

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121323\
 Data File : BP018854.D
 Acq On : 14 Dec 2023 21:32
 Operator : MA/JU
 Sample : 05646-16DL 10X
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AS3DL

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: BP018854.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.687	99	102	111	rBV	50705	83388	0.18%	0.058%
2	5.140	341	349	359	rBV	1533454	2213645	4.89%	1.550%
3	7.005	660	666	681	rVB2	99493	215492	0.48%	0.151%
4	7.169	689	694	705	rVB	93965	149525	0.33%	0.105%
5	7.363	720	727	739	rBV	101714	165131	0.36%	0.116%
6	7.722	779	788	798	rVB	1088328	1777391	3.93%	1.244%
7	7.881	800	815	821	rBV	21675623	42833815	94.65%	29.991%
8	8.187	859	867	875	rVB	5939854	9872746	21.82%	6.913%
9	8.546	922	928	933	rBV	112812	182111	0.40%	0.128%
10	8.604	933	938	945	rVB	63408	105695	0.23%	0.074%
11	8.993	997	1004	1010	rBV	87458	148881	0.33%	0.104%
12	9.551	1092	1099	1103	rBV	180388	301694	0.67%	0.211%
13	9.599	1103	1107	1121	rVB3	175761	486665	1.08%	0.341%
14	9.710	1121	1126	1131	rBV	54314	93114	0.21%	0.065%
15	9.769	1131	1136	1142	rVB2	61810	99848	0.22%	0.070%
16	10.004	1166	1176	1183	rBV	49113	94549	0.21%	0.066%
17	10.251	1211	1218	1225	rBV	106991	184465	0.41%	0.129%
18	10.504	1250	1261	1268	rBV	22687686	45255280	100.00%	31.687%
19	10.634	1276	1283	1292	rVB	1094359	1776851	3.93%	1.244%
20	10.769	1299	1306	1319	rBV	107440	193606	0.43%	0.136%
21	11.016	1339	1348	1362	rVB	9178785	15595584	34.46%	10.920%
22	12.193	1540	1548	1556	rVB	46793	79196	0.17%	0.055%
23	12.369	1571	1578	1583	rBV2	56515	98113	0.22%	0.069%
24	12.622	1612	1621	1622	rBV	230766	322195	0.71%	0.226%
25	12.651	1622	1626	1633	rVV	835970	1289607	2.85%	0.903%
26	12.722	1633	1638	1645	rVB3	26115	45350	0.10%	0.032%
27	13.263	1721	1730	1743	rBV	3173874	4961047	10.96%	3.474%
28	13.875	1829	1834	1841	rBV	166325	242126	0.54%	0.170%
29	14.151	1876	1881	1888	rBV	184473	274649	0.61%	0.192%
30	14.463	1927	1934	1941	rBV	1405899	2120985	4.69%	1.485%
31	14.775	1981	1987	2001	rVB	376639	553254	1.22%	0.387%
32	15.451	2096	2102	2115	rBV	243914	377432	0.83%	0.264%
33	15.575	2118	2123	2134	rVB4	24375	47828	0.11%	0.033%
34	17.210	2394	2401	2412	rBV	1599769	2399605	5.30%	1.680%
35	17.310	2412	2418	2427	rVB	251294	390313	0.86%	0.273%
36	18.098	2547	2552	2562	rBV3	52400	83994	0.19%	0.059%

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ClientSampleId :
C0AS3DL

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

37	19.545	2793	2798	2803	rVB	380170	497507	1.10%	0.348%
38	19.880	2851	2855	2859	rBV	50049	59522	0.13%	0.042%
39	20.357	2933	2936	2940	rBV3	54490	64709	0.14%	0.045%
40	21.298	3091	3096	3103	rVB	2477472	3083752	6.81%	2.159%
41	23.498	3466	3470	3476	rVB	267693	456119	1.01%	0.319%
42	23.657	3490	3497	3513	rVB	1758609	3545052	7.83%	2.482%

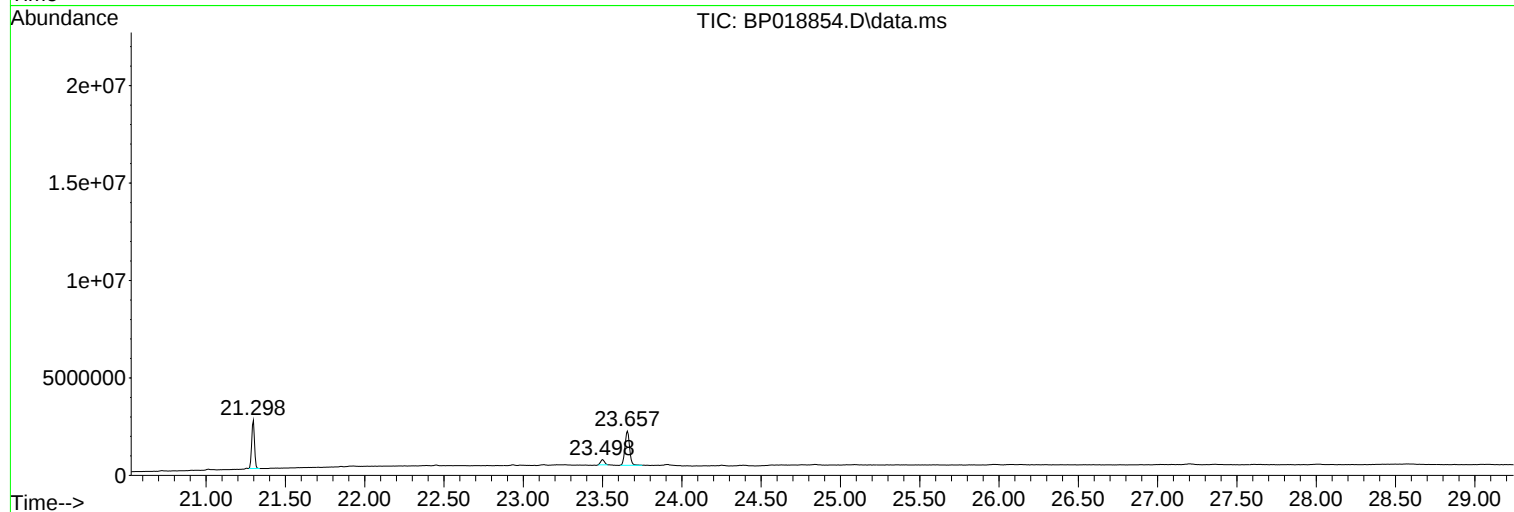
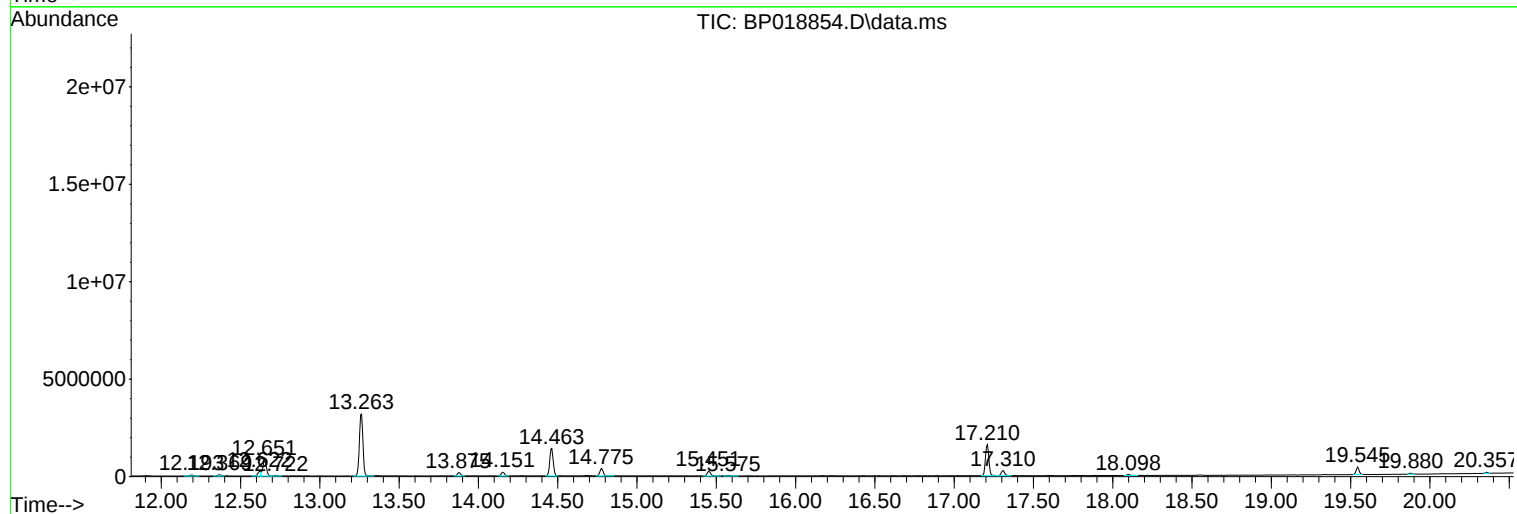
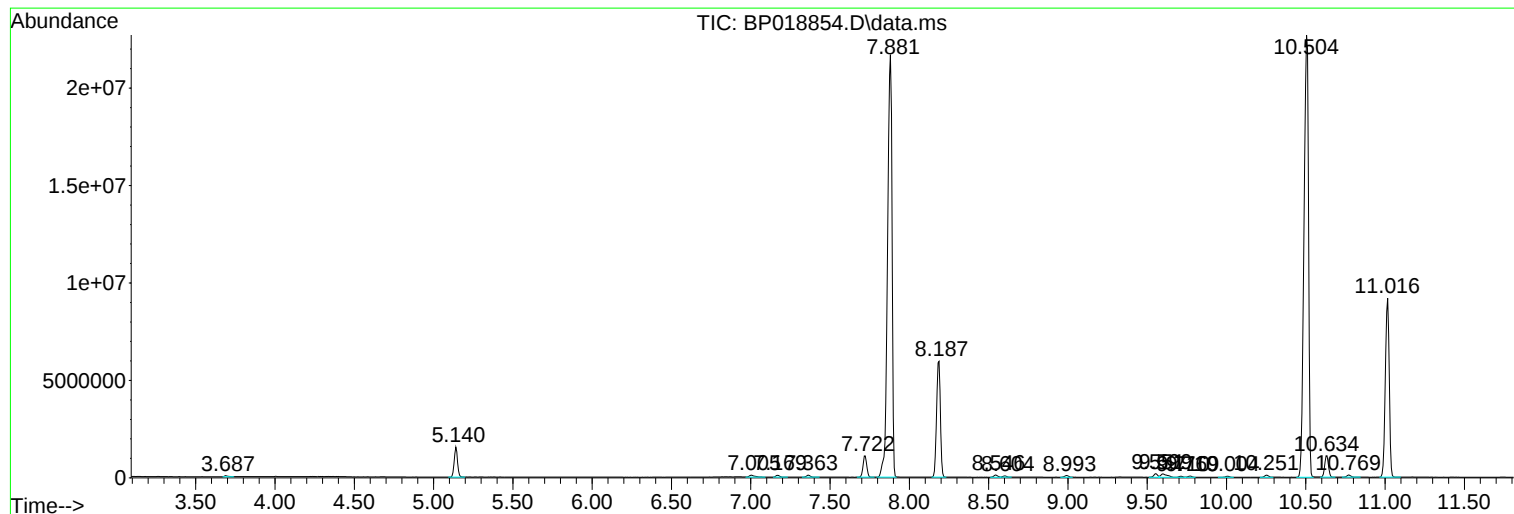
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AS3DL

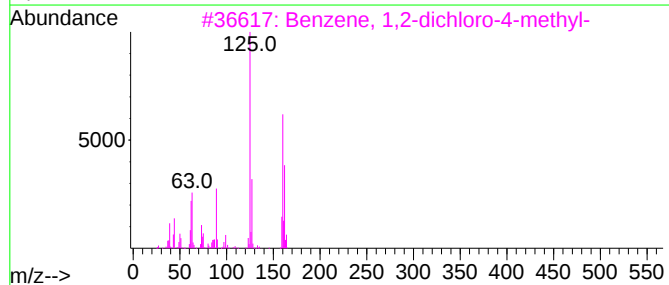
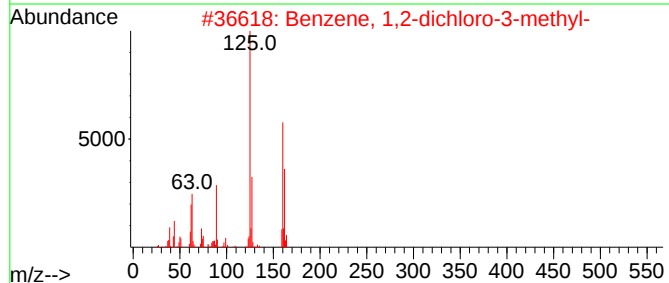
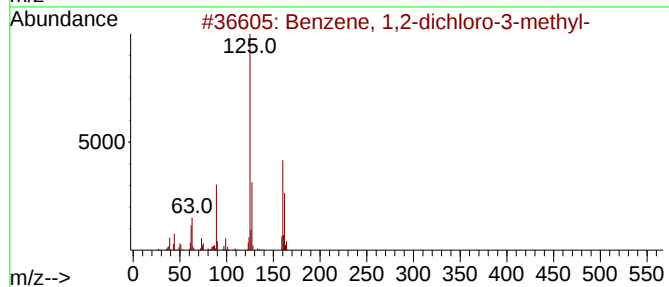
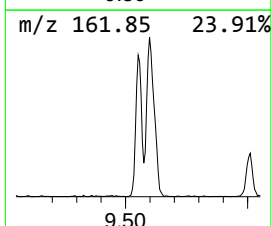
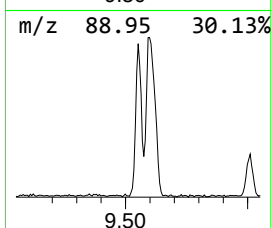
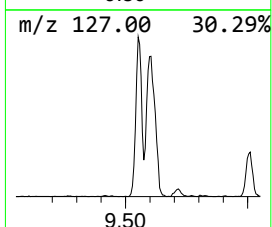
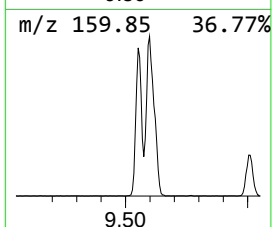
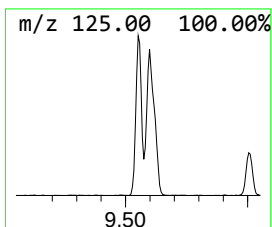
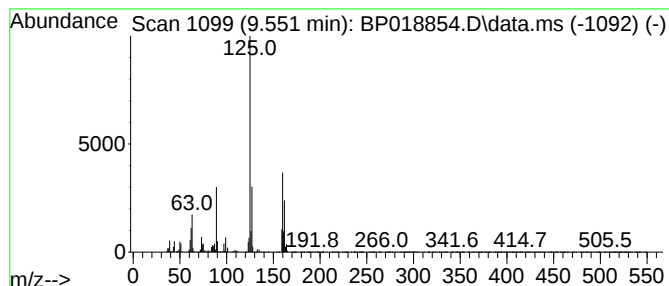
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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.
9.551	3.40 ng/ul	301694	Naphthalene-d8	10.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
3			Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97
4			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
5			Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97



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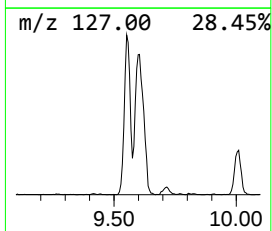
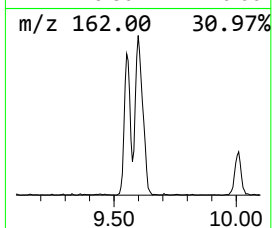
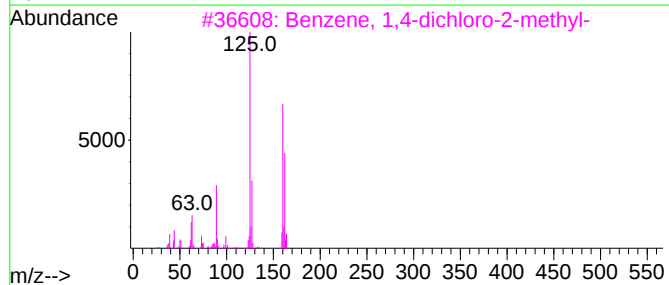
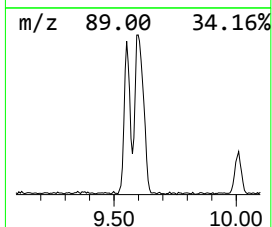
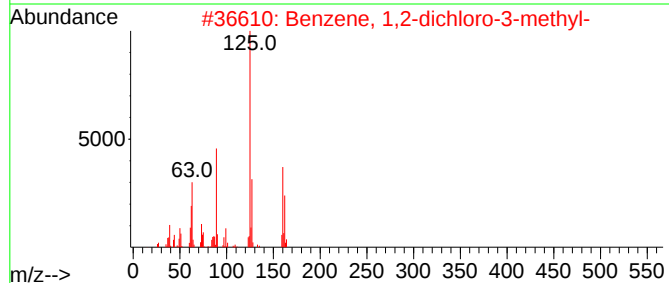
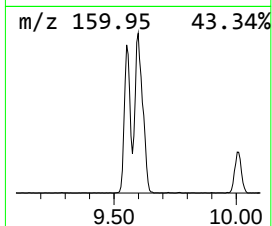
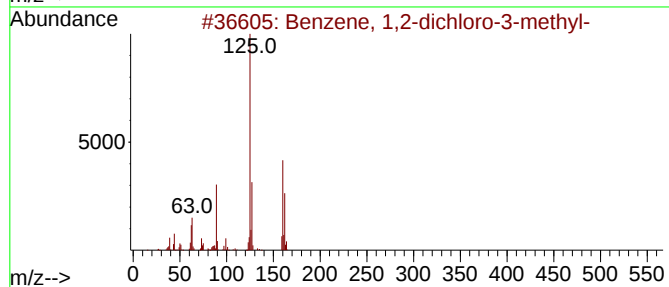
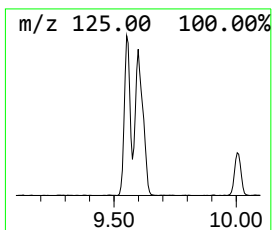
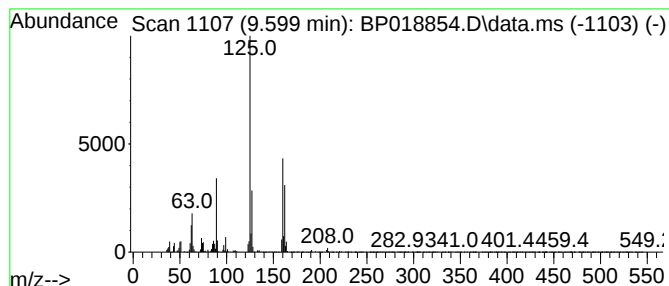
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.599	5.48 ng/ul	486665	Naphthalene-d8	10.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
3			Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
4			Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	96
5			Benzene, 1-chloro-2-(chloromethyl)-	160	C7H6Cl2	000611-19-8	96



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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,2-di...	9.551	3.4	ng/ul	301694	2	10.634	1776850	20.0
Benzene, 1,4-di...	9.599	5.5	ng/ul	486665	2	10.634	1776850	20.0