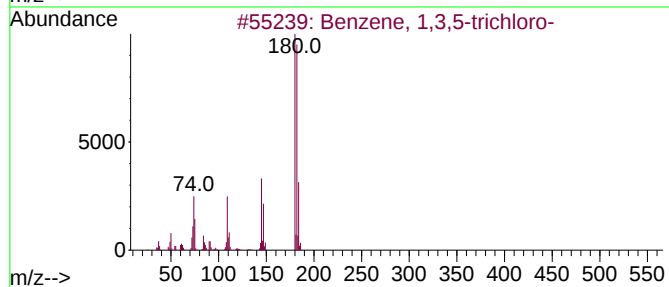
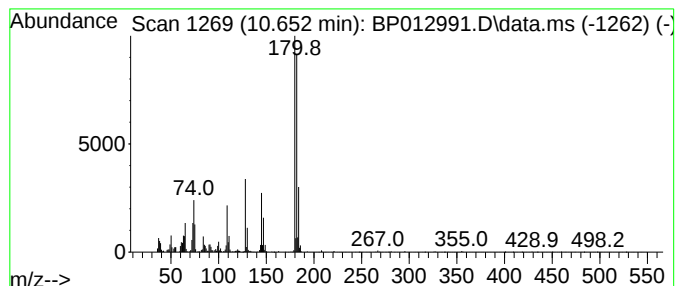
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,3,5-trichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.652	2.05 ng/ul	569703	Naphthalene-d8	10.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
4			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	96



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3,5-...	10.652	2.0	ng/u1	569703	2	10.793	5563650	20.0