

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
 Data File : BN034408.D
 Acq On : 05 Oct 2024 03:12
 Operator : RC/JU
 Sample : P4227-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP(20240926)

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_N\Methods\SFAM-EPA-BN091124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN034408.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.999	15	19	21	rBV	36934	49122	0.13%	0.032%
2	3.028	21	24	36	rVB	55111	87514	0.23%	0.057%
3	3.387	81	85	98	rBV	209714	367246	0.97%	0.240%
4	3.687	132	136	142	rVB	33919	47402	0.13%	0.031%
5	4.764	312	319	338	rVV	10372235	16419407	43.49%	10.751%
6	4.928	342	347	354	rVB2	31188	57280	0.15%	0.038%
7	6.581	618	628	640	rBV	368800	607571	1.61%	0.398%
8	6.734	647	654	666	rBV	614591	964706	2.56%	0.632%
9	6.922	679	686	698	rBV	769746	1281525	3.39%	0.839%
10	7.275	738	746	757	rBV	9812836	16085032	42.61%	10.532%
11	7.434	758	773	778	rBV	17496158	37750422	100.00%	24.717%
12	7.722	815	822	834	rVB	1385641	2128472	5.64%	1.394%
13	8.093	878	885	896	rBV	813712	1327046	3.52%	0.869%
14	8.357	924	930	942	rVB	220118	341409	0.90%	0.224%
15	8.522	952	958	968	rVB	701217	1121293	2.97%	0.734%
16	9.181	1064	1070	1074	rBV	165938	266557	0.71%	0.175%
17	9.234	1074	1079	1087	rVV	613887	995177	2.64%	0.652%
18	9.769	1160	1170	1177	rBV	1024155	1768142	4.68%	1.158%
19	9.834	1177	1181	1191	rVB	164100	281638	0.75%	0.184%
20	10.004	1205	1210	1213	rBV	31976	53471	0.14%	0.035%
21	10.063	1213	1220	1225	rVV	55423	121972	0.32%	0.080%
22	10.134	1225	1232	1238	rVV	974233	1596127	4.23%	1.045%
23	10.187	1238	1241	1247	rVB	41863	61283	0.16%	0.040%
24	10.310	1247	1262	1277	rBV2	2096310	5021971	13.30%	3.288%
25	11.628	1480	1486	1488	rBV2	23154	38759	0.10%	0.025%
26	12.210	1581	1585	1592	rVB2	34493	58434	0.15%	0.038%
27	13.063	1722	1730	1735	rBV4	36971	77780	0.21%	0.051%
28	13.269	1759	1765	1771	rBV	59858	102224	0.27%	0.067%
29	13.375	1774	1783	1791	rVV4	22688	57393	0.15%	0.038%
30	13.463	1791	1798	1807	rVB	1827571	2508408	6.64%	1.642%
31	13.716	1834	1841	1855	rBV2	1892274	2900990	7.68%	1.899%
32	14.034	1888	1895	1911	rBV	1500633	2287722	6.06%	1.498%
33	14.269	1931	1935	1944	rVB	198843	321556	0.85%	0.211%
34	14.604	1983	1992	1997	rBV	13434423	24304567	64.38%	15.914%
35	14.739	2007	2015	2022	rVB	354670	529063	1.40%	0.346%
36	15.039	2059	2066	2077	rBV	2540409	3768283	9.98%	2.467%

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
 Data File : BN034408.D
 Acq On : 05 Oct 2024 03:12
 Operator : RC/JU
 Sample : P4227-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP(20240926)

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_N\Methods\SFAM-EPA-BN091124.MA.M
 Title : SVOA CALIBRATION

37	15.169	2082	2088	2100	rBV	941136	1292543	3.42%	0.846%
38	15.345	2114	2118	2122	rVB2	48709	71989	0.19%	0.047%
39	15.416	2126	2130	2136	rVB	40981	59480	0.16%	0.039%
40	15.698	2174	2178	2181	rBV	24830	38057	0.10%	0.025%
41	15.863	2203	2206	2214	rVB	24281	40760	0.11%	0.027%
42	16.451	2302	2306	2316	rVB2	25957	54937	0.15%	0.036%
43	16.792	2358	2364	2371	rBV	1906251	2682676	7.11%	1.757%
44	16.886	2374	2380	2395	rVB	2897785	4243524	11.24%	2.778%
45	17.175	2425	2429	2439	rVB	75432	127026	0.34%	0.083%
46	17.316	2449	2453	2459	rVB	78343	109743	0.29%	0.072%
47	17.722	2518	2522	2529	rBV2	180303	294722	0.78%	0.193%
48	18.757	2695	2698	2703	rVB	46149	57213	0.15%	0.037%
49	18.827	2707	2710	2714	rVV	43557	52899	0.14%	0.035%
50	19.022	2739	2743	2752	rBV2	131191	215043	0.57%	0.141%
51	19.198	2767	2773	2781	rBV2	3893986	5335558	14.13%	3.493%
52	21.022	3078	3083	3093	rBV	2457032	3275058	8.68%	2.144%
53	23.545	3503	3512	3525	rBV	2435469	5492914	14.55%	3.597%
54	23.727	3535	3543	3560	rVB	1558262	3527238	9.34%	2.309%

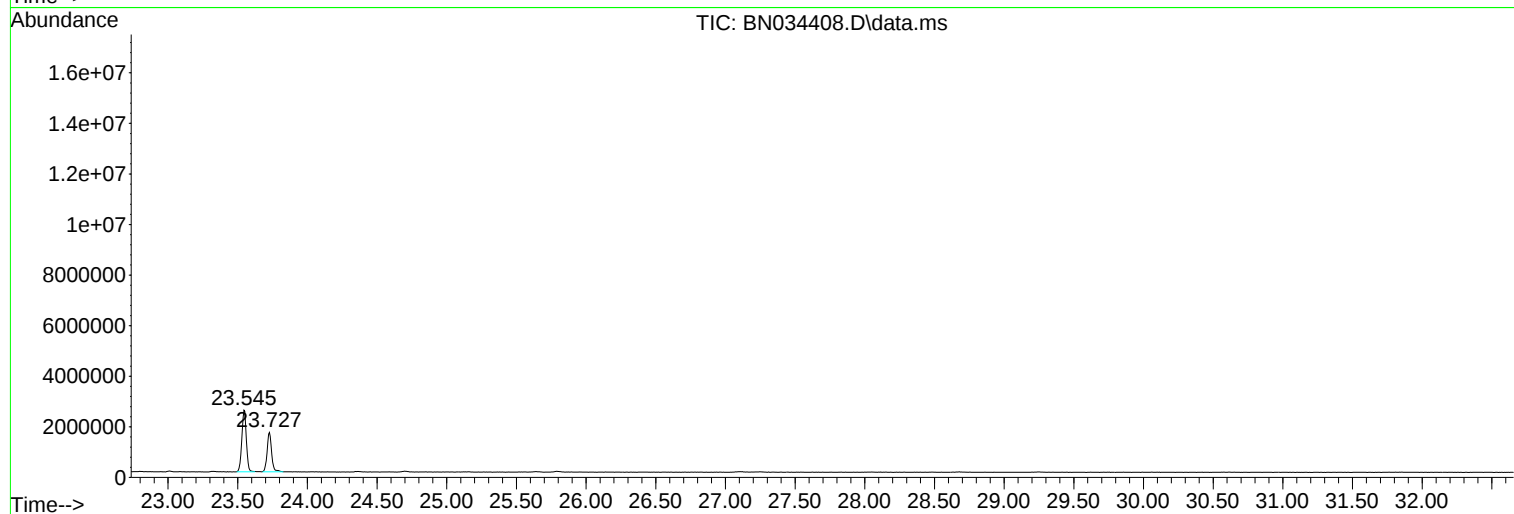
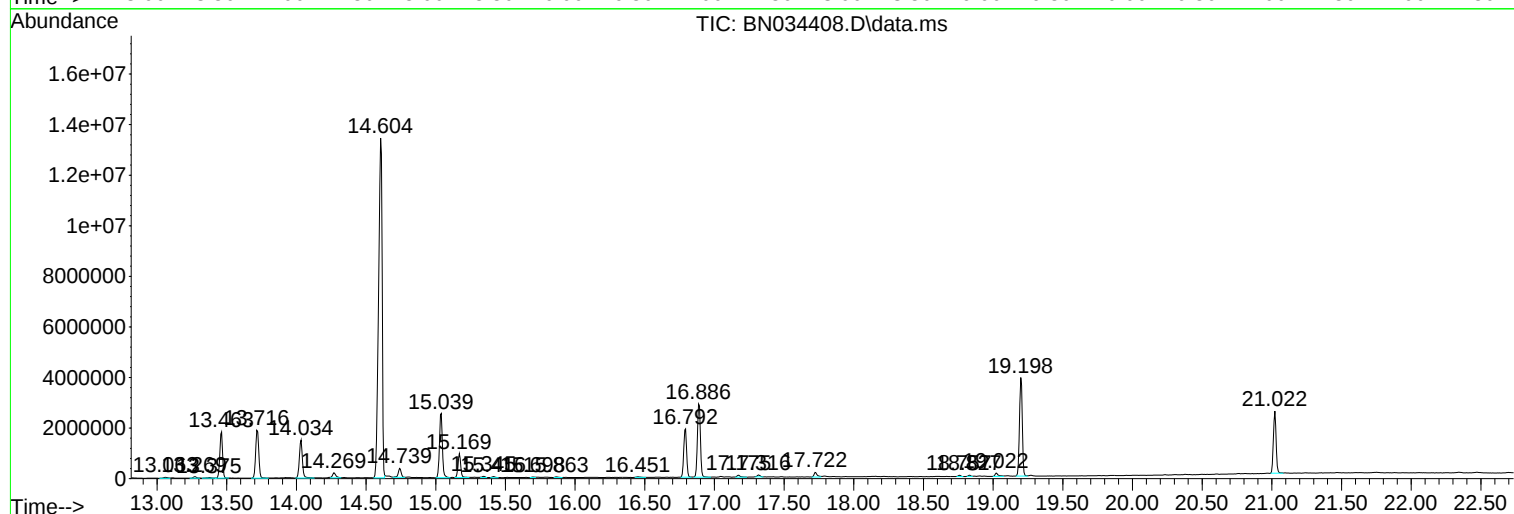
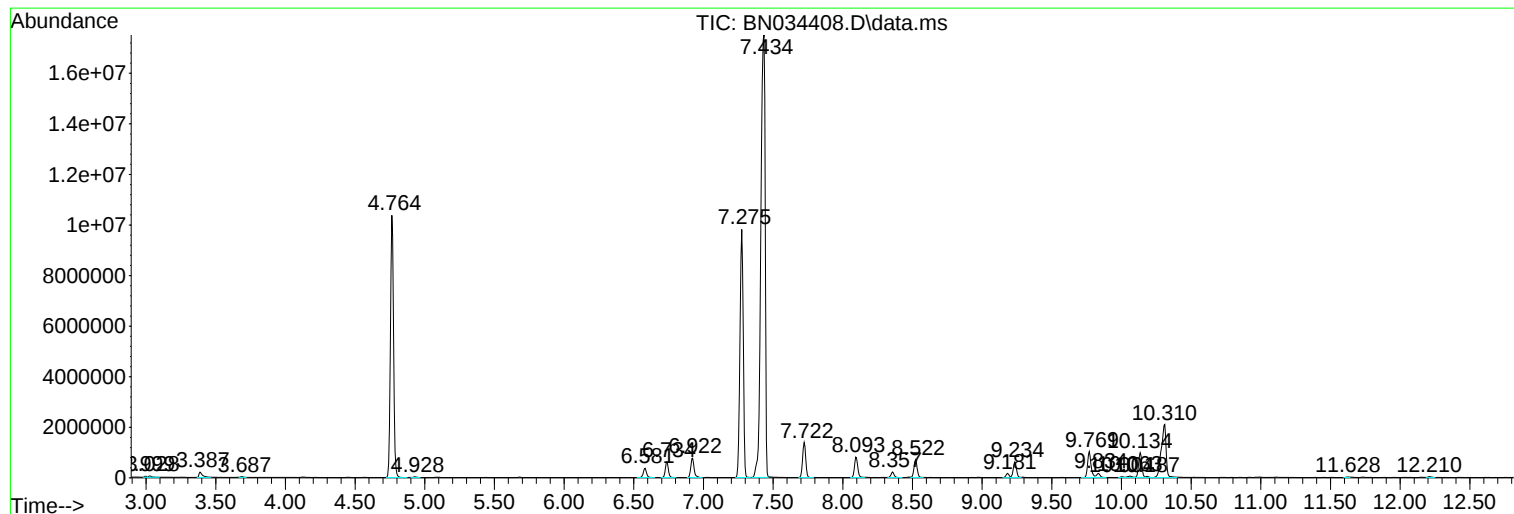
Sum of corrected areas: 152728344

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

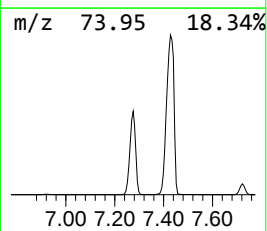
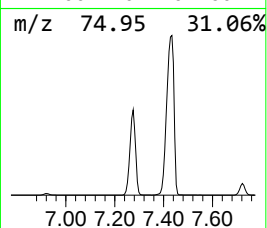
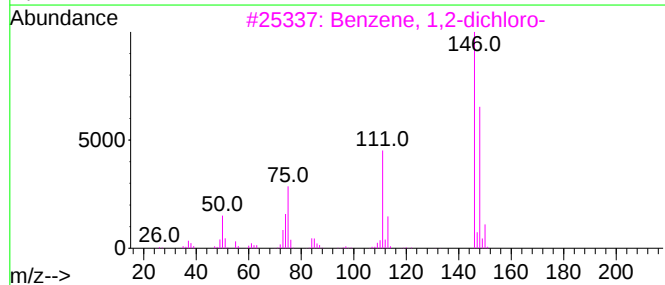
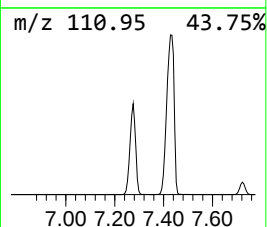
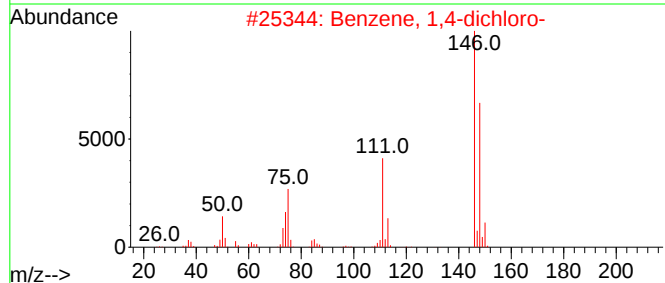
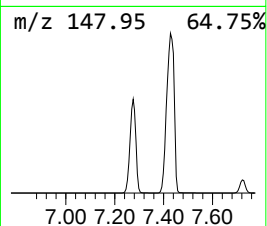
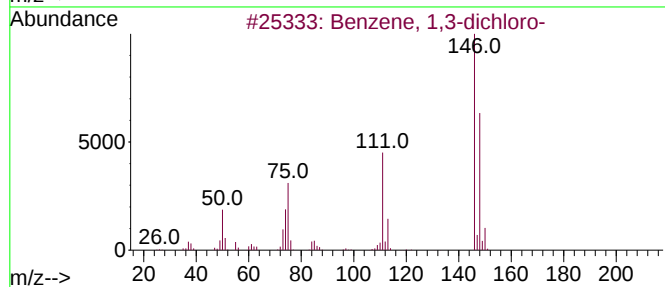
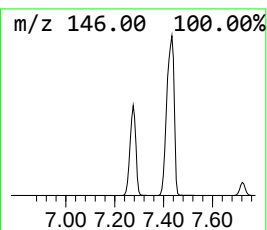
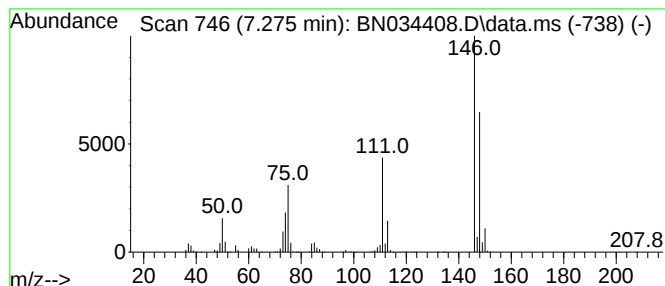
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.275	8.52 ng/ul	16085000	1,4-Dichlorobenzene-d4	7.387

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

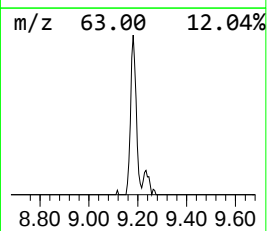
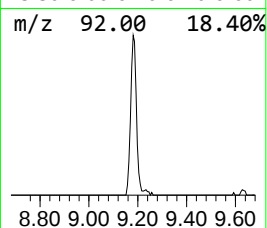
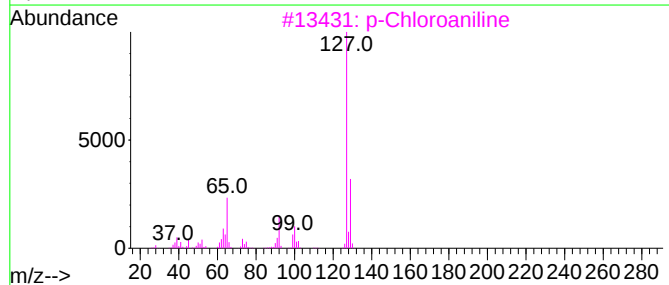
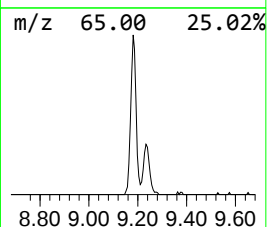
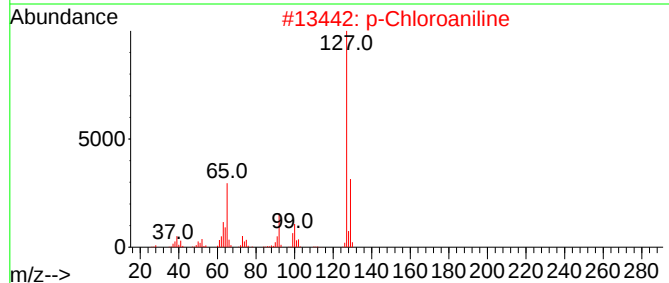
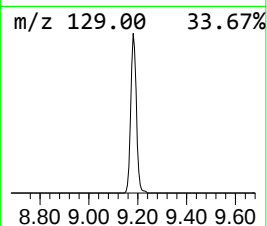
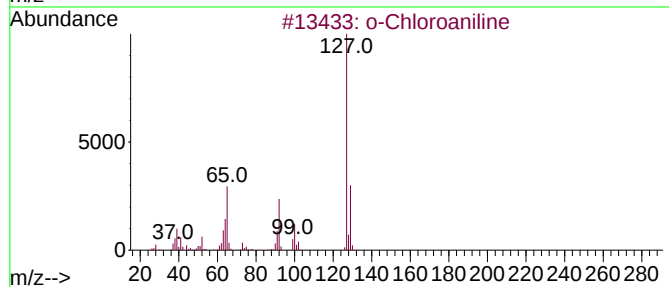
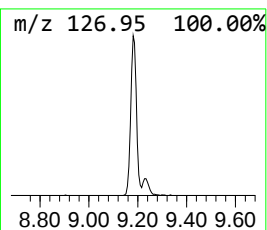
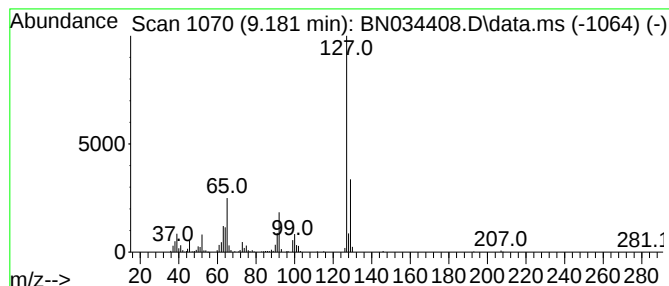
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 o-Chloroaniline Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.181	3.34 ng/ul	266557	Naphthalene-d8	10.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Chloroaniline	127	C6H6ClN	000095-51-2	97
2			p-Chloroaniline	127	C6H6ClN	000106-47-8	97
3			p-Chloroaniline	127	C6H6ClN	000106-47-8	97
4			o-Chloroaniline	127	C6H6ClN	000095-51-2	96
5			o-Chloroaniline	127	C6H6ClN	000095-51-2	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

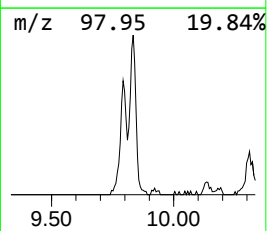
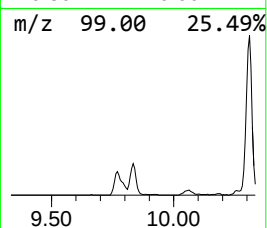
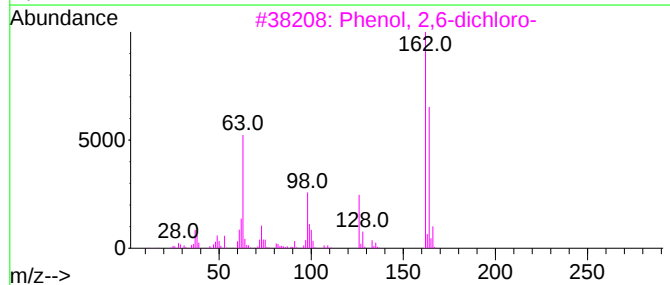
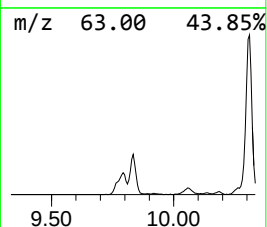
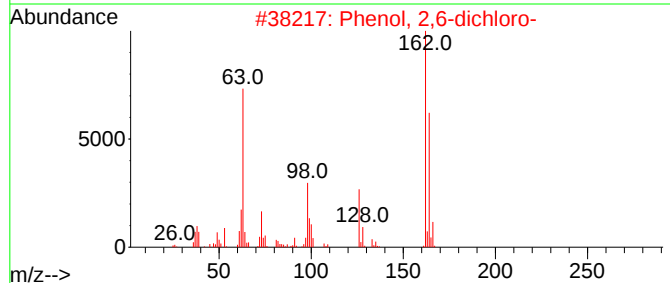
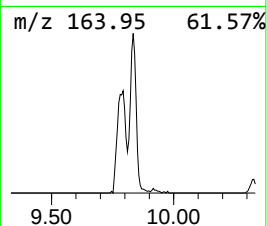
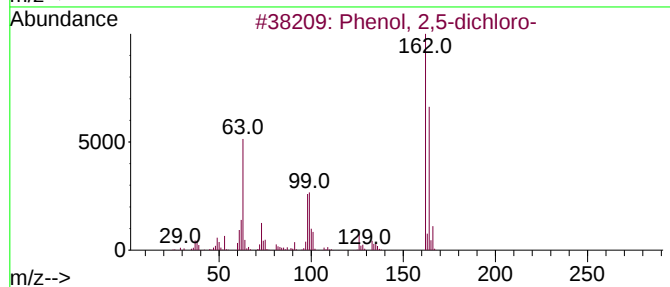
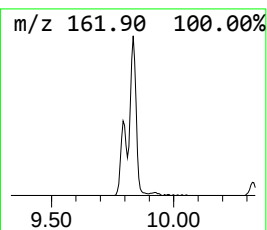
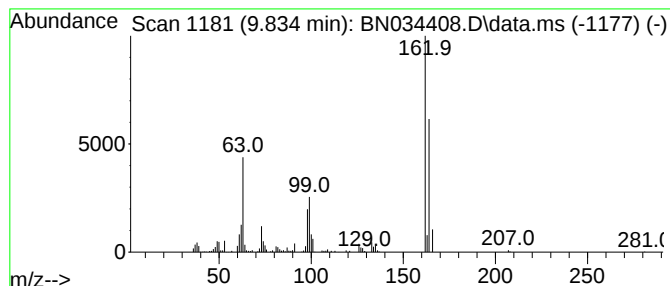
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 2,5-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.834	3.53 ng/ul	281638	Naphthalene-d8	10.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,5-dichloro-	162	C6H4Cl2O	000583-78-8	98
2			Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	94
3			Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	94
4			Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	94
5			Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

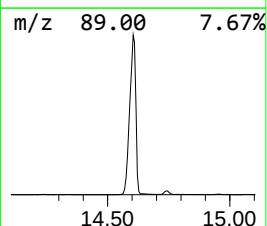
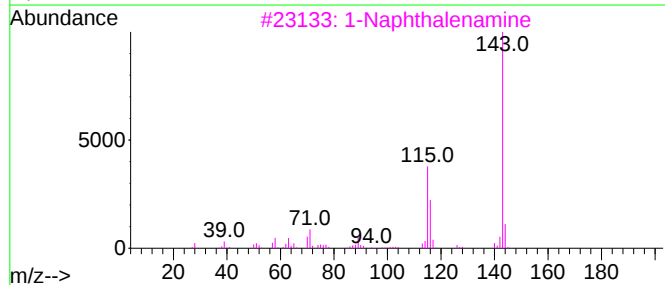
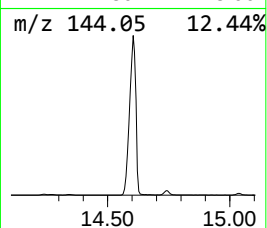
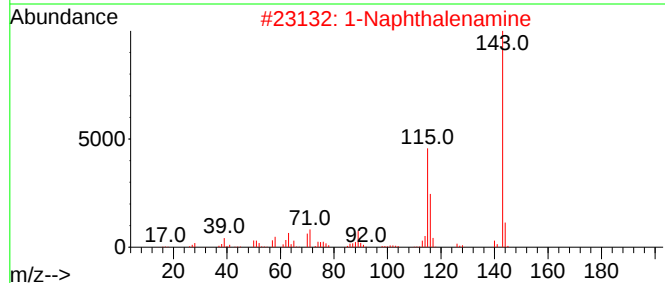
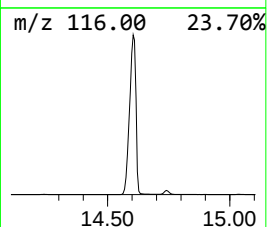
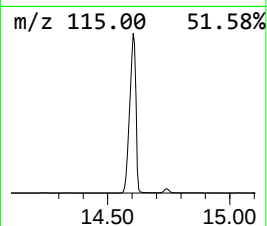
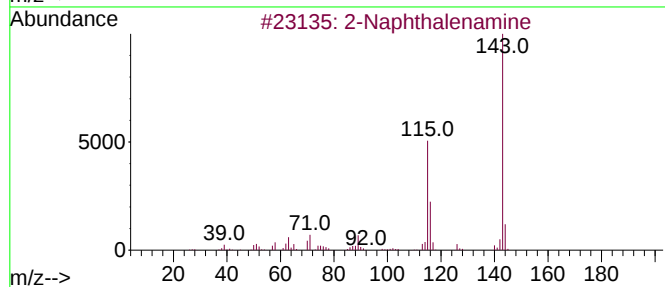
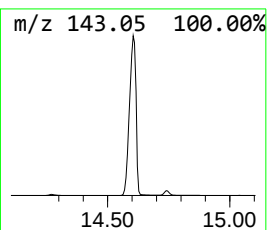
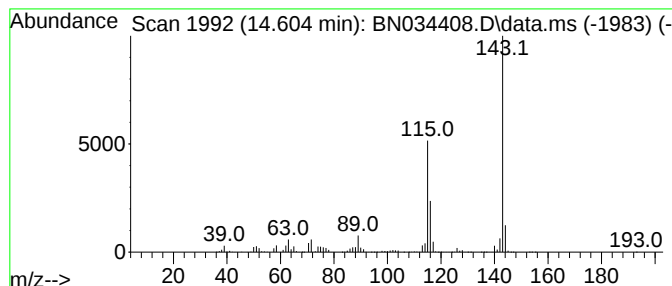
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Naphthalenamine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.604	212.48 ng/ul	24304600	Acenaphthene-d10	14.034

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenamine			143	C10H9N	000091-59-8	97
2	1-Naphthalenamine			143	C10H9N	000134-32-7	97
3	1-Naphthalenamine			143	C10H9N	000134-32-7	97
4	2-Naphthalenamine			143	C10H9N	000091-59-8	95
5	1-Naphthalenamine			143	C10H9N	000134-32-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

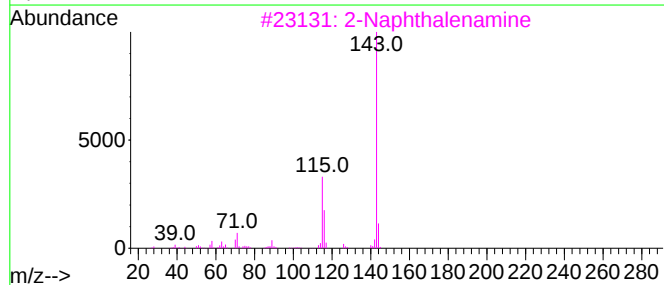
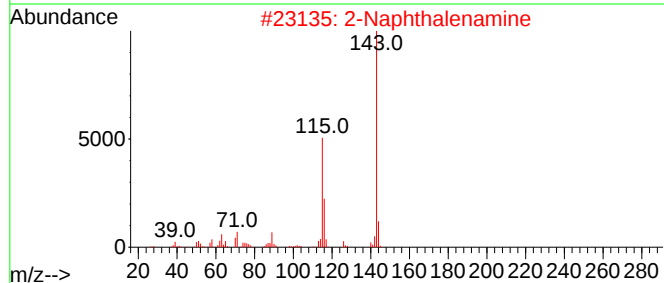
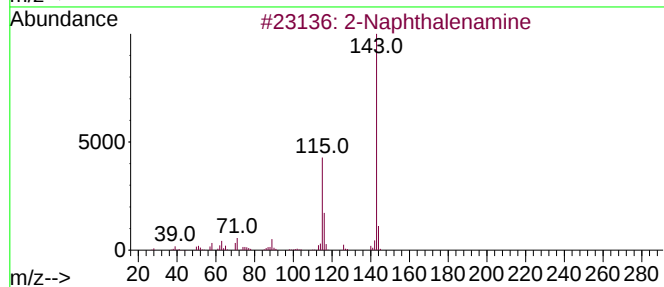
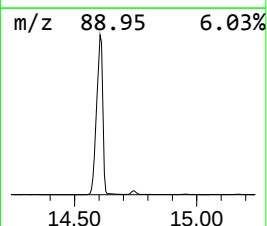
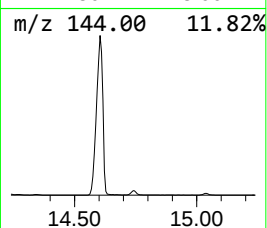
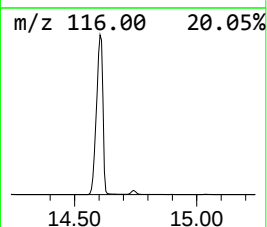
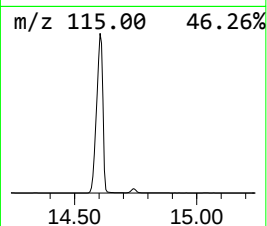
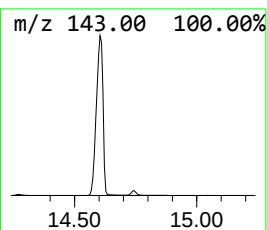
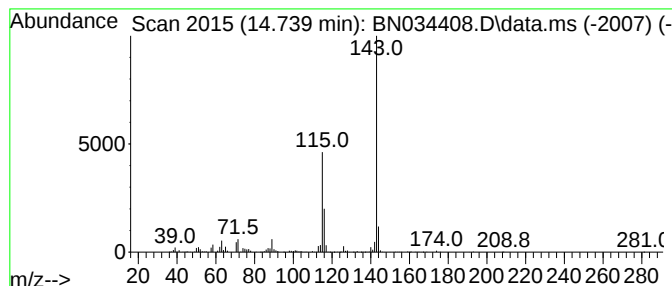
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Naphthalenamine Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.740	4.63 ng/ul	529063	Acenaphthene-d10	14.034

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenamine			143	C10H9N	000091-59-8	97
2	2-Naphthalenamine			143	C10H9N	000091-59-8	96
3	2-Naphthalenamine			143	C10H9N	000091-59-8	95
4	1-Naphthalenamine			143	C10H9N	000134-32-7	95
5	1-Naphthalenamine			143	C10H9N	000134-32-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

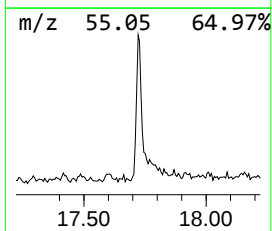
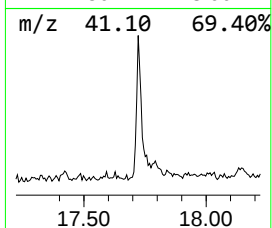
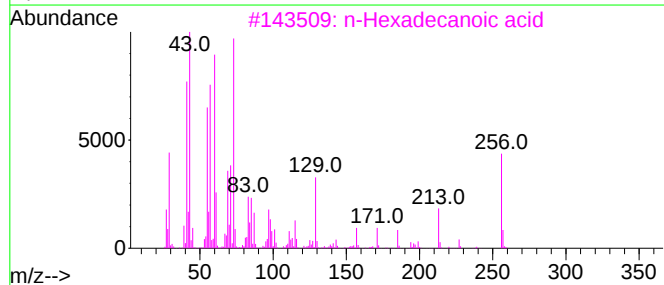
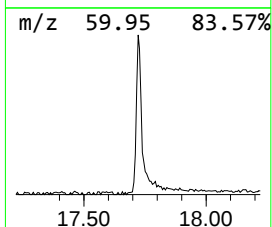
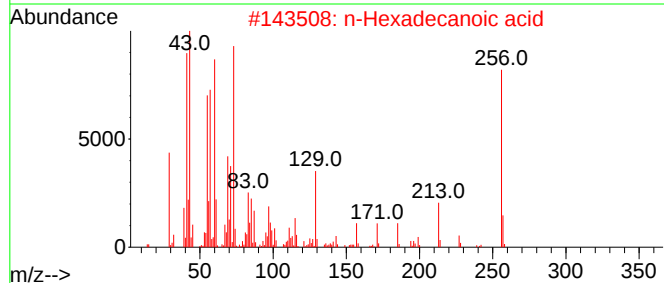
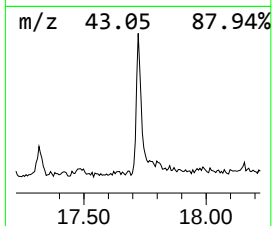
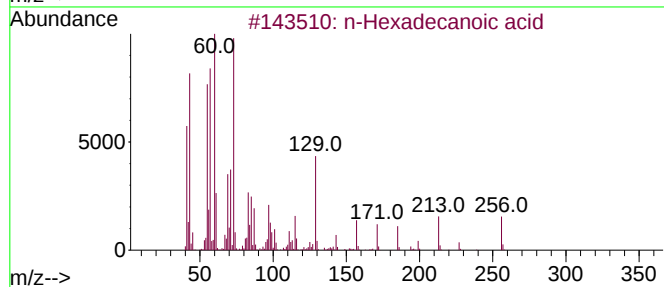
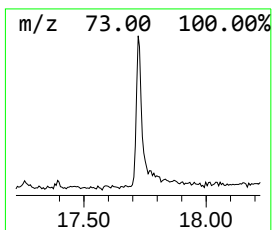
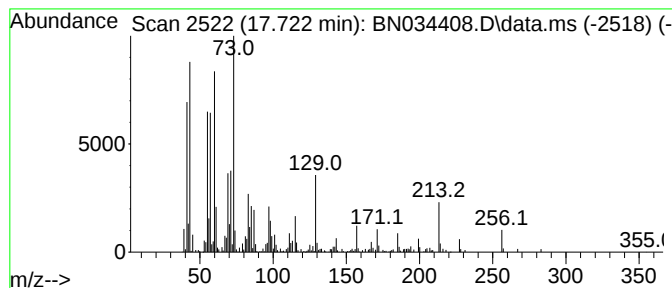
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.722	2.20 ng/ul	294722	Phenanthrene-d10	16.792

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100424\
Data File : BN034408.D
Acq On : 05 Oct 2024 03:12
Operator : RC/JU
Sample : P4227-07
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DUP(20240926)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3-di...	7.275	8.5	ng/ul	16085000	1	7.387	37750400	20.0
o-Chloroaniline	9.181	3.3	ng/ul	266557	2	10.134	1596130	20.0
Phenol, 2,5-dic...	9.834	3.5	ng/ul	281638	2	10.134	1596130	20.0
2-Naphthalenamine	14.604	212.5	ng/ul	24304600	3	14.034	2287720	20.0
1-Naphthalenamine	14.740	4.6	ng/ul	529063	3	14.034	2287720	20.0
n-Hexadecanoic ...	17.722	2.2	ng/ul	294722	4	16.792	2682680	20.0