

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121523\
 Data File : BP018895.D
 Acq On : 16 Dec 2023 00:01
 Operator : MA/JU
 Sample : 05646-14DL 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AS1DL

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018895.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.140	342	349	363	rVB	5435133	8490717	9.41%	4.814%
2	6.999	660	665	674	rVB2	62662	124378	0.14%	0.071%
3	7.169	688	694	704	rVB	93303	152684	0.17%	0.087%
4	7.363	720	727	735	rBV	91834	163754	0.18%	0.093%
5	7.716	779	787	798	rVB	1158804	1962691	2.17%	1.113%
6	7.899	798	818	823	rBV2	27166929	90247678	100.00%	51.173%
7	8.181	859	866	877	rBV	4902000	7812185	8.66%	4.430%
8	8.540	920	927	933	rBV	88201	143836	0.16%	0.082%
9	8.987	995	1003	1013	rBV	84497	145848	0.16%	0.083%
10	9.552	1092	1099	1102	rBV	182706	291734	0.32%	0.165%
11	9.593	1102	1106	1119	rVV	171330	406259	0.45%	0.230%
12	9.769	1131	1136	1144	rVB	68296	113242	0.13%	0.064%
13	10.004	1168	1176	1182	rBV	51186	91846	0.10%	0.052%
14	10.246	1210	1217	1224	rBV	95028	167111	0.19%	0.095%
15	10.493	1249	1259	1274	rBV	15832334	28316914	31.38%	16.056%
16	10.622	1274	1281	1291	rVB	1197207	1963834	2.18%	1.114%
17	10.769	1299	1306	1316	rVB	79915	155510	0.17%	0.088%
18	11.010	1338	1347	1360	rBV	7519061	12899736	14.29%	7.314%
19	12.363	1570	1577	1585	rBV2	62387	102520	0.11%	0.058%
20	12.616	1612	1620	1622	rBV	262604	400461	0.44%	0.227%
21	12.645	1622	1625	1632	rVV	575182	901483	1.00%	0.511%
22	13.257	1721	1729	1742	rBV	2613534	3929204	4.35%	2.228%
23	13.875	1828	1834	1842	rBV	174086	251623	0.28%	0.143%
24	14.151	1875	1881	1889	rBV2	209409	306070	0.34%	0.174%
25	14.457	1925	1933	1948	rBV	1649511	2463699	2.73%	1.397%
26	14.775	1980	1987	1993	rVV	383571	553301	0.61%	0.314%
27	15.451	2096	2102	2113	rBV	305455	461089	0.51%	0.261%
28	17.204	2394	2400	2412	rBV	2175858	3053447	3.38%	1.731%
29	17.304	2412	2417	2434	rVB	304549	479269	0.53%	0.272%
30	19.539	2792	2797	2804	rVV	491766	657314	0.73%	0.373%
31	20.357	2932	2936	2940	rBV4	66755	93008	0.10%	0.053%
32	21.292	3090	3095	3106	rBV	3201080	4076364	4.52%	2.311%
33	23.498	3465	3470	3476	rBV2	314370	555959	0.62%	0.315%
34	23.651	3489	3496	3509	rVB	2237068	4424270	4.90%	2.509%

Sum of corrected areas: 176359038

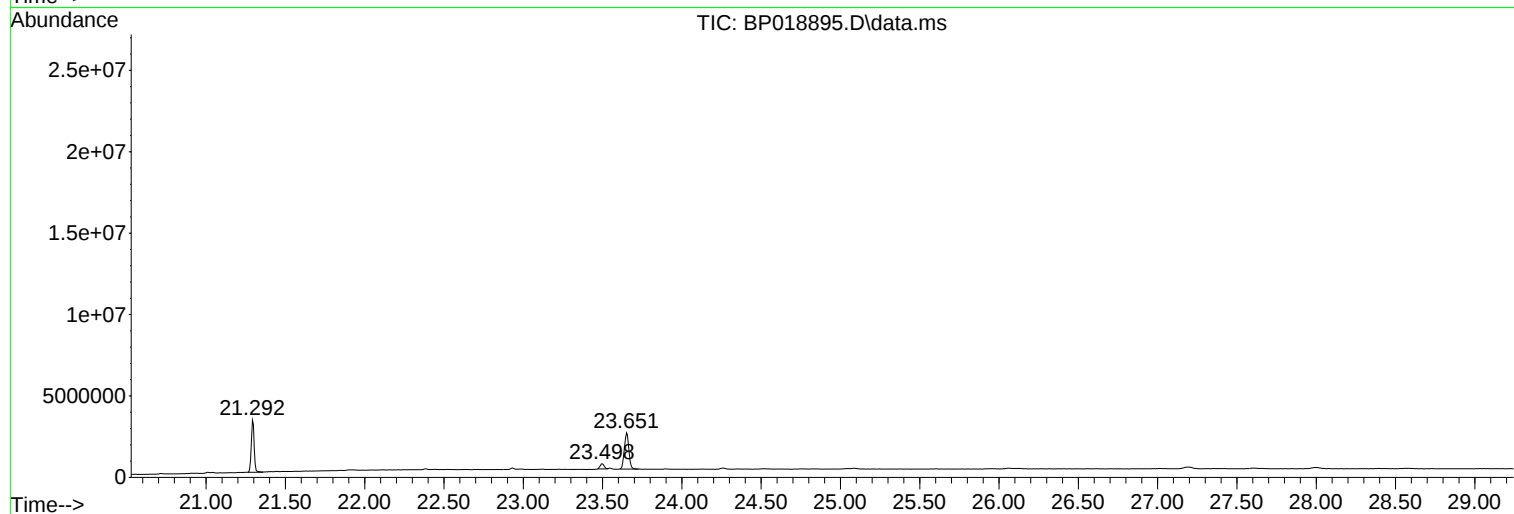
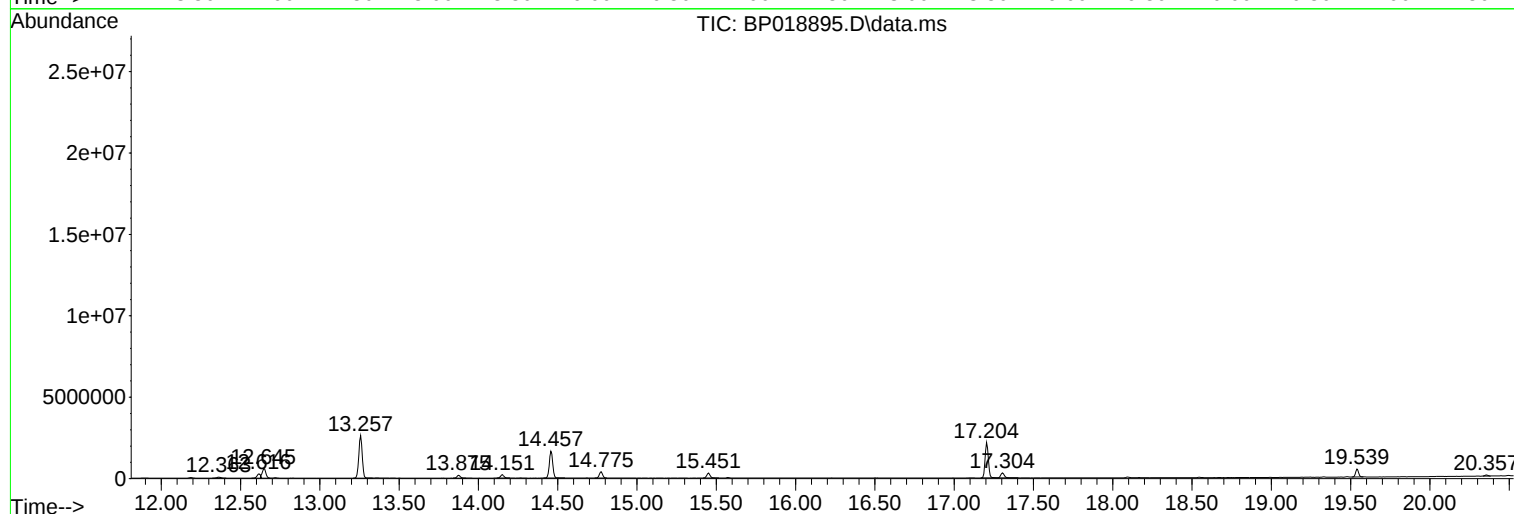
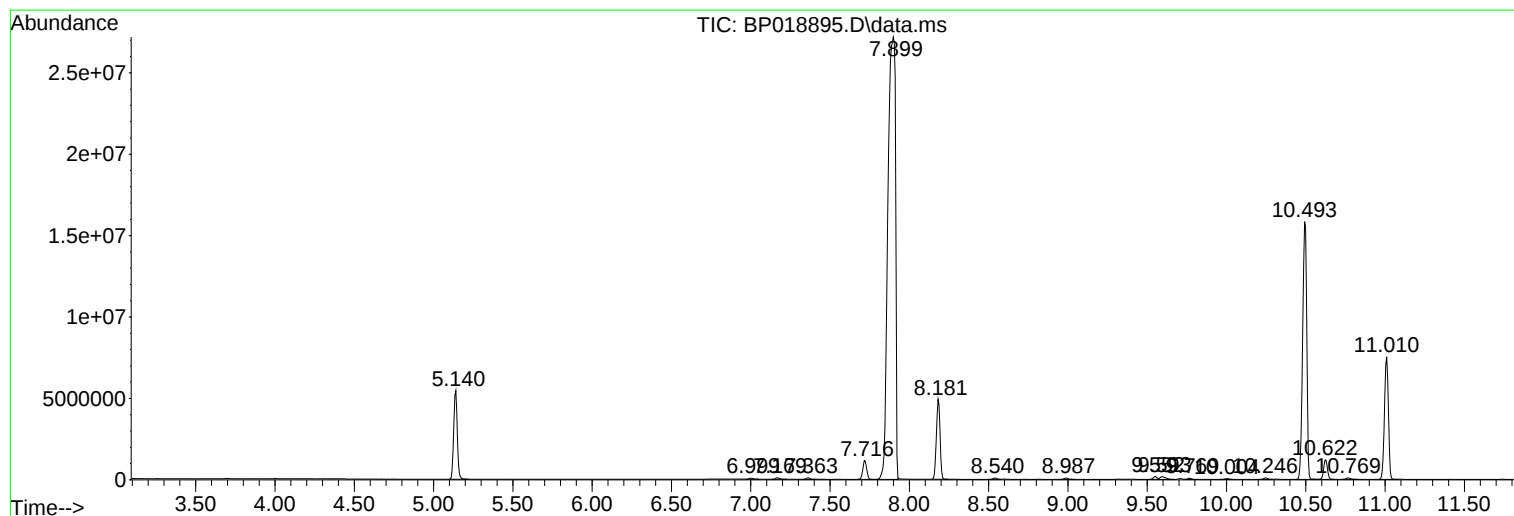
Instrument :
BNA_P
ClientSampleId :
C0AS1DL

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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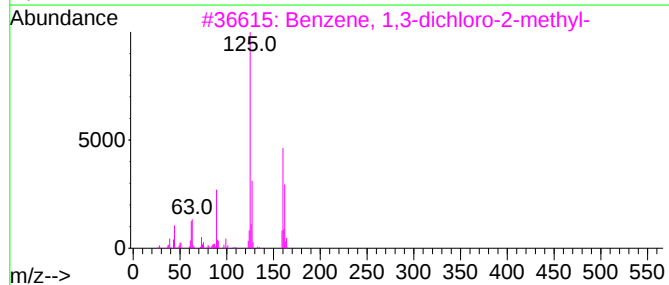
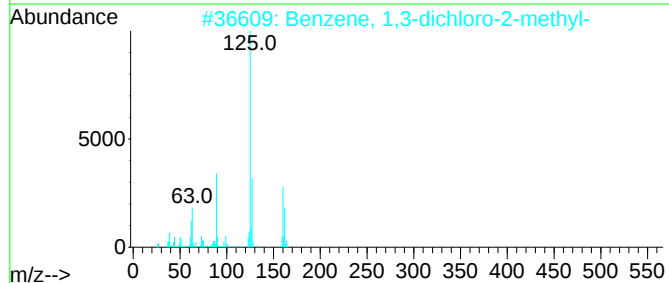
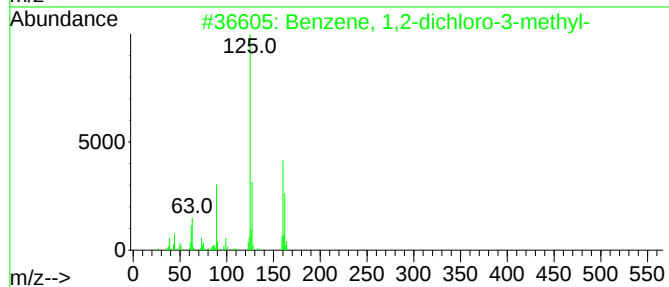
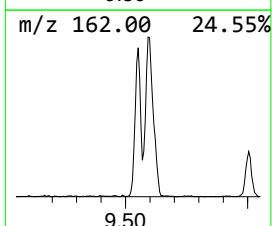
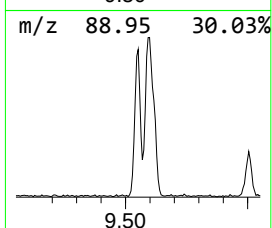
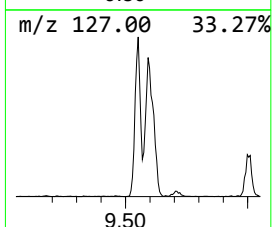
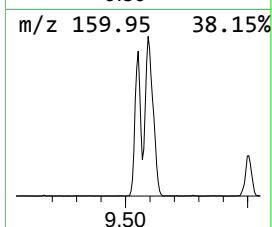
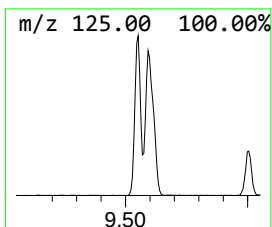
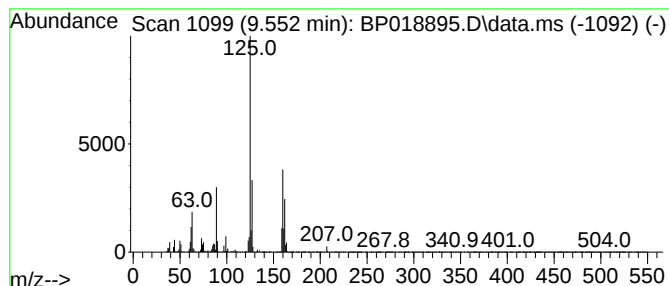
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2-dichloro-3-met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.552	2.97 ng/ul	291734	Naphthalene-d8	10.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
3			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
4			Benzene, 2,4-dichloro-1-methyl-	160	C7H6Cl2	000095-73-8	97
5			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97



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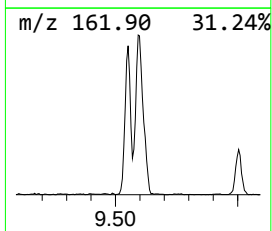
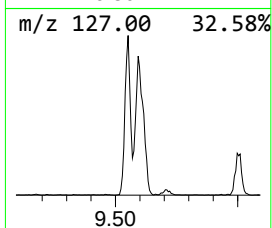
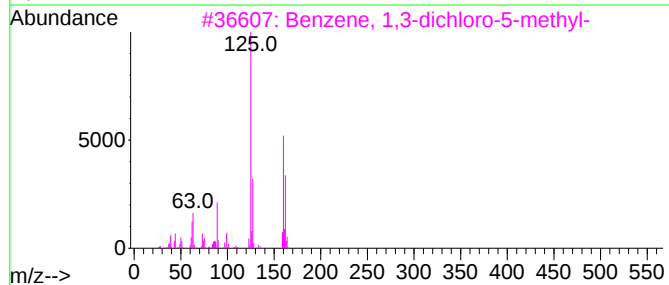
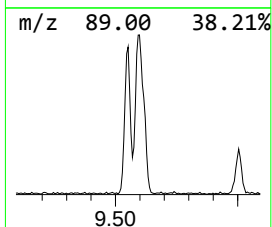
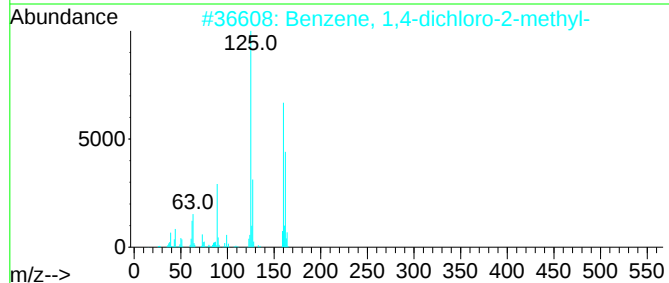
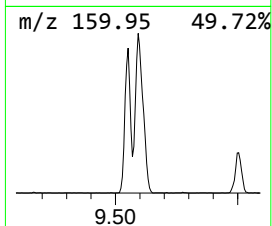
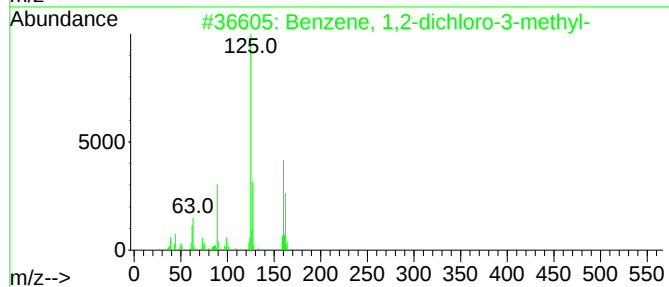
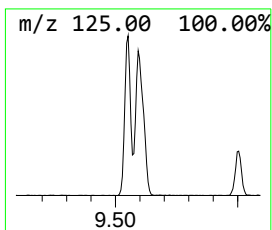
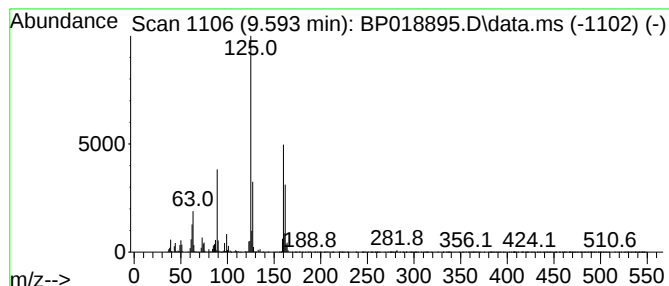
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,4-dichloro-2-met... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.593	4.14 ng/ul	406259	Naphthalene-d8	10.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2			Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
3			Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	96
4			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96
5			Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96



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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, 1,2-di...	9.552	3.0	ng/u1	291734	2	10.622	1963830	20.0
Benzene, 1,4-di...	9.593	4.1	ng/u1	406259	2	10.622	1963830	20.0